

### **ATTACHMENT 3: LABORATORY ANALYTICAL REPORTS (SUMMARY SHEETS)**

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This attachment contains the summary sheets from the laboratory analytical reports prepared by Laucks and CAS. Complete analytical reports are available upon request.

# **SAMPLE DATA**

**SDG# JPL28**

**Volatiles**



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-001  
 Lab File ID: M0410021.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:10  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.49	J
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	5.4	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.26	J
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-001

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 15:10

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	3.1	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-001  
 Lab File ID: M0410021.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:10  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-21-5

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL28-001

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/28/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/10/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
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30				

Comments:

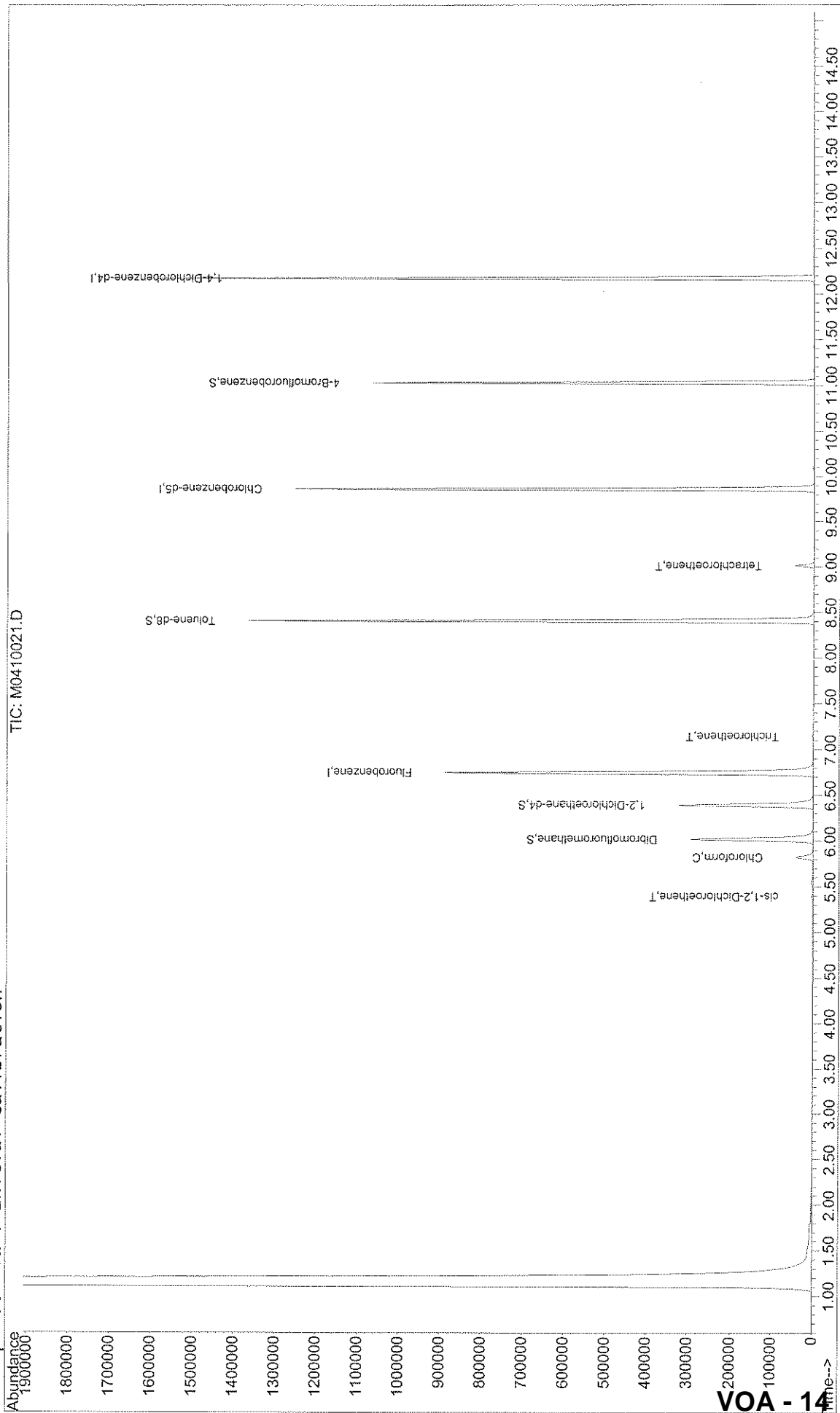
Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410021.D  
Acq On : 10 Apr 2007 15:10  
Sample : JPL28-001  
Misc : #1 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:24 2007

Vial: 58  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00

Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 14

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410021.D  
 Acq On : 10 Apr 2007 15:10  
 Sample : JPL28-001  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:24 2007

Vial: 58  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)

Title : VOA 524- 5ML Calibration 5973M

Last Update : Tue Apr 10 10:47:40 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar)
1) Fluorobenzene	6.75	96	888114	50.00	ug/l	0.00	84.80%
50) Chlorobenzene-d5	9.86	82	371828	50.00	ug/l	0.00	82.08%
70) 1,4-Dichlorobenzene-d4	12.17	152	378256	50.00	ug/l	0.00	80.94%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	214616	50.79	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.58%	
37) 1,2-Dichloroethane-d4	6.39	65	252518	51.80	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	103.60%	
51) Toluene-d8	8.41	98	897944	51.95	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.90%	
72) 4-Bromofluorobenzene	11.03	95	316101	51.09	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	102.18%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	d
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.05	76	118		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	d
19) trans-1,2-Dichloroethene	3.86	96	55		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	3.95	73	129		N.D.	
22) 1,1-Dichloroethane	4.56	63	955		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.40	96	1913	0.49	ug/l	84
27) 2-Butanone	5.54	43	135		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.83	83	38076	5.37	ug/l	95
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.02	75	57		N.D.	
38) Benzene	6.42	78	359		N.D.	

*4/11/07*

Quantitation Report

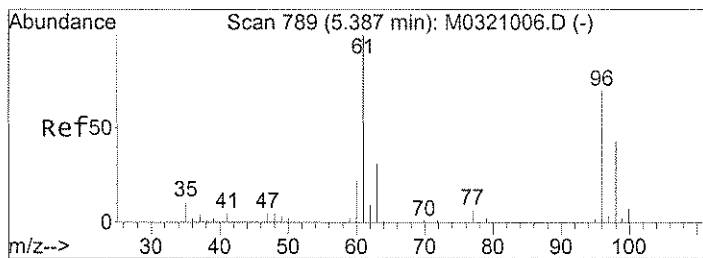
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 Sample : JPL28-001  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:24 2007

Vial: 58  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

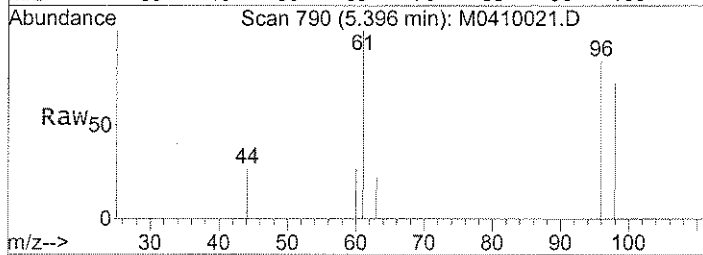
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	7.14	130	1096	0.26 ug/l #	51
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	7.71	83	278	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	8.47	92	279	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	9.02	166	13076	3.10 ug/l	95
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.89	112	200	N.D.	
62) 1-Chlorohexane	9.86	91	1290	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.99	91	828	N.D.	
65) m,p-xylene	10.10	106	1506	N.D.	
66) o-xylene	10.49	106	323	N.D.	
67) Styrene	0.00	104	0	N.D. d	
68) Bromoform	0.00	173	0	N.D.	
69) Isopropylbenzene	10.85	105	143	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	11.04	156	70	N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	282	N.D.	
75) 1,2,3-Trichloropropane	11.03	110	186	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	
77) 2-Chlorotoluene	11.26	91	249	N.D.	
78) 4-Chlorotoluene	11.47	91	206	N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	116	N.D.	
80) tert-Butylbenzene	0.00	119	0	N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0	N.D. d	
82) sec-butylbenzene	0.00	105	0	N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0	N.D. d	
84) 4-Isopropyltoluene	0.00	119	0	N.D. d	
85) 1,4-Dichlorobenzene	12.19	146	362	N.D.	
86) 1,2-Dichlorobenzene	12.57	146	470	N.D.	
87) n-Butylbenzene	12.52	91	357	N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0	N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	564	N.D.	
90) Hexachlorobutadiene	14.30	225	55	N.D.	
91) Naphthalene	0.00	128	0	N.D. d	
92) 1,2,3-Trichlorobenzene	14.66	180	548	N.D.	

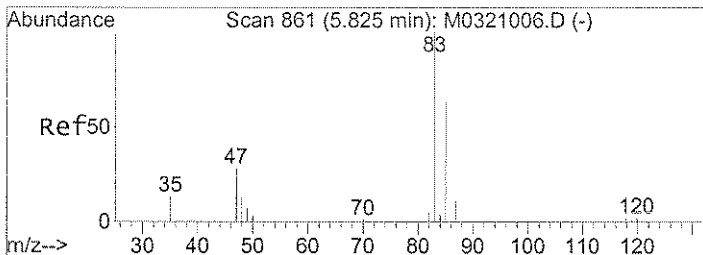
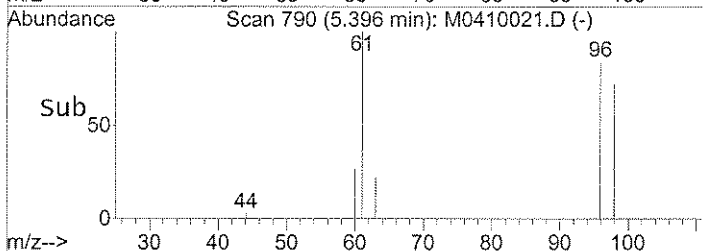
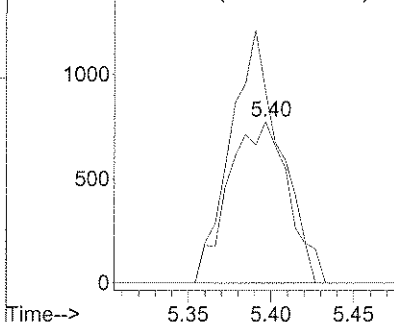


#26  
 cis-1,2-Dichloroethene  
 Concen: 0.49 ug/l  
 RT: 5.40 min Scan# 790  
 Delta R.T. 0.01 min  
 Lab File: M0410021.D  
 Acq: 10 Apr 2007 15:10

Tgt Ion: 96 Resp: 1913  
 Ion Ratio Lower Upper  
 96 100  
 61 134.4 93.4 140.2

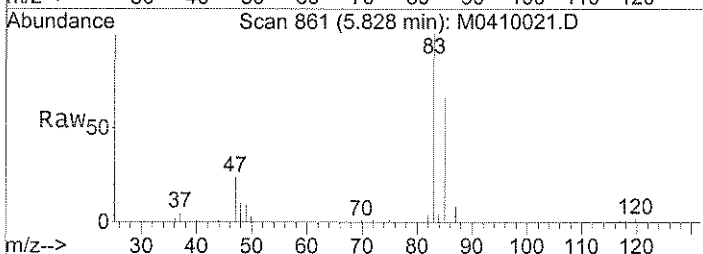


Abundance Ion 95.95 (95.65 to 96.65): M0  
 Ion 61.05 (60.75 to 61.75): M0

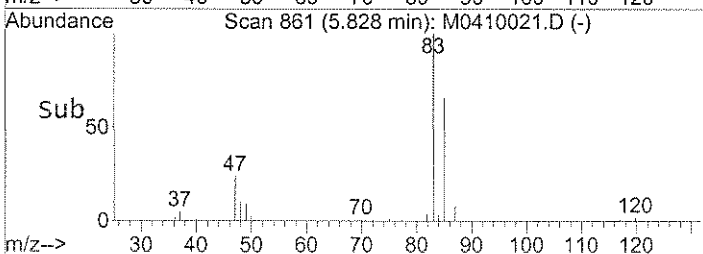
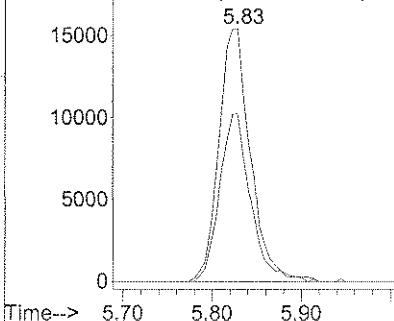


#31  
 Chloroform  
 Concen: 5.37 ug/l  
 RT: 5.83 min Scan# 861  
 Delta R.T. 0.00 min  
 Lab File: M0410021.D  
 Acq: 10 Apr 2007 15:10

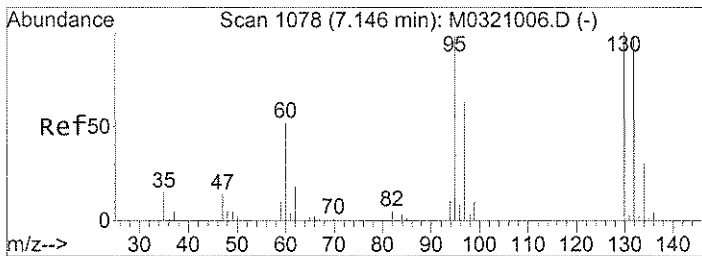
Tgt Ion: 83 Resp: 38076  
 Ion Ratio Lower Upper  
 83 100  
 85 64.9 41.2 81.2



Abundance Ion 83.00 (82.70 to 83.70): M0  
 Ion 85.00 (84.70 to 85.70): M0

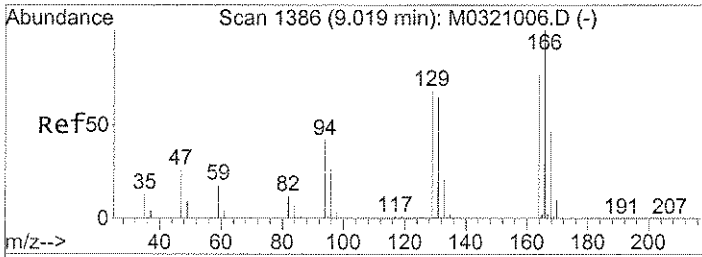
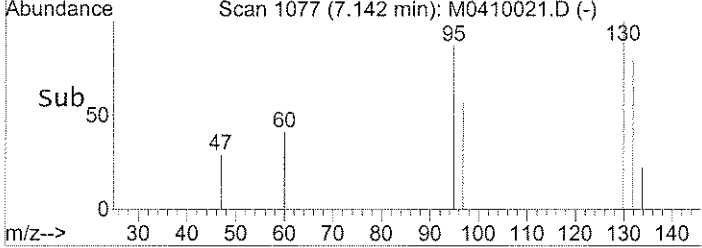
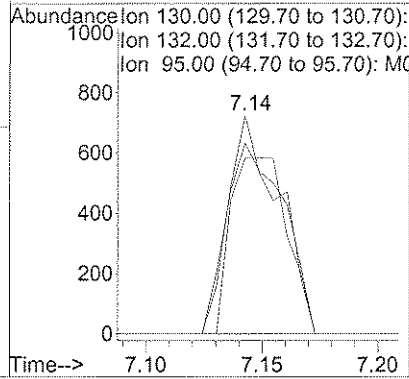
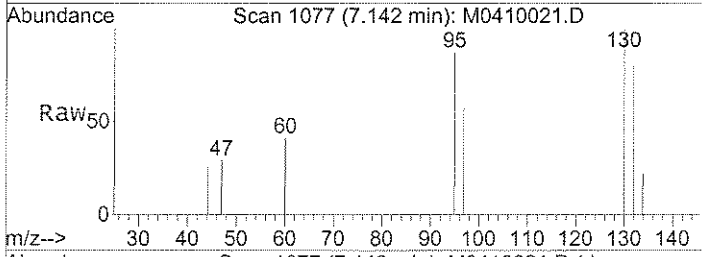






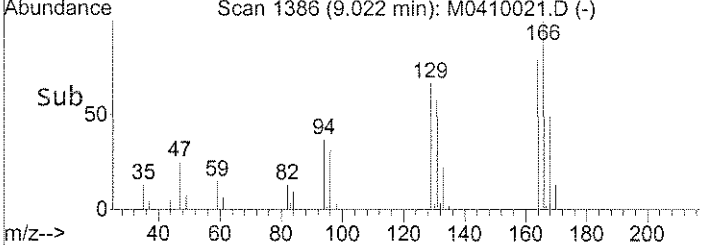
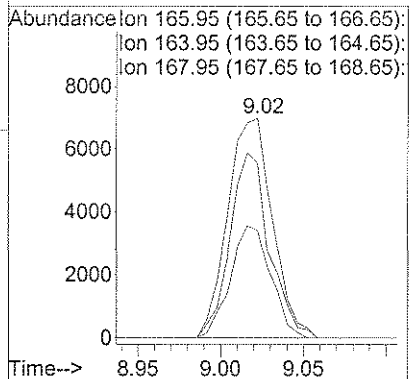
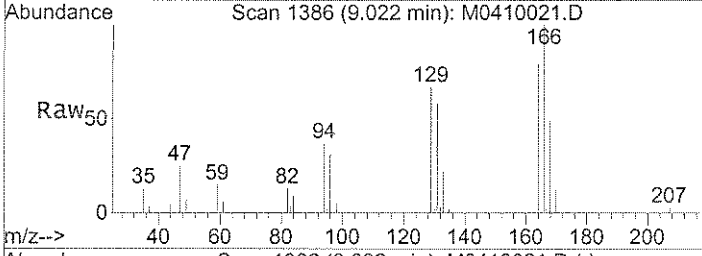
#41  
 Trichloroethene  
 Concen: 0.26 ug/l  
 RT: 7.14 min Scan# 1077  
 Delta R.T. -0.00 min  
 Lab File: M0410021.D  
 Acq: 10 Apr 2007 15:10

Tgt Ion	Resp	Ion Ratio	Lower	Upper
130	1096	100		
132		90.5	75.0	115.0
95		0.0	69.4	109.4#



#56  
 Tetrachloroethene  
 Concen: 3.10 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. 0.00 min  
 Lab File: M0410021.D  
 Acq: 10 Apr 2007 15:10

Tgt Ion	Resp	Ion Ratio	Lower	Upper
166	13076	100		
164		74.1	63.3	94.9
168		46.0	39.6	59.4



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410021.D                   Vial: 58  
Acq On    : 10 Apr 2007 15:10                   Operator: LPM  
Sample    : JPL28-001                         Inst     : MOBY  
Misc      : #1 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410021.D 524.M    wed Apr 11 08:32:00 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-002  
 Lab File ID: M0410022.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:32  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.59	
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	4.6	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 15:32

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	2.6	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-002  
 Lab File ID: M0410022.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:32  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-21-4

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL28-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/28/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/10/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
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17					
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19					
20					
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22					
23					
24					
25					
26					
27					
28					
29					
30					

Comments:

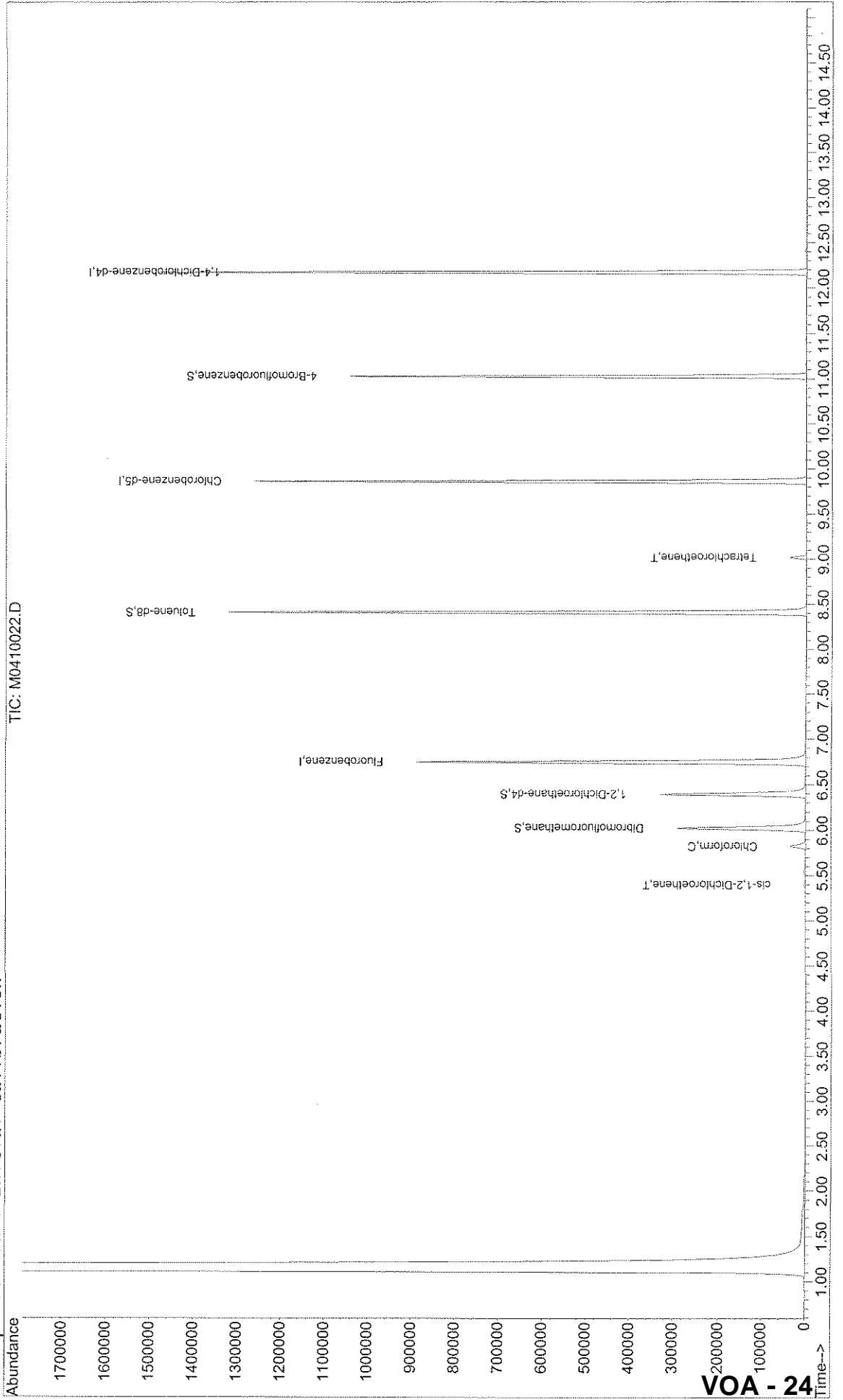
Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410022.D  
Acq On : 10 Apr 2007 15:32  
Sample : JPL28-002  
Misc : #2 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:25 2007

Vial: 53  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00

Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 24

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410022.D  
 Acq On : 10 Apr 2007 15:32  
 Sample : JPL28-002  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:25 2007

Vial: 53  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	881752	50.00	ug/l	0.00 84.20%
50) Chlorobenzene-d5	9.86	82	370119	50.00	ug/l	0.00 81.70%
70) 1,4-Dichlorobenzene-d4	12.18	152	369068	50.00	ug/l	0.00 78.97%

System Monitoring Compounds						
33) Dibromofluoromethane	6.03	111	212064	50.55	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.10%
37) 1,2-Dichloroethane-d4	6.40	65	252552	52.18	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	104.36%
51) Toluene-d8	8.41	98	891495	51.82	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.64%
72) 4-Bromofluorobenzene	11.03	95	310040	51.36	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	102.72%

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	1.46	50	62		N.D.	
4) Vinyl chloride	1.57	62	60		N.D.	
5) Bromomethane	1.88	96	58		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	d
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.06	76	76		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	3.93	73	65		N.D.	
22) 1,1-Dichloroethane	4.57	63	492		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.40	96	2282	0.59	ug/l	99
27) 2-Butanone	5.57	43	66		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.83	83	32193	4.57	ug/l	96
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.14	75	60		N.D.	
38) Benzene	6.41	78	152		N.D.	

*Handwritten signature*



Quantitation Report

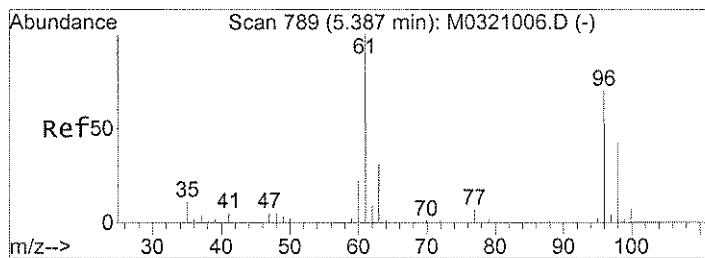
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 Sample : JPL28-002  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:25 2007

Vial: 53  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

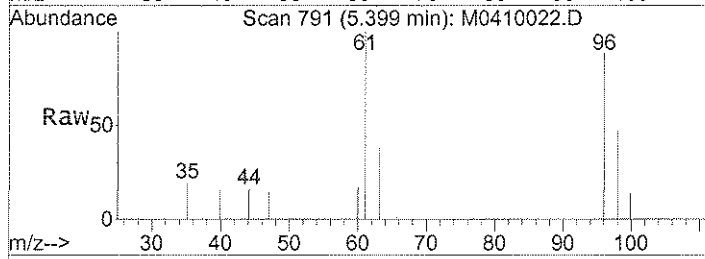
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	6.58	43	110	Below Cal #	22
41) Trichloroethene	7.15	130	941	N.D.	
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D. d	
46) Bromodichloromethane	7.72	83	208	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	8.48	92	79	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	9.02	166	11051	2.63 ug/l	96
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.89	112	154	N.D.	
62) 1-Chlorohexane	9.86	91	1436	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	10.10	91	1899	N.D.	
65) m,p-Xylene	10.10	106	910	N.D.	
66) o-xylene	10.50	106	137	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D. d	
69) Isopropylbenzene	11.03	105	349	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	0.00	156	0	N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	256	N.D.	
75) 1,2,3-Trichloropropane	11.03	110	135	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	
77) 2-Chlorotoluene	11.26	91	60	N.D.	
78) 4-Chlorotoluene	11.26	91	60	N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	204	N.D.	
80) tert-Butylbenzene	11.75	119	73	N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0	N.D. d	
82) sec-butylbenzene	0.00	105	0	N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0	N.D. d	
84) 4-Isopropyltoluene	0.00	119	0	N.D. d	
85) 1,4-Dichlorobenzene	12.20	146	368	N.D.	
86) 1,2-Dichlorobenzene	12.57	146	504	N.D.	
87) n-Butylbenzene	12.52	91	250	N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0	N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	177	N.D.	
90) Hexachlorobutadiene	14.31	225	55	N.D.	
91) Naphthalene	0.00	128	0	N.D. d	
92) 1,2,3-Trichlorobenzene	14.66	180	418	N.D.	

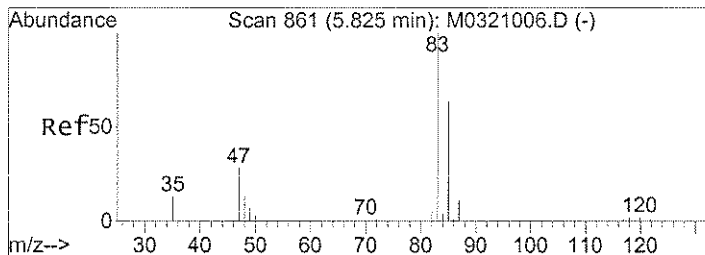
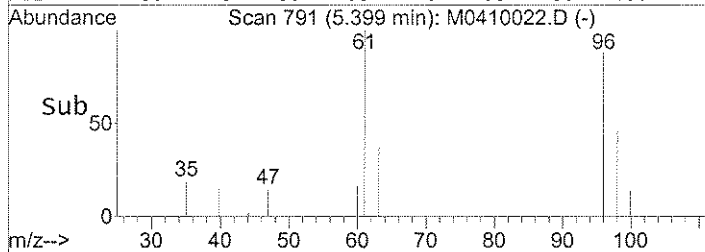
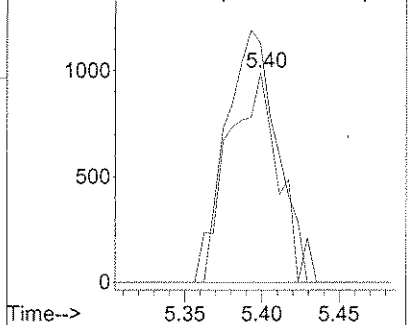


#26  
 cis-1,2-Dichloroethene  
 Concen: 0.59 ug/l  
 RT: 5.40 min Scan# 791  
 Delta R.T. 0.01 min  
 Lab File: M0410022.D  
 Acq: 10 Apr 2007 15:32

Tgt Ion: 96 Resp: 2282  
 Ion Ratio Lower Upper  
 96 100  
 61 117.9 93.4 140.2

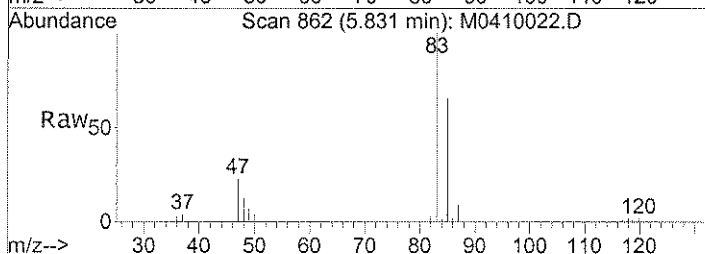


Abundance Ion 95.95 (95.65 to 96.65): M0  
 Ion 61.05 (60.75 to 61.75): M0

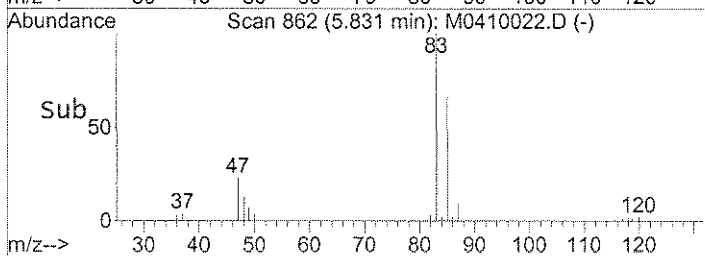
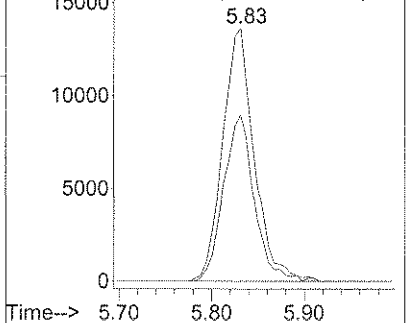


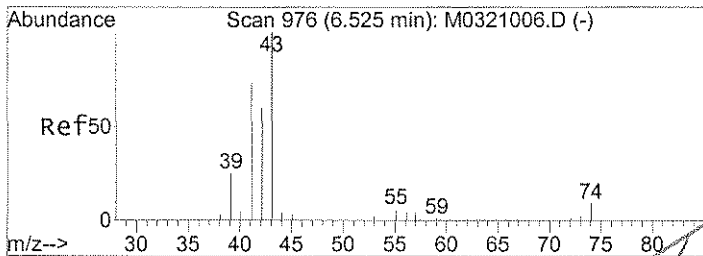
#31  
 Chloroform  
 Concen: 4.57 ug/l  
 RT: 5.83 min Scan# 862  
 Delta R.T. 0.00 min  
 Lab File: M0410022.D  
 Acq: 10 Apr 2007 15:32

Tgt Ion: 83 Resp: 32193  
 Ion Ratio Lower Upper  
 83 100  
 85 64.1 41.2 81.2



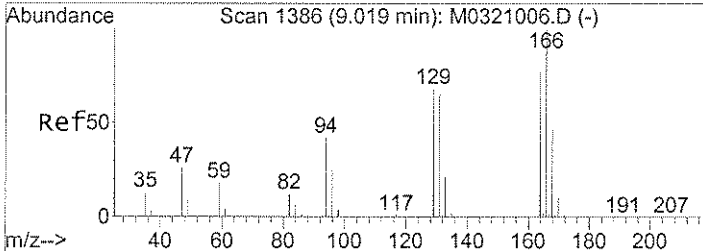
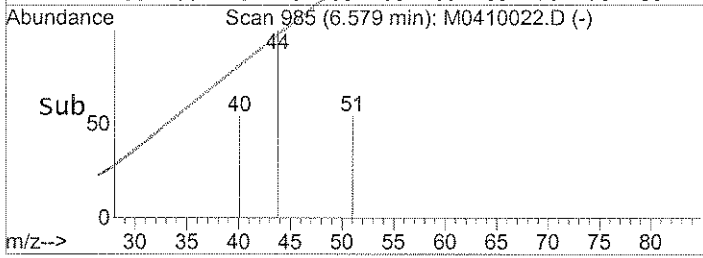
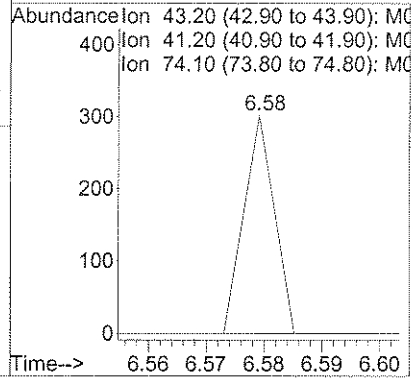
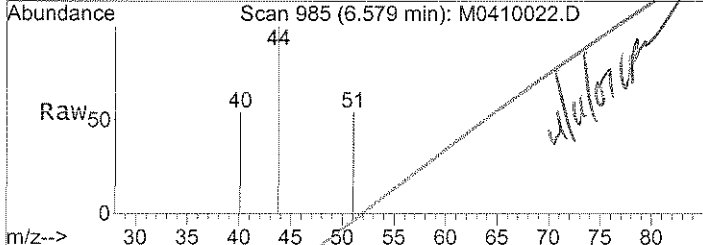
Abundance Ion 83.00 (82.70 to 83.70): M0  
 Ion 85.00 (84.70 to 85.70): M0





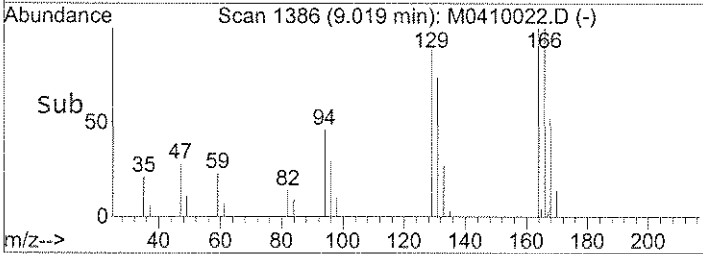
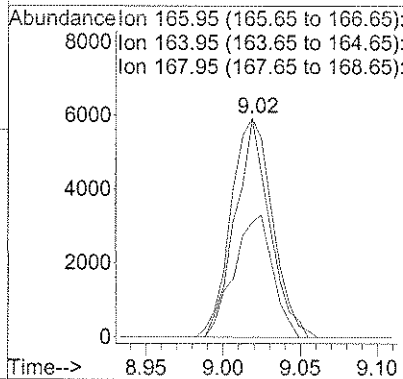
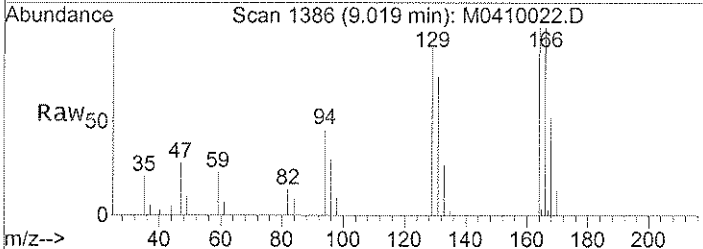
#40  
 Isobutanol  
 Concen: Below Cal  
 RT: 6.58 min Scan# 985  
 Delta R.T. 0.07 min  
 Lab File: M0410022.D  
 Acq: 10 Apr 2007 15:32

Tgt Ion	43	Resp	110
Ion Ratio	100	Lower	Upper
41	0.0	55.8	83.8#
74	0.0	8.0	12.0#



#56  
 Tetrachloroethene  
 Concen: 2.63 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. -0.00 min  
 Lab File: M0410022.D  
 Acq: 10 Apr 2007 15:32

Tgt Ion	166	Resp	11051
Ion Ratio	100	Lower	Upper
164	82.8	63.3	94.9
168	52.0	39.6	59.4



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410022.D                   Vial: 53  
Acq On    : 10 Apr 2007 15:32                   Operator: LPM  
Sample    : JPL28-002                         Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410022.D 524.M    wed Apr 11 08:32:08 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-003  
 Lab File ID: M0410023.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:55  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.88	
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	2.4	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.2	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-003  
 Lab File ID: M0410023.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:55  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	5.5	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-003  
 Lab File ID: M0410023.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 15:55  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-21-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL28  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-003  
 Lab File ID: M0410023.D  
 Date Collected: 03/28/2007  
 Date Analyzed: 04/10/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
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26				
27				
28				
29				
30				

Comments:



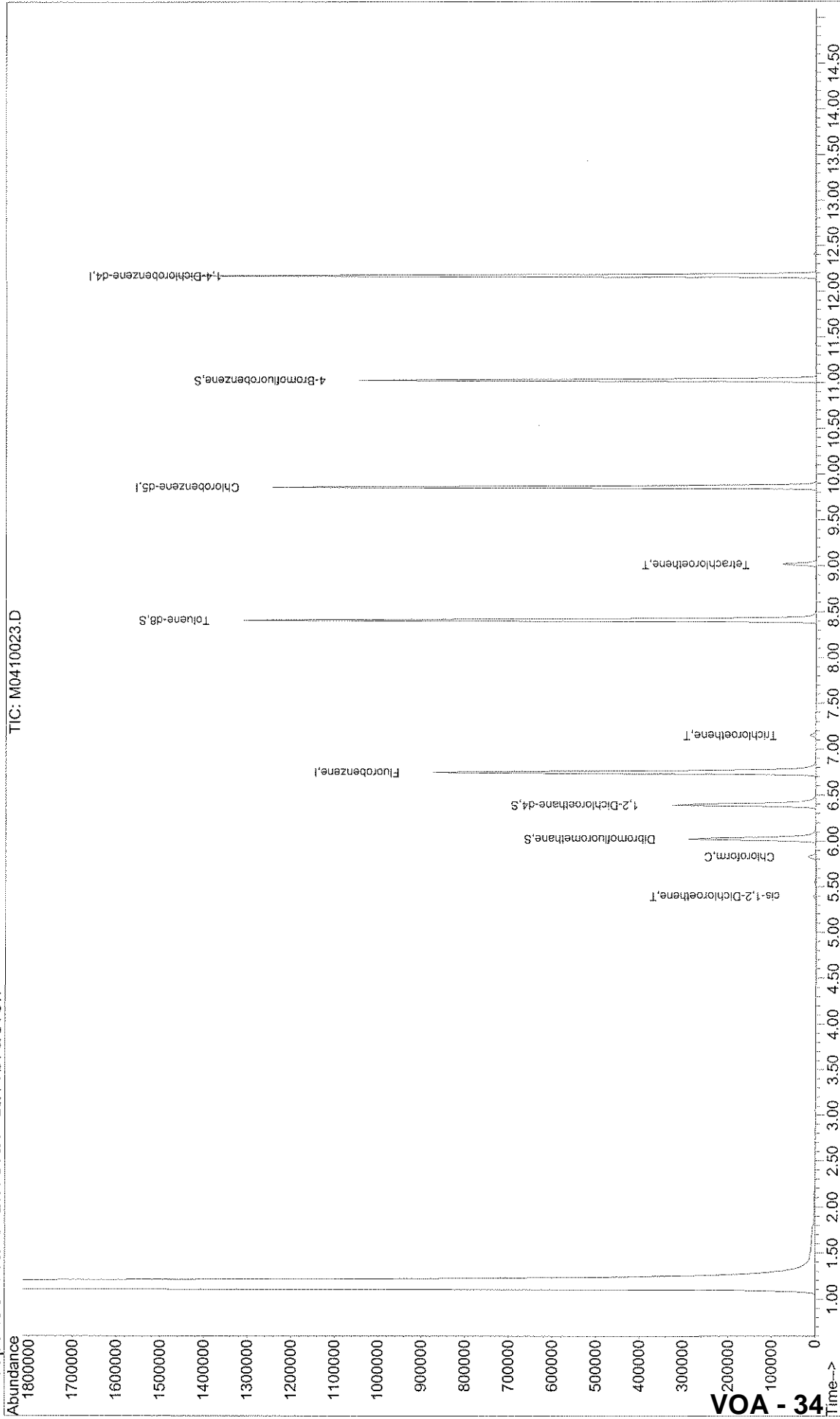
Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410023.D  
Acq On : 10 Apr 2007 15:55  
Sample : JPL28-003  
Misc : #2 5ml +IS/ss (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:26 2007

Vial: 53  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00

Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 34

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410023.D  
 Acq On : 10 Apr 2007 15:55  
 Sample : JPL28-003  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:26 2007

Vial: 53  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)

Title : VOA 524- 5ML Calibration 5973M

Last Update : Tue Apr 10 10:47:40 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	877670	50.00	ug/l	0.00 83.81%
50) Chlorobenzene-d5	9.86	82	368923	50.00	ug/l	0.00 81.44%
70) 1,4-Dichlorobenzene-d4	12.17	152	368972	50.00	ug/l	0.00 78.95%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	206226	49.38	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	98.76%		
37) 1,2-Dichloroethane-d4	6.39	65	248781	51.64	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	103.28%		
51) Toluene-d8	8.41	98	890487	51.93	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	103.86%		
72) 4-Bromofluorobenzene	11.03	95	306829	50.84	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery =	101.68%		

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.46	50	57	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.	d	
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	3.04	43	115	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.05	76	62	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.90	96	247	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	3.97	73	64	N.D.		
22) 1,1-Dichloroethane	4.58	63	1409	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.40	96	3389	0.88	ug/l	91
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.83	83	16592	2.37	ug/l	95
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	6.03	75	57	N.D.		
38) Benzene	6.41	78	74	N.D.		

*4/11/07*

Quantitation Report

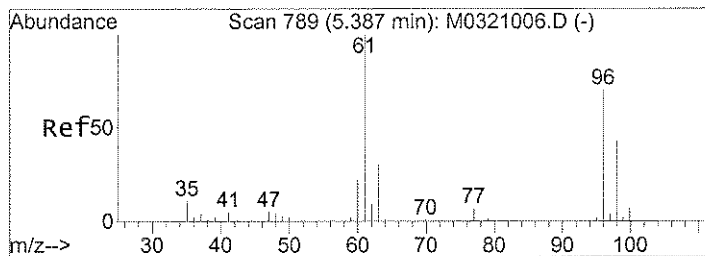
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 Acq On : 10 Apr 2007 15:55  
 Sample : JPL28-003  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:26 2007

Vial: 53  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

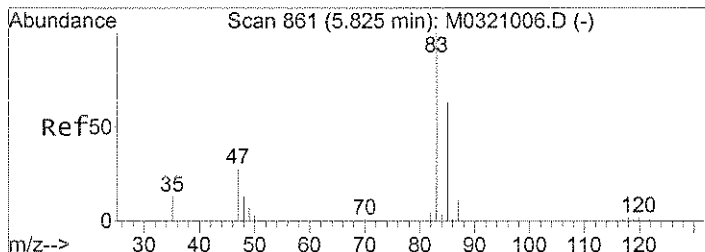
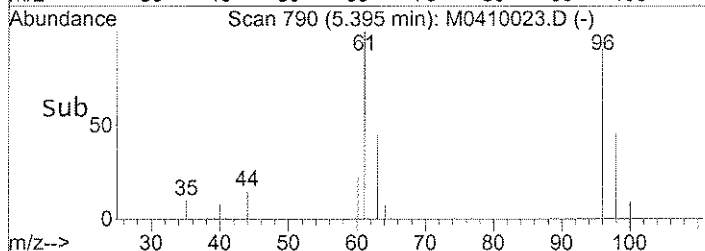
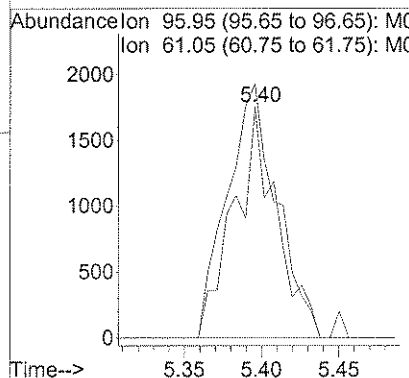
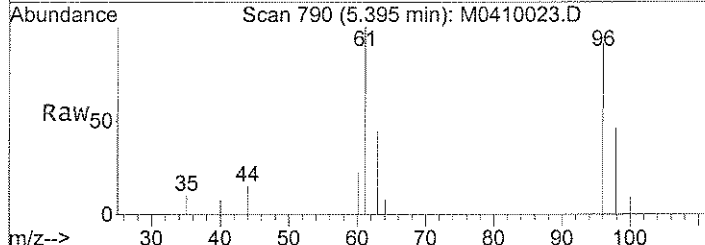
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	4845	1.16	ug/l	95
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.73	83	132		N.D.	
47) 2-Chloroethyl vinyl ether	8.11	63	57		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.47	92	295		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	23012	5.50	ug/l	98
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	9.05	43	70		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	242		N.D.	
62) 1-chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.98	91	822		N.D.	
65) m,p-xylene	10.09	106	885		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D. d	
69) Isopropylbenzene	10.86	105	78		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	293		N.D.	
75) 1,2,3-Trichloropropane	11.02	110	71		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.27	91	275		N.D.	
78) 4-Chlorotoluene	11.27	91	275		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D. d	
82) sec-butylbenzene	0.00	105	0		N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D. d	
84) 4-Isopropyltoluene	0.00	119	0		N.D. d	
85) 1,4-Dichlorobenzene	12.20	146	181		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	393		N.D.	
87) n-Butylbenzene	12.53	91	346		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	225		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D. d	
92) 1,2,3-Trichlorobenzene	14.66	180	472		N.D.	



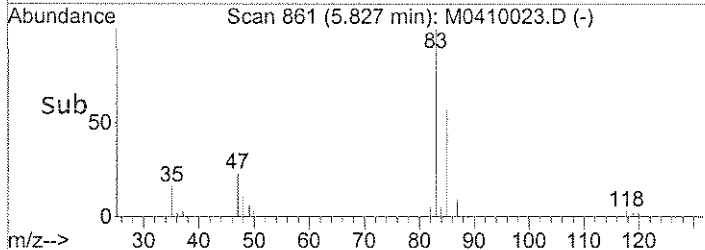
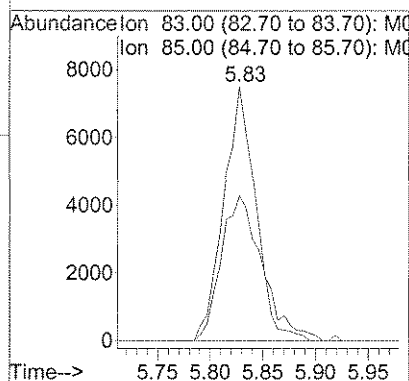
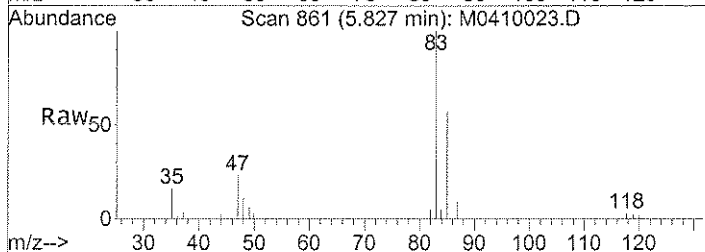
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 cis-1,2-Dichloroethene  
 Concen: 0.88 ug/l  
 RT: 5.40 min Scan# 790  
 Delta R.T. 0.01 min  
 Lab File: M0410023.D  
 Acq: 10 Apr 2007 15:55

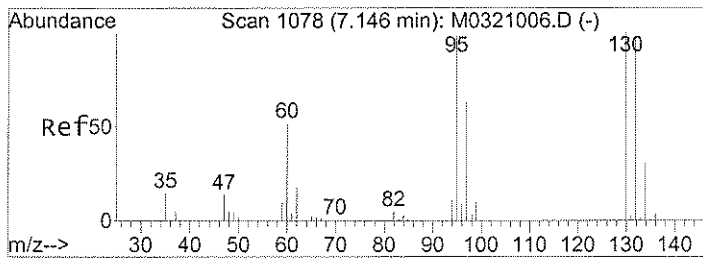
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 Ion Ratio Lower Upper  
 96 100  
 61 127.1 93.4 140.2



#31  
 Chloroform  
 Concen: 2.37 ug/l  
 RT: 5.83 min Scan# 861  
 Delta R.T. 0.00 min  
 Lab File: M0410023.D  
 Acq: 10 Apr 2007 15:55

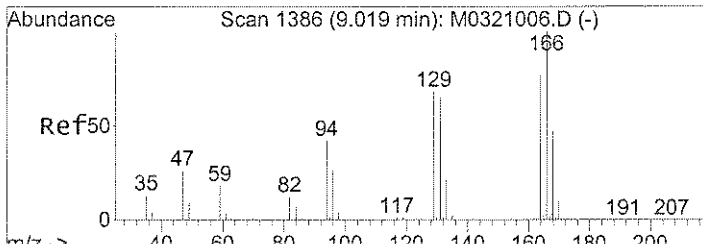
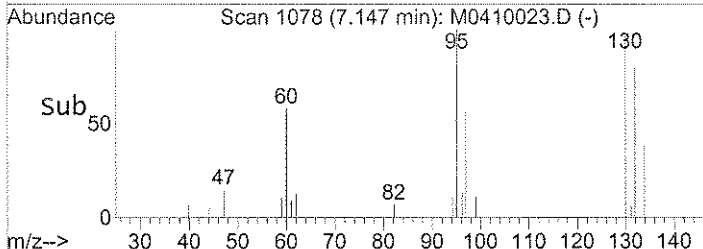
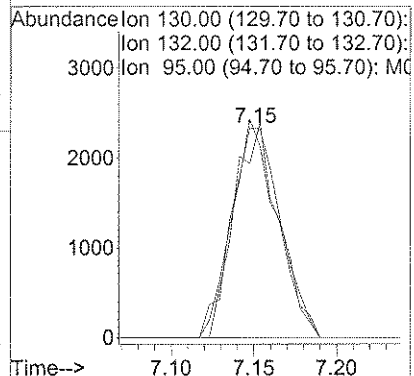
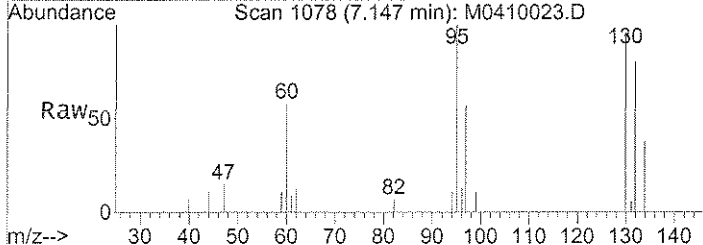
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 Ion Ratio Lower Upper  
 83 100  
 85 65.1 41.2 81.2





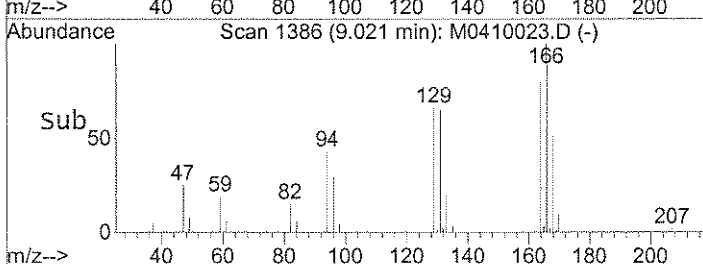
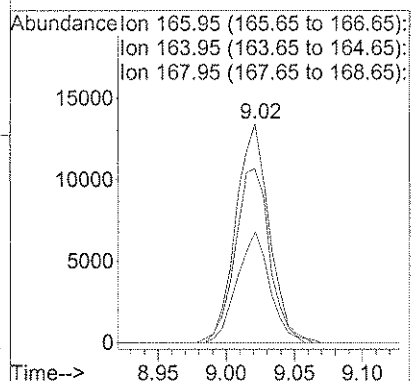
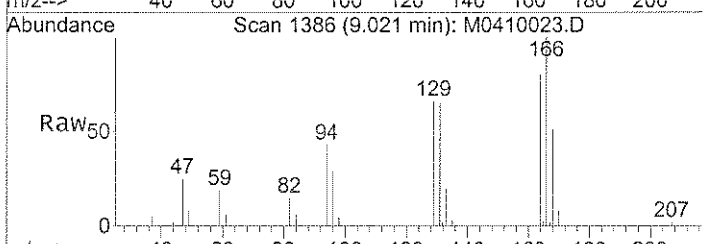
#41  
 Trichloroethene  
 Concen: 1.16 ug/l  
 RT: 7.15 min Scan# 1078  
 Delta R.T. 0.00 min  
 Lab File: M0410023.D  
 Acq: 10 Apr 2007 15:55

Tgt Ion	Resp	Lower	Upper
130	4845		
130	100		
132	91.5	75.0	115.0
95	94.9	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 5.50 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. 0.00 min  
 Lab File: M0410023.D  
 Acq: 10 Apr 2007 15:55

Tgt Ion	Resp	Lower	Upper
166	23012		
166	100		
164	81.3	63.3	94.9
168	49.1	39.6	59.4



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410023.D                   Vial: 53  
Acq On    : 10 Apr 2007  15:55                   Operator: LPM  
Sample    : JPL28-003                           Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410023.D  524.M    Wed Apr 11 08:32:15 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 16:17

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	1.6	
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	1.1	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.1	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZE-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-004  
 Lab File ID: M0410024.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 16:17  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	7.3	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-004  
 Lab File ID: M0410024.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 16:17  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-21-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL28  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-004  
 Lab File ID: M0410024.D  
 Date Collected: 03/28/2007  
 Date Analyzed: 04/10/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
10				
11				
12				
13				
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26				
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28				
29				
30				

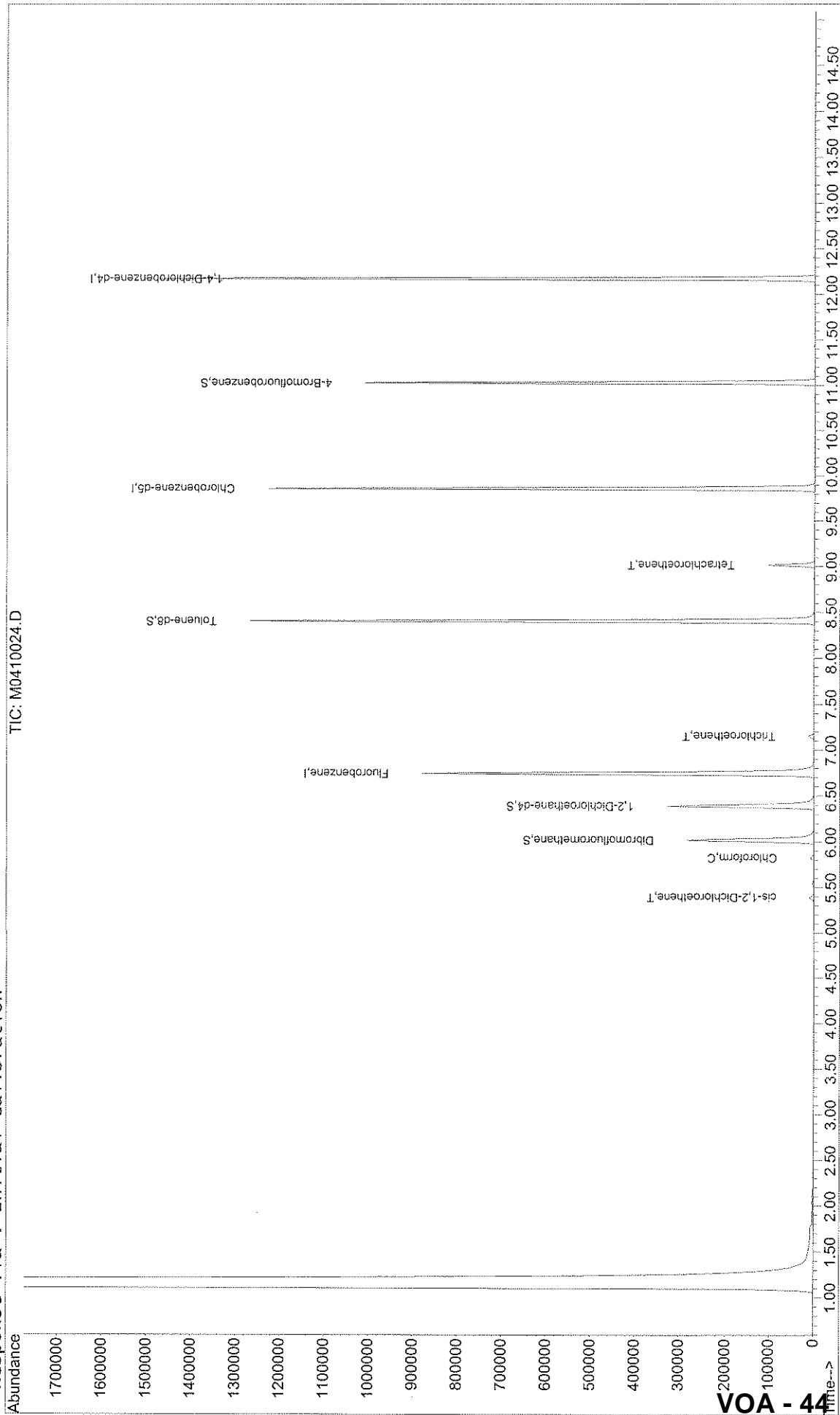
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410024.D  
Acq On : 10 Apr 2007 16:17  
Sample : JPL28-004  
Misc : #3 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:28 2007

Vial: 54  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 44

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410024.D  
 Acq On : 10 Apr 2007 16:17  
 Sample : JPL28-004  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:28 2007

Vial: 54  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	864755	50.00	ug/l	0.00 82.57%
50) Chlorobenzene-d5	9.86	82	365074	50.00	ug/l	0.00 80.59%
70) 1,4-Dichlorobenzene-d4	12.17	152	368668	50.00	ug/l	0.00 78.89%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	208187	50.60	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.20%
37) 1,2-Dichloroethane-d4	6.39	65	249317	52.53	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	105.06%
51) Toluene-d8	8.41	98	869976	51.27	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	102.54%
72) 4-Bromofluorobenzene	11.03	95	307979	51.07	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	102.14%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.45	50	64	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	3.02	43	85	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.05	76	64	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.90	96	119	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	3.92	73	743	N.D.		
22) 1,1-Dichloroethane	4.56	63	1206	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.39	96	5857	1.55	ug/l	98
27) 2-Butanone	5.55	43	74	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.82	83	7359	1.07	ug/l	95
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.41	78	222	N.D.		

*4/11/07*

Quantitation Report

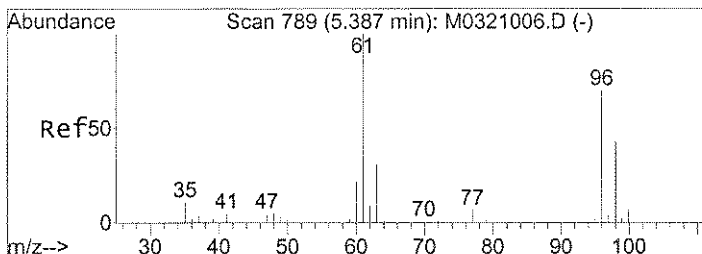
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 Sample : JPL28-004  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:28 2007

Vial: 54  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

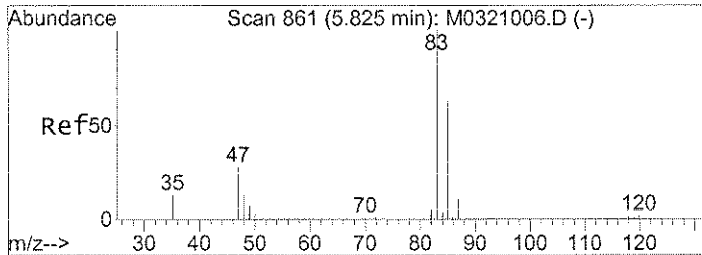
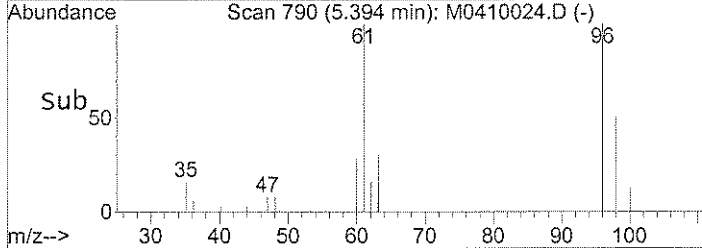
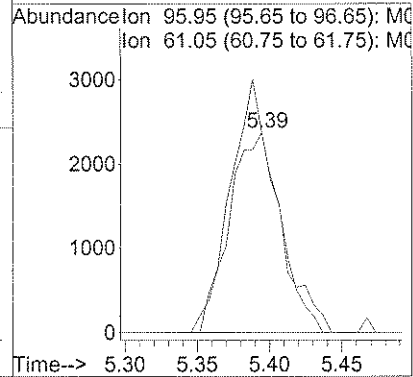
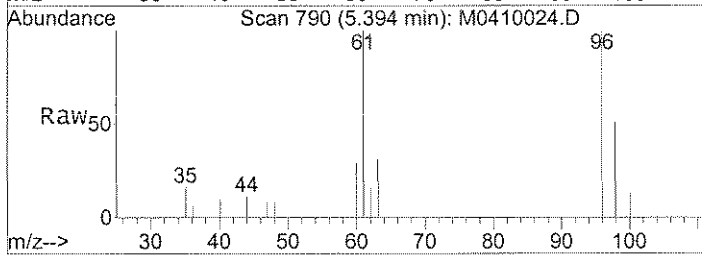
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	4399	1.07	ug/l	92
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	120		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	9.01	97	57		N.D.	
56) Tetrachloroethene	9.02	166	30010	7.25	ug/l	96
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	278		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	900		N.D.	
65) m,p-Xylene	10.10	106	506		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	223		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	137		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-chlorotoluene	11.36	91	134		N.D.	
78) 4-Chlorotoluene	11.47	91	55		N.D.	
79) 1,3,5-Trimethylbenzene	11.43	105	55		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D. d	
84) 4-Isopropyltoluene	0.00	119	0		N.D. d	
85) 1,4-Dichlorobenzene	12.20	146	149		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	260		N.D.	
87) n-Butylbenzene	12.53	91	279		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	145		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D. d	
92) 1,2,3-Trichlorobenzene	14.65	180	126		N.D.	



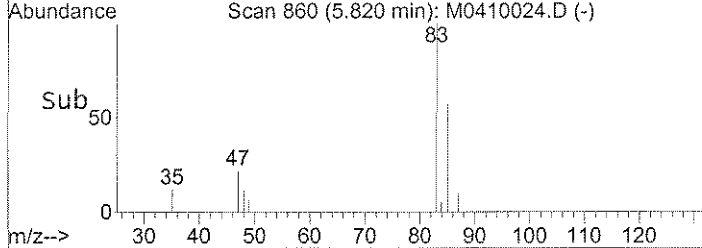
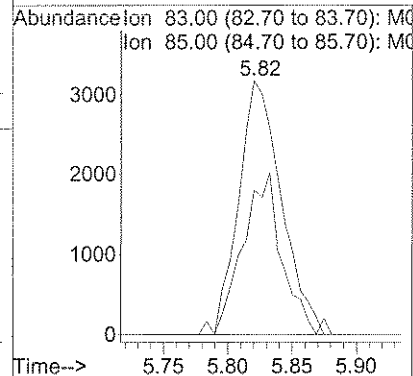
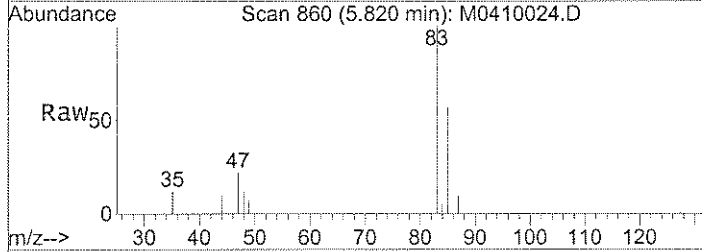
#26  
 cis-1,2-Dichloroethene  
 Concen: 1.55 ug/l  
 RT: 5.39 min Scan# 790  
 Delta R.T. 0.01 min  
 Lab File: M0410024.D  
 Acq: 10 Apr 2007 16:17

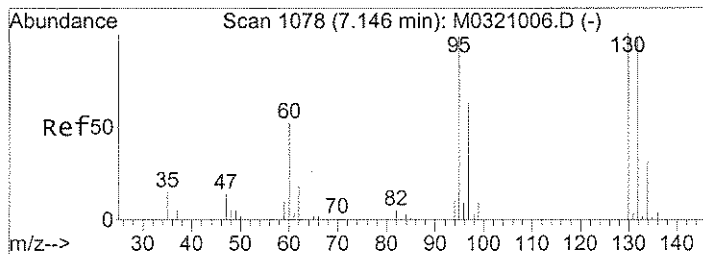
Tgt Ion	Resp	Lower	Upper
96	5857		
96	100		
61	114.8	93.4	140.2



#31  
 Chloroform  
 Concen: 1.07 ug/l  
 RT: 5.82 min Scan# 860  
 Delta R.T. -0.01 min  
 Lab File: M0410024.D  
 Acq: 10 Apr 2007 16:17

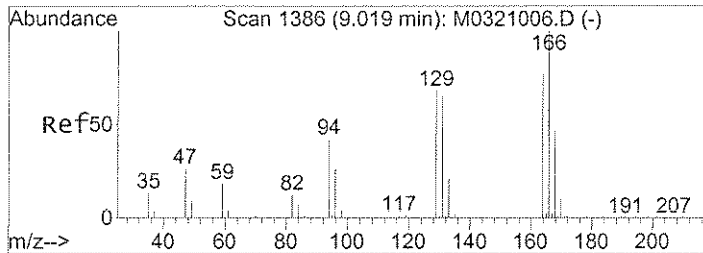
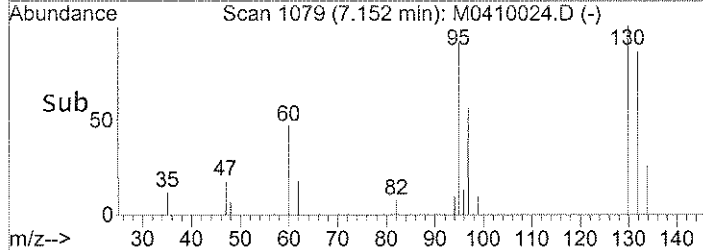
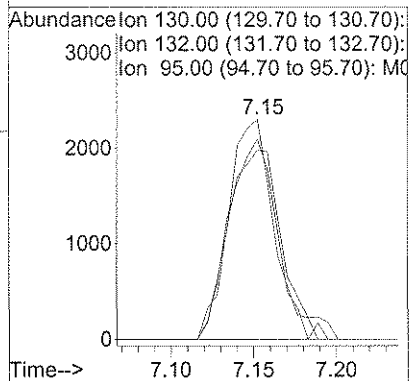
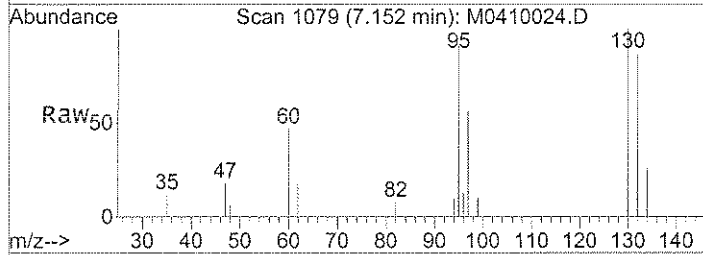
Tgt Ion	Resp	Lower	Upper
83	7359		
83	100		
85	57.1	41.2	81.2





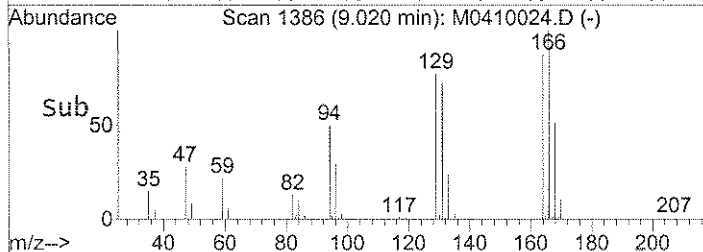
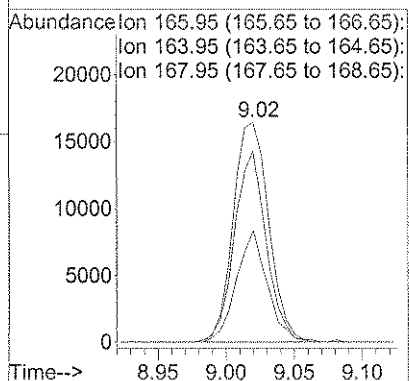
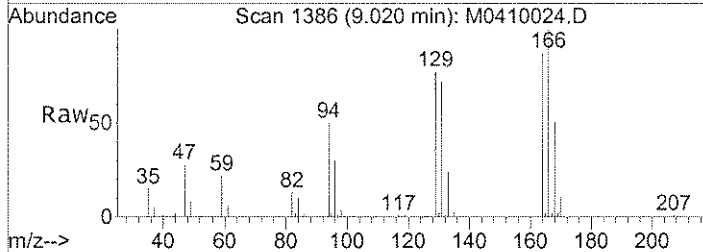
#41  
 Trichloroethene  
 Concen: 1.07 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0410024.D  
 Acq: 10 Apr 2007 16:17

Tgt Ion	Resp	Lower	Upper
130	4399		
130	100		
132	103.1	75.0	115.0
95	96.0	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 7.25 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. 0.00 min  
 Lab File: M0410024.D  
 Acq: 10 Apr 2007 16:17

Tgt Ion	Resp	Lower	Upper
166	30010		
166	100		
164	77.6	63.3	94.9
168	44.2	39.6	59.4



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410024.D                   Vial: 54  
Acq On    : 10 Apr 2007 16:17                   Operator: LPM  
Sample    : JPL28-004                           Inst     : MOBY  
Misc      : #3 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410024.D 524.M    wed Apr 11 08:32:23 2007



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 16:39

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.47	J
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.43	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 16:39

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.33	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-21-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-005  
 Lab File ID: M0410025.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 16:39  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-21-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL28  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-005  
 Lab File ID: M0410025.D  
 Date Collected: 03/28/2007  
 Date Analyzed: 04/10/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

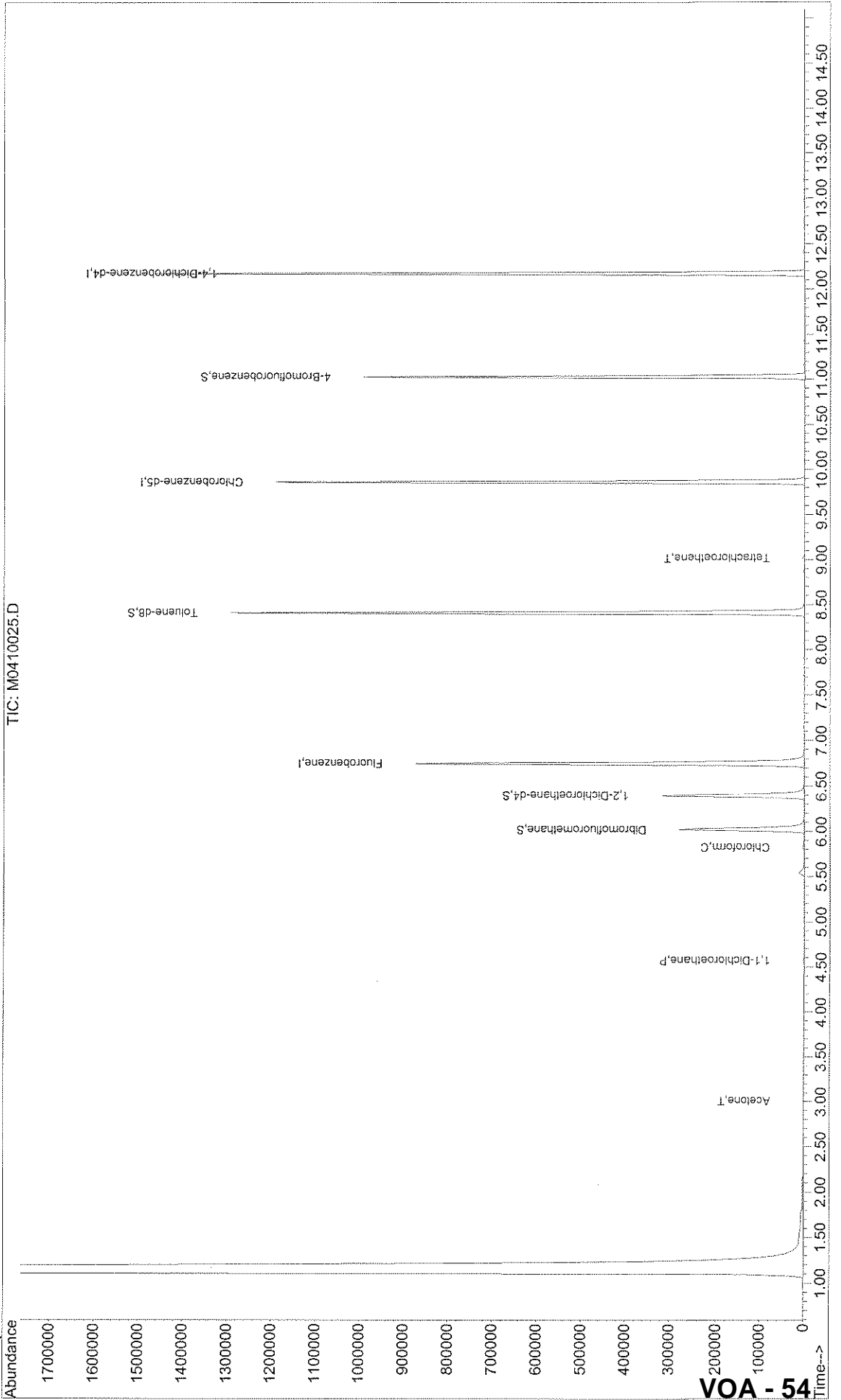
CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
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11				
12				
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29				
30				

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410025.D  
Acq On : 10 Apr 2007 16:39 Vial: 55  
Sample : JPL28-005 Operator: LPM  
Misc : #1 5ml +IS/SS (524) Inst : MOBY  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 11 7:30 2007 Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 54

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410025.D  
 Acq On : 10 Apr 2007 16:39  
 Sample : JPL28-005  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:30 2007

Vial: 55  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.75	96	858089	50.00	ug/l	0.00 81.94%
50) Chlorobenzene-d5	9.86	82	358237	50.00	ug/l	0.00 79.08%
70) 1,4-Dichlorobenzene-d4	12.18	152	364547	50.00	ug/l	0.00 78.01%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	204737	50.15	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.30%
37) 1,2-Dichloroethane-d4	6.40	65	246329	52.30	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	104.60%
51) Toluene-d8	8.41	98	866986	52.06	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.12%
72) 4-Bromofluorobenzene	11.03	95	302111	50.67	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	101.34%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.46	50	67	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D. d		
11) Acetone	3.01	43	721	0.79	ug/l #	62
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.06	76	114	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D. d		
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.57	63	2926	0.47	ug/l	87
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	0.00	43	0	N.D. d		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	5.87	41	324	N.D.		
31) Chloroform	5.82	83	2959	0.43	ug/l	93
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	6.18	75	72	N.D.		
38) Benzene	6.40	78	55	N.D.		

*4/11/07*

Quantitation Report

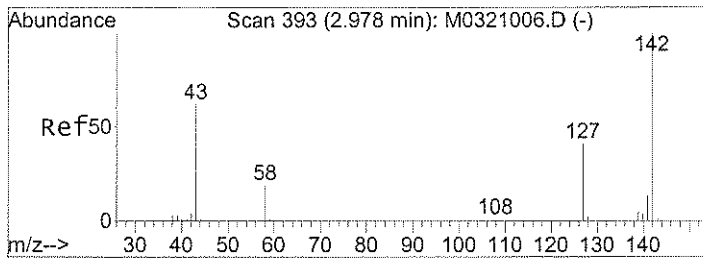
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 Acq On : 10 Apr 2007 16:39  
 Sample : JPL28-005  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:30 2007

Vial: 55  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

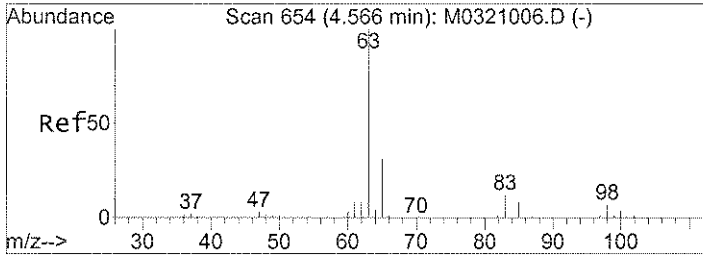
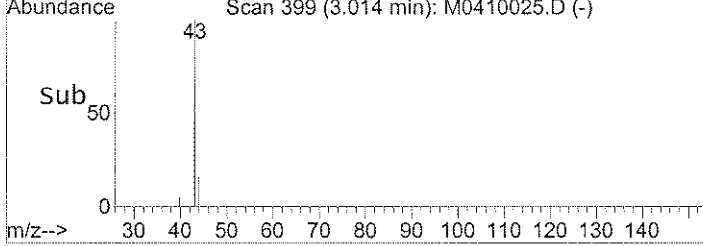
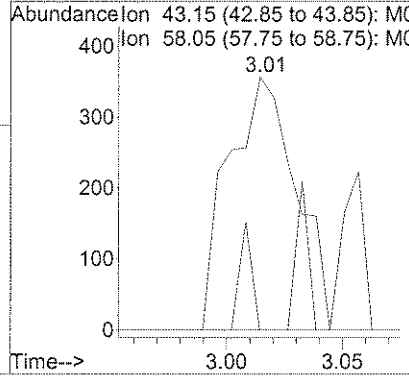
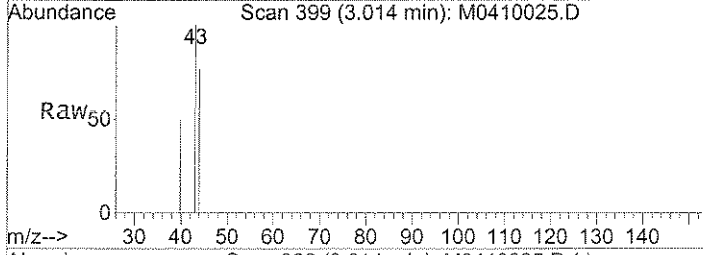
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	711		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.71	83	58		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	56		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.01	166	1358	0.33	ug/l	97
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	55		N.D.	
62) 1-chlorohexane	9.86	91	1219		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.98	91	234		N.D.	
65) m,p-Xylene	10.10	106	541		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	d
69) Isopropylbenzene	11.03	105	250		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	115		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	77		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.26	91	174		N.D.	
78) 4-Chlorotoluene	11.26	91	174		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	d
82) sec-butylbenzene	0.00	105	0		N.D.	d
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	12.54	91	57		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	14.66	180	121		N.D.	



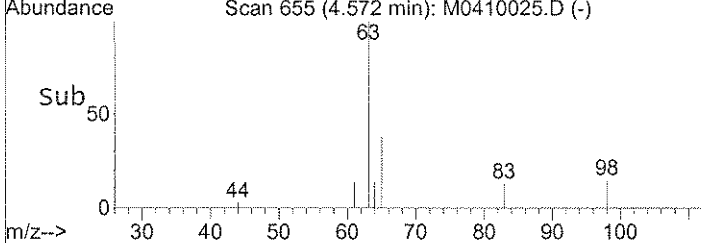
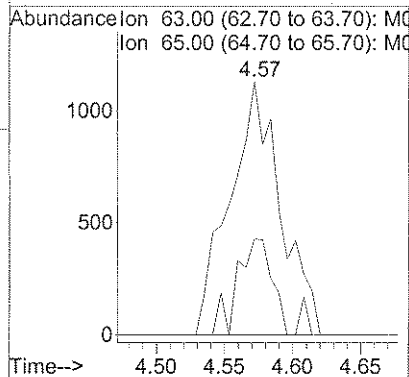
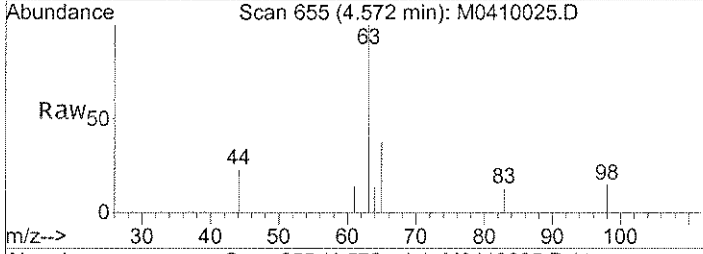
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 Acetone  
 Concen: 0.79 ug/l  
 RT: 3.01 min Scan# 399  
 Delta R.T. 0.05 min  
 Lab File: M0410025.D  
 Acq: 10 Apr 2007 16:39

Tgt Ion: 43 Resp: 721  
 Ion Ratio Lower Upper  
 43 100  
 58 7.8 21.9 32.9#

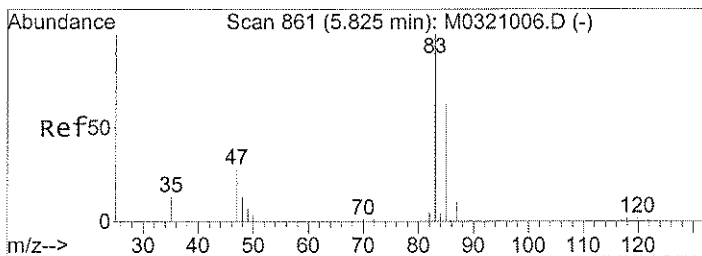


#22  
 1,1-Dichloroethane  
 Concen: 0.47 ug/l  
 RT: 4.57 min Scan# 655  
 Delta R.T. 0.01 min  
 Lab File: M0410025.D  
 Acq: 10 Apr 2007 16:39

Tgt Ion: 63 Resp: 2926  
 Ion Ratio Lower Upper  
 63 100  
 65 26.4 13.7 53.7

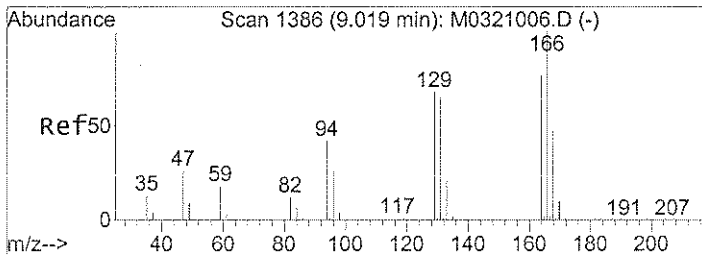
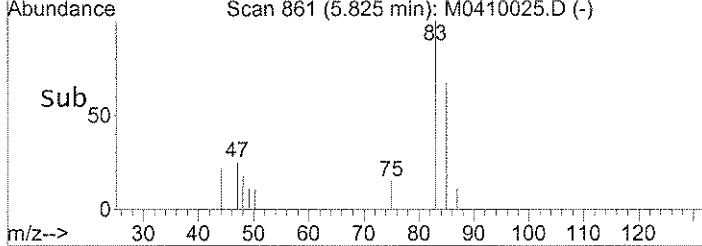
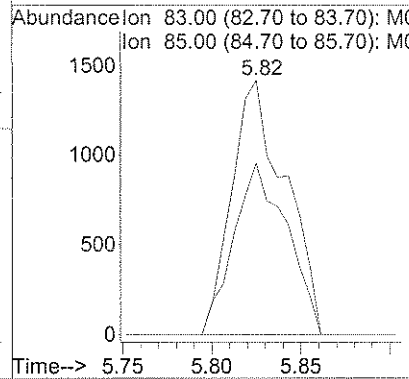
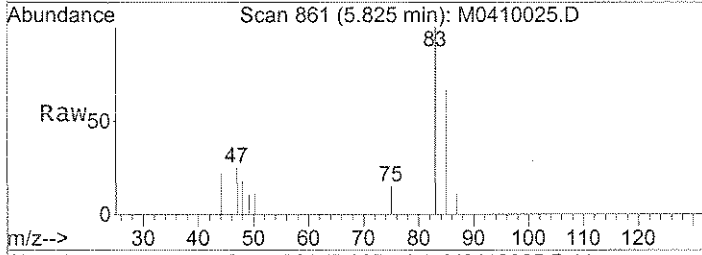






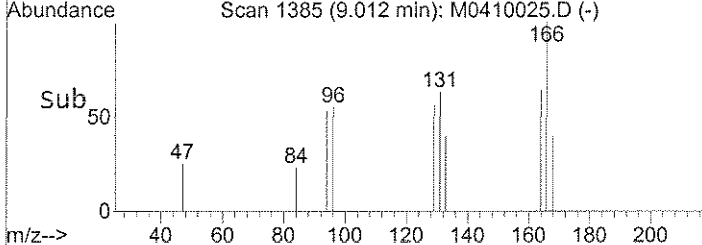
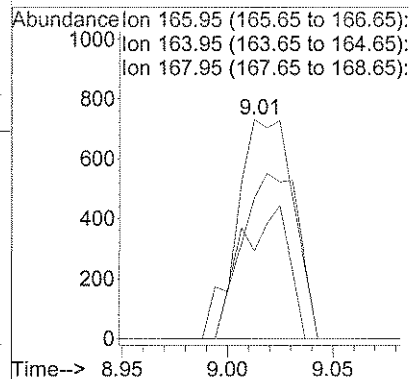
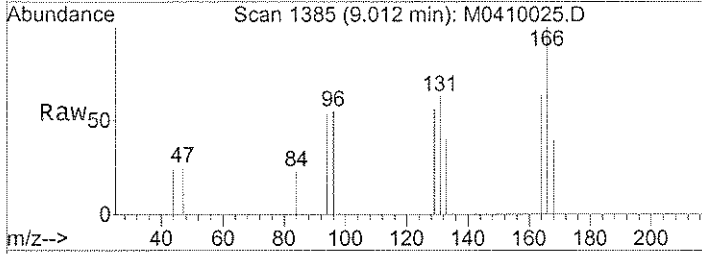
#31  
 Chloroform  
 Concen: 0.43 ug/l  
 RT: 5.82 min Scan# 861  
 Delta R.T. -0.00 min  
 Lab File: M0410025.D  
 Acq: 10 Apr 2007 16:39

Tgt Ion: 83 Resp: 2959  
 Ion Ratio Lower Upper  
 83 100  
 85 66.8 41.2 81.2



#56  
 Tetrachloroethene  
 Concen: 0.33 ug/l  
 RT: 9.01 min Scan# 1385  
 Delta R.T. -0.01 min  
 Lab File: M0410025.D  
 Acq: 10 Apr 2007 16:39

Tgt Ion: 166 Resp: 1358  
 Ion Ratio Lower Upper  
 166 100  
 164 75.3 63.3 94.9  
 168 50.9 39.6 59.4



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410025.D                   Vial: 55  
Acq On    : 10 Apr 2007 16:39                   Operator: LPM  
Sample    : JPL28-005                         Inst     : MOBY  
Misc      : #1 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0410025.D 524.M    wed Apr 11 08:32:31 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 17:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-006  
 Lab File ID: M0410026.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 17:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-006  
 Lab File ID: M0410026.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 17:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL28-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/28/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/10/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
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30				

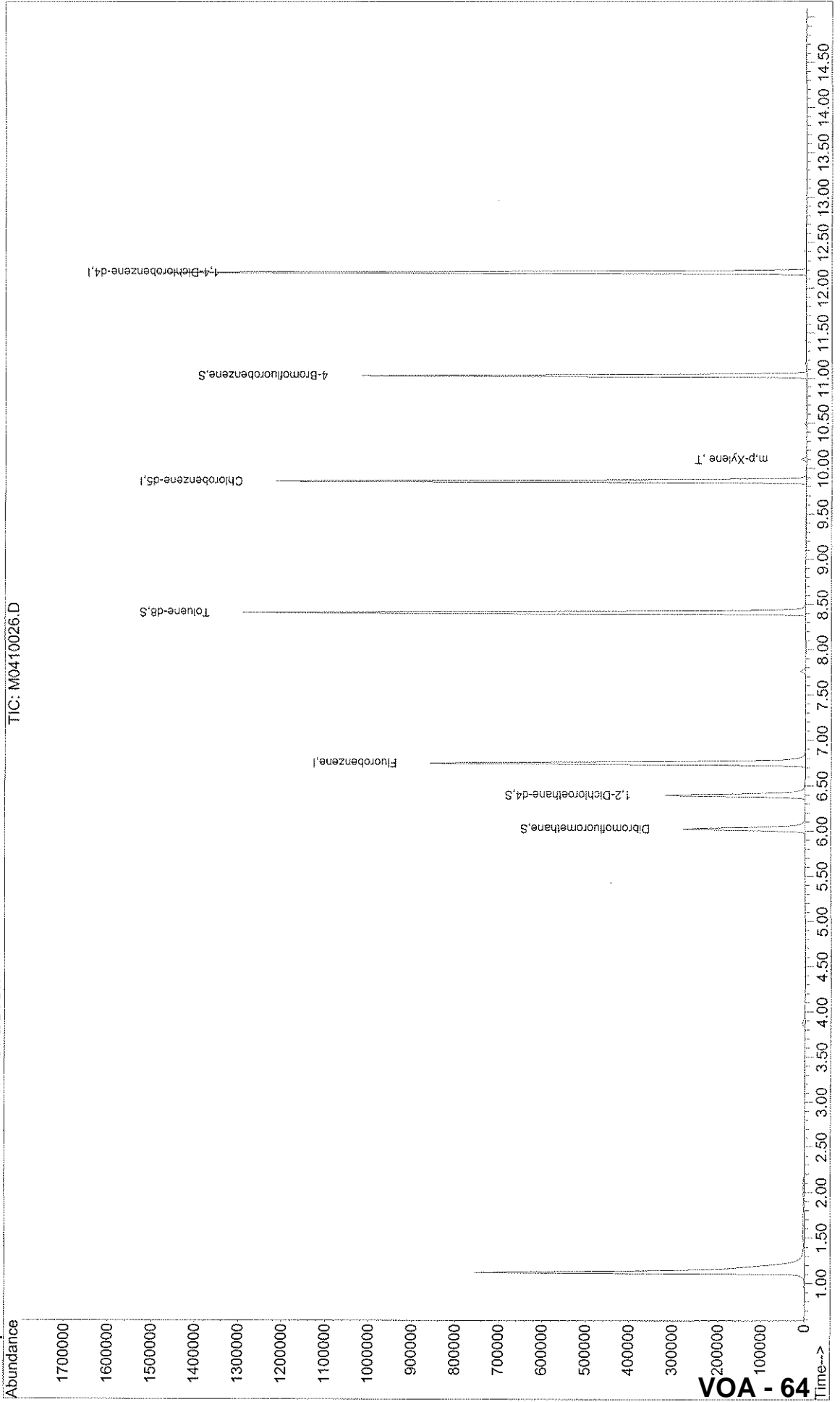
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410026.D  
Acq On : 10 Apr 2007 17:01  
Sample : JPL28-006  
Misc : #1 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:32 2007

Vial: 56  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 64

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410026.D  
 Acq On : 10 Apr 2007 17:01  
 Sample : JPL28-006  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:32 2007

Vial: 56  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.75	96	851008	50.00	ug/l	0.00 81.26%
50) Chlorobenzene-d5	9.86	82	362755	50.00	ug/l	0.00 80.08%
70) 1,4-Dichlorobenzene-d4	12.18	152	360550	50.00	ug/l	0.00 77.15%

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
33) Dibromofluoromethane	6.03	111	204152	50.42	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	100.84%	
37) 1,2-Dichloroethane-d4	6.40	65	245483	52.55	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	105.10%	
51) Toluene-d8	8.41	98	860893	51.05	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	102.10%	
72) 4-Bromofluorobenzene	11.03	95	303646	51.49	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery	=	102.98%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.06	76	60	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	3.87	53	56	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.55	43	178	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.82	83	125	N.D.		
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.	d	
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.40	78	56	N.D.		

*Handwritten signature*



Quantitation Report

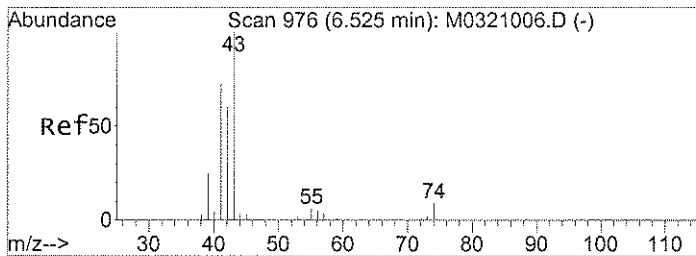
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 Acq On : 10 Apr 2007 17:01  
 Sample : JPL28-006  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:32 2007

Vial: 56  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

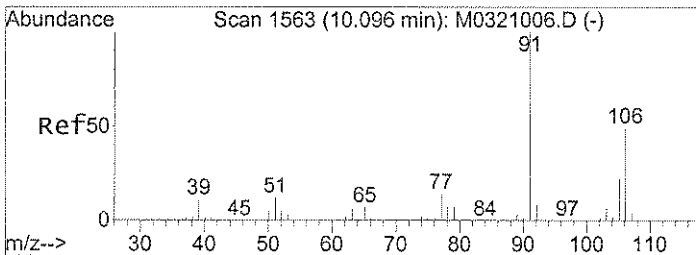
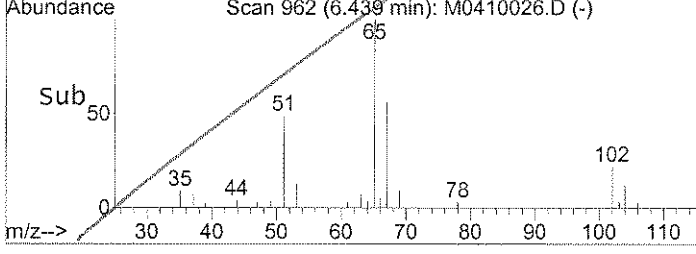
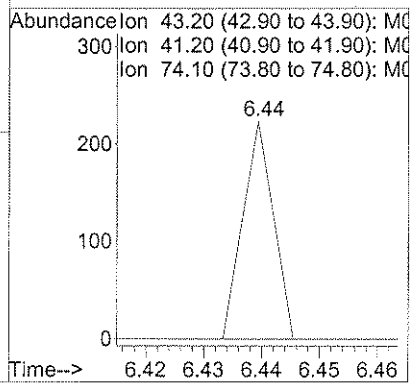
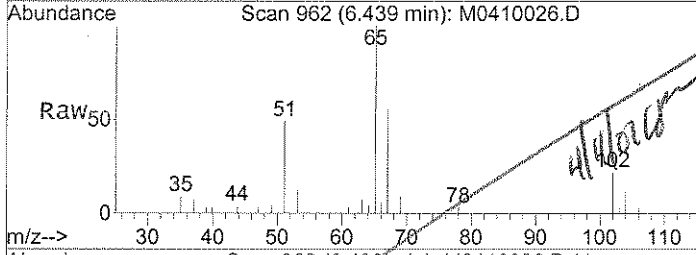
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	6.44	43	82	Below Cal #	22
41) Trichloroethene	7.24	130	64	N.D.	
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	8.48	92	179	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	0.00	166	0	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	8.98	43	412	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.90	112	57	N.D.	
62) 1-Chlorohexane	9.86	91	1435	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.99	91	2328	N.D.	
65) m,p-Xylene	10.10	106	3398	0.43 ug/l #	86
66) o-xylene	10.50	106	1264	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D. d	
69) Isopropylbenzene	10.86	105	122	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	0.00	156	0	N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	140	N.D.	
75) 1,2,3-Trichloropropane	11.03	110	210	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	
77) 2-Chlorotoluene	11.46	91	59	N.D.	
78) 4-Chlorotoluene	11.46	91	59	N.D.	
79) 1,3,5-Trimethylbenzene	11.20	105	67	N.D.	
80) tert-Butylbenzene	11.74	119	70	N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0	N.D. d	
82) sec-butylbenzene	0.00	105	0	N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0	N.D.	
84) 4-Isopropyltoluene	0.00	119	0	N.D. d	
85) 1,4-Dichlorobenzene	0.00	146	0	N.D.	
86) 1,2-Dichlorobenzene	12.57	146	68	N.D.	
87) n-Butylbenzene	12.52	91	148	N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0	N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	63	N.D.	
90) Hexachlorobutadiene	0.00	225	0	N.D.	
91) Naphthalene	0.00	128	0	N.D. d	
92) 1,2,3-Trichlorobenzene	0.00	180	0	N.D.	



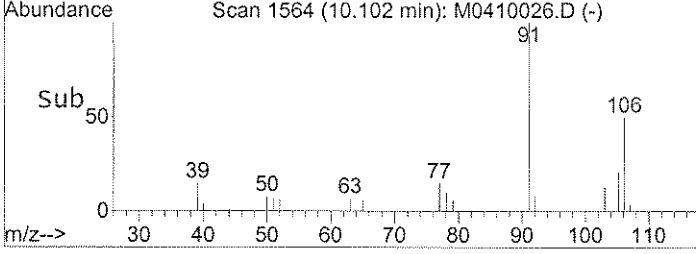
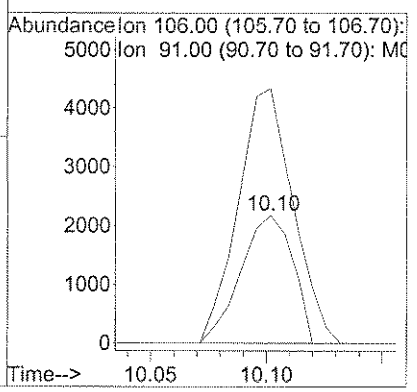
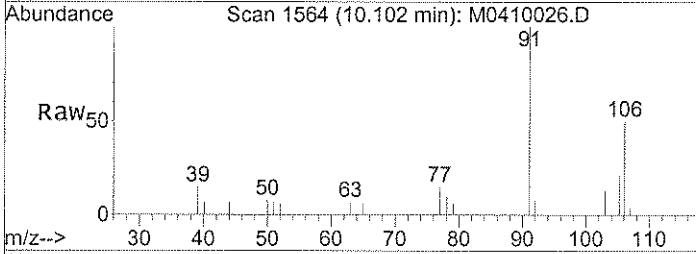
#40  
 Isobutanol  
 Concen: Below Cal  
 RT: 6.44 min Scan# 962  
 Delta R.T. -0.07 min  
 Lab File: M0410026.D  
 Acq: 10 Apr 2007 17:01

Tgt Ion	Resp	Lower	Upper
43	100		
41	0.0	55.8	83.8#
74	0.0	8.0	12.0#



#65  
 m,p-xylene  
 Concen: 0.43 ug/l  
 RT: 10.10 min Scan# 1564  
 Delta R.T. -0.00 min  
 Lab File: M0410026.D  
 Acq: 10 Apr 2007 17:01

Tgt Ion	Resp	Lower	Upper
106	100		
91	210.6	169.6	209.6#



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410026.D                   Vial: 56  
Acq On    : 10 Apr 2007 17:01                   Operator: LPM  
Sample    : JPL28-006                         Inst     : MOBY  
Misc      : #1 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410026.D 524.M    Wed Apr 11 08:32:38 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL28  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-007  
 Lab File ID: M0410020.D  
 Date Collected: 03/27/2007  
 Date/Time Analyzed: 04/10/2007 14:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 14:48

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	0.50		U
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	0.50		U
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL28

Run Sequence: R016728

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL28-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0410020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/27/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/10/2007 14:48

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-1-3/27/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL28  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016728  
 Lab Sample ID: JPL28-007  
 Lab File ID: M0410020.D  
 Date Collected: 03/28/2007  
 Date Analyzed: 04/10/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
10				
11				
12				
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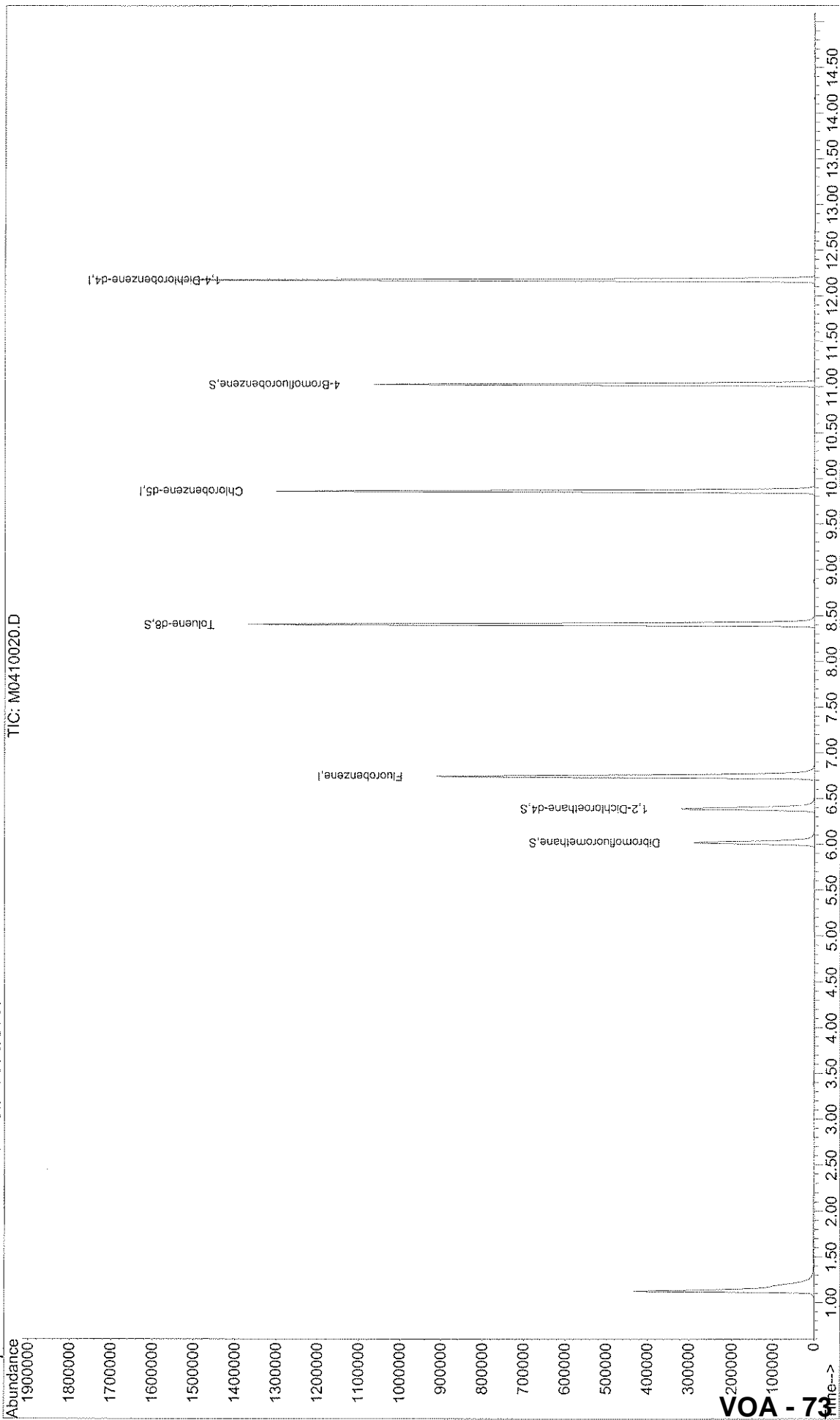
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410020.D  
Acq On : 10 Apr 2007 14:48  
Sample : JPL28-007  
Misc : #1 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 11 7:23 2007

Vial: 57  
Operator: LPM  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 73



Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410020.D  
 Acq On : 10 Apr 2007 14:48  
 Sample : JPL28-007  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:23 2007

Vial: 57  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar)
1) Fluorobenzene	6.75	96	905574	50.00	ug/l	0.00 86.47%
50) Chlorobenzene-d5	9.86	82	379343	50.00	ug/l	0.00 83.74%
70) 1,4-Dichlorobenzene-d4	12.17	152	382691	50.00	ug/l	0.00 81.89%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	215508	50.02	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.04%
37) 1,2-Dichloroethane-d4	6.39	65	258150	51.94	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	103.88%
51) Toluene-d8	8.41	98	909739	51.59	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.18%
72) 4-Bromofluorobenzene	11.03	95	319778	51.09	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	102.18%

Target Compounds

				Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.
3) Chloromethane	0.00	50	0	N.D.
4) Vinyl chloride	0.00	62	0	N.D.
5) Bromomethane	0.00	96	0	N.D.
6) Chloroethane	0.00	64	0	N.D.
7) Trichlorofluoromethane	0.00	101	0	N.D.
8) Acrolein	0.00	56	0	N.D.
9) 1,1-Dichloroethene	0.00	96	0	N.D.
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D. d
11) Acetone	0.00	43	0	N.D. d
12) Iodomethane	0.00	142	0	N.D.
13) Bromoethane	0.00	108	0	N.D.
14) Carbon Disulfide	3.05	76	587	N.D.
15) Allyl chloride	0.00	76	0	N.D.
16) Acetonitrile	0.00	40	0	N.D. d
17) Methyl Acetate	0.00	43	0	N.D.
18) Methylene Chloride	0.00	84	0	N.D. d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.
20) Acrylonitrile	0.00	53	0	N.D.
21) Methyl tert-butyl ether	0.00	73	0	N.D.
22) 1,1-Dichloroethane	0.00	63	0	N.D.
23) Vinyl acetate	0.00	43	0	N.D.
24) Chloroprene	0.00	53	0	N.D.
25) 2,2-Dichloropropane	0.00	77	0	N.D.
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.
27) 2-Butanone	0.00	43	0	N.D.
28) Propionitrile	0.00	54	0	N.D.
29) Bromochloromethane	0.00	128	0	N.D.
30) Methacrylonitrile	0.00	41	0	N.D.
31) Chloroform	0.00	83	0	N.D.
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.
34) Cyclohexane	0.00	56	0	N.D.
35) Carbon Tetrachloride	0.00	117	0	N.D.
36) 1,1-Dichloropropene	0.00	75	0	N.D.
38) Benzene	6.41	78	62	N.D.

*4/11/07*

Quantitation Report

Data File : X:\MSVOA\MOBY\041007\M0410020.D  
 Acq On : 10 Apr 2007 14:48  
 Sample : JPL28-007  
 Misc : #1 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 11 7:23 2007

Vial: 57  
 Operator: LPM  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	d
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.49	92	226		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	69		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	158		N.D.	
62) 1-Chlorohexane	9.86	91	1326		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.98	91	136		N.D.	
65) m,p-Xylene	10.10	106	284		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	d
68) Bromoform	0.00	173	0		N.D.	d
69) Isopropylbenzene	10.85	105	242		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	58		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	173		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	11.43	120	58		N.D.	
77) 2-Chlorotoluene	11.36	91	323		N.D.	
78) 4-Chlorotoluene	11.48	91	310		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	129		N.D.	
80) tert-Butylbenzene	11.75	119	239		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	d
82) sec-butylbenzene	0.00	105	0		N.D.	d
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	d
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	504		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	382		N.D.	
87) n-Butylbenzene	12.52	91	814		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.16	180	572		N.D.	
90) Hexachlorobutadiene	14.31	225	188		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	14.65	180	846		N.D.	

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041007\M0410020.D                   Vial: 57  
Acq On    : 10 Apr 2007 14:48                   Operator: LPM  
Sample    : JPL28-007                         Inst     : MOBY  
Misc      : #1 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0410020.D 524.M    Wed Apr 11 08:31:52 2007

**Metals Data**

**JPL28**

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-21-5	JPL28-001
MW-21-4	JPL28-002
MW-21-3	JPL28-003
MW-21-2	JPL28-004
MW-21-1	JPL28-005
EB-1-3/27/07	JPL28-006
EB-1-3/27/07MS	JPL28-006MS
EB-1-3/27/07MSD	JPL28-006MSD

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

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I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Cheronne Oeirs

Date: 04/16/2007

Title: Metals Lead

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-21-5

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-001

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.78			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-21-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-002

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.39			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-21-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-003

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.36			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-21-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-004

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.04			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-21-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-005

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-1-3/27/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL28

Matrix (soil/water): Water

Lab Sample ID: JPL28-006

Level (low/med): LOW

Date Received: 03/28/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL28**

**COVER PAGE - INORGANIC ANALYSES DATA PACKAGE**

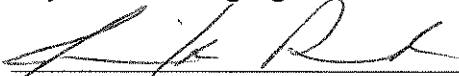
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL28

Client Identification	Lab Sample Work Order Number
MW-21-5	JPL28-001DL
MW-21-4	JPL28-002DL
MW-21-3	JPL28-003DL
MW-21-2	JPL28-004DL
MW-21-1	JPL28-005DL
EB-1-3/27/07	JPL28-006
EB-1-3/27/07MS	JPL28-006MS

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Penner

Date: 4-16-07

Title: Inorganics Lead

## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL28  
Sample Number: MW-21-5 Date/Time Collected: 03/27/2007 07:57  
Lab Sample ID: JPL28-001 Date/Time Received: 03/28/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/05/2007	04/06/2007	R016553



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL28  
**Sample Number:** MW-21-4 **Date/Time Collected:** 03/27/2007 08:32  
**Lab Sample ID:** JPL28-002 **Date/Time Received:** 03/28/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/05/2007	04/06/2007	R016553

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL28  
Sample Number: MW-21-3 Date/Time Collected: 03/27/2007 09:06  
Lab Sample ID: JPL28-003 Date/Time Received: 03/28/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/05/2007	04/06/2007	R016553

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL28  
**Sample Number:** MW-21-2 **Date/Time Collected:** 03/27/2007 09:35  
**Lab Sample ID:** JPL28-004 **Date/Time Received:** 03/28/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/05/2007	04/06/2007	R016553

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL28  
**Sample Number:** MW-21-1 **Date/Time Collected:** 03/27/2007 10:09  
**Lab Sample ID:** JPL28-005 **Date/Time Received:** 03/28/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/05/2007	04/06/2007	R016553

**Laucks Testing Laboratories, Inc.**

**Final Results**

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL28  
**Sample Number:** EB-1-3/27/07 **Date/Time Collected:** 03/27/2007 09:59  
**Lab Sample ID:** JPL28-006 **Date/Time Received:** 03/28/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/05/2007	04/06/2007	R016553

**SAMPLE DATA**

SDG JPL29

VOLATILES ANALYSIS

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-001  
 Lab File ID: M0411016.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-001  
 Lab File ID: M0411016.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-001  
 Lab File ID: M0411016.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

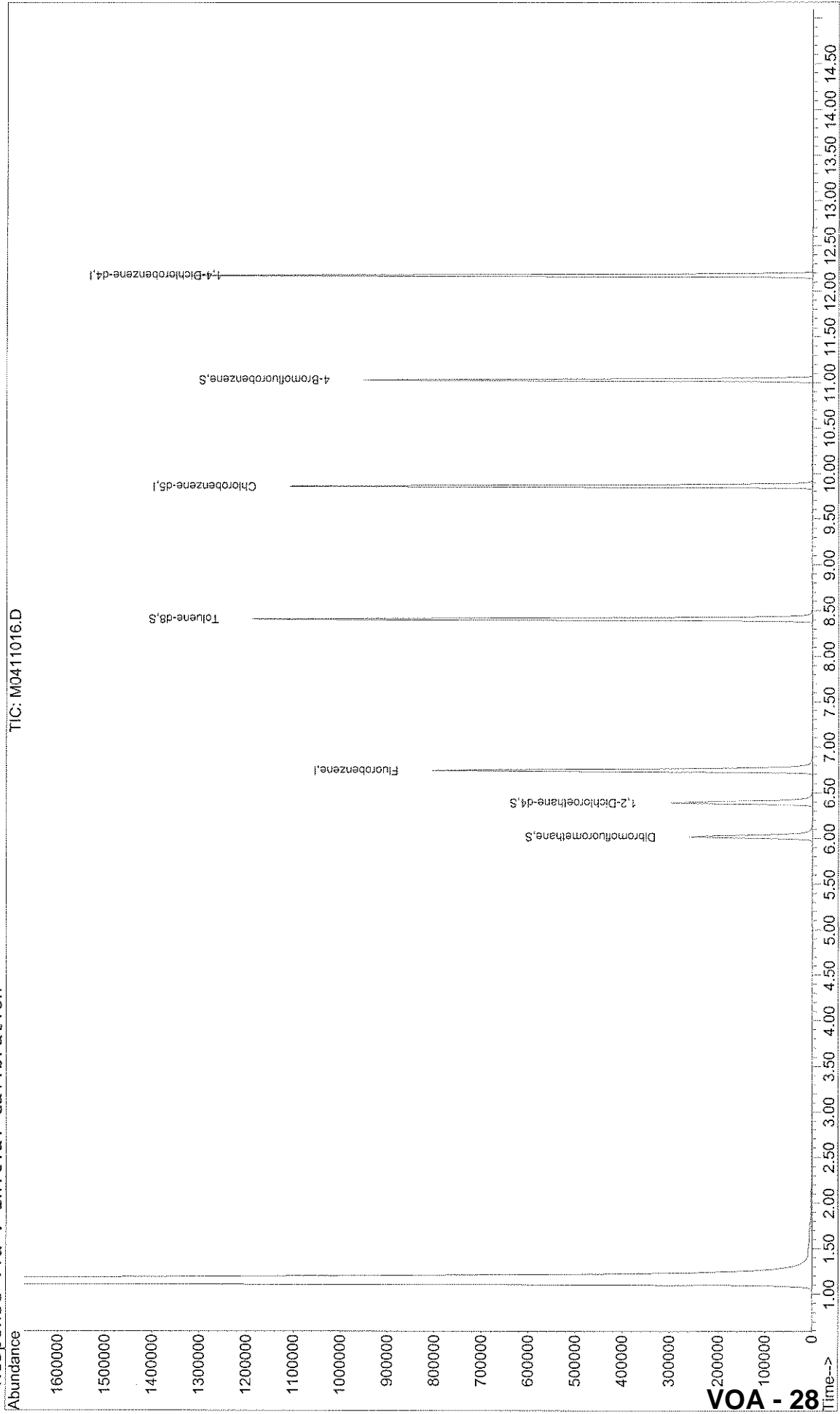
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411016.D  
Acq On : 11 Apr 2007 13:25  
Sample : JPL29-001  
Misc : #3 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:11 2007

Vial: 56  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 28

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411016.D  
 Acq On : 11 Apr 2007 13:25  
 Sample : JPL29-001  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:11 2007

Vial: 56  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)

Title : VOA 524- 5ML Calibration 5973M

Last Update : Tue Apr 10 10:47:40 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	788319	50.00	ug/l	0.00 75.27%
50) Chlorobenzene-d5	9.86	82	332222	50.00	ug/l	0.00 73.34%
70) 1,4-Dichlorobenzene-d4	12.18	152	330534	50.00	ug/l	0.00 70.73%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	187795	50.07	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	100.14%		
37) 1,2-Dichloroethane-d4	6.39	65	229384	53.01	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	106.02%		
51) Toluene-d8	8.41	98	797702	51.65	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	103.30%		
72) 4-Bromofluorobenzene	11.03	95	279979	51.79	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery =	103.58%		

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	3.01	43	56	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.06	76	197	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	3.49	84	57	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	0.00	83	0	N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	0.00	78	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0411016.D 524.M Thu Apr 12 09:12:03 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411016.D  
 Acq On : 11 Apr 2007 13:25  
 Sample : JPL29-001  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:11 2007

Vial: 56  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	141		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.86	91	1238		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	910		N.D.	
65) m,p-Xylene	10.10	106	425		N.D.	
66) o-xylene	10.49	106	58		N.D.	
67) Styrene	0.00	104	0		N.D.	d
68) Bromoform	0.00	173	0		N.D.	d
69) Isopropylbenzene	11.03	105	159		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	75		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	340		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	184		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.76	119	62		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	d
83) 1,3-Dichlorobenzene	12.19	146	150		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	12.19	146	150		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	12.52	91	63		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	14.66	180	83		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-002  
 Lab File ID: M0411017.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:47  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-002  
 Lab File ID: M0411017.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:47  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-002  
 Lab File ID: M0411017.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 13:47  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

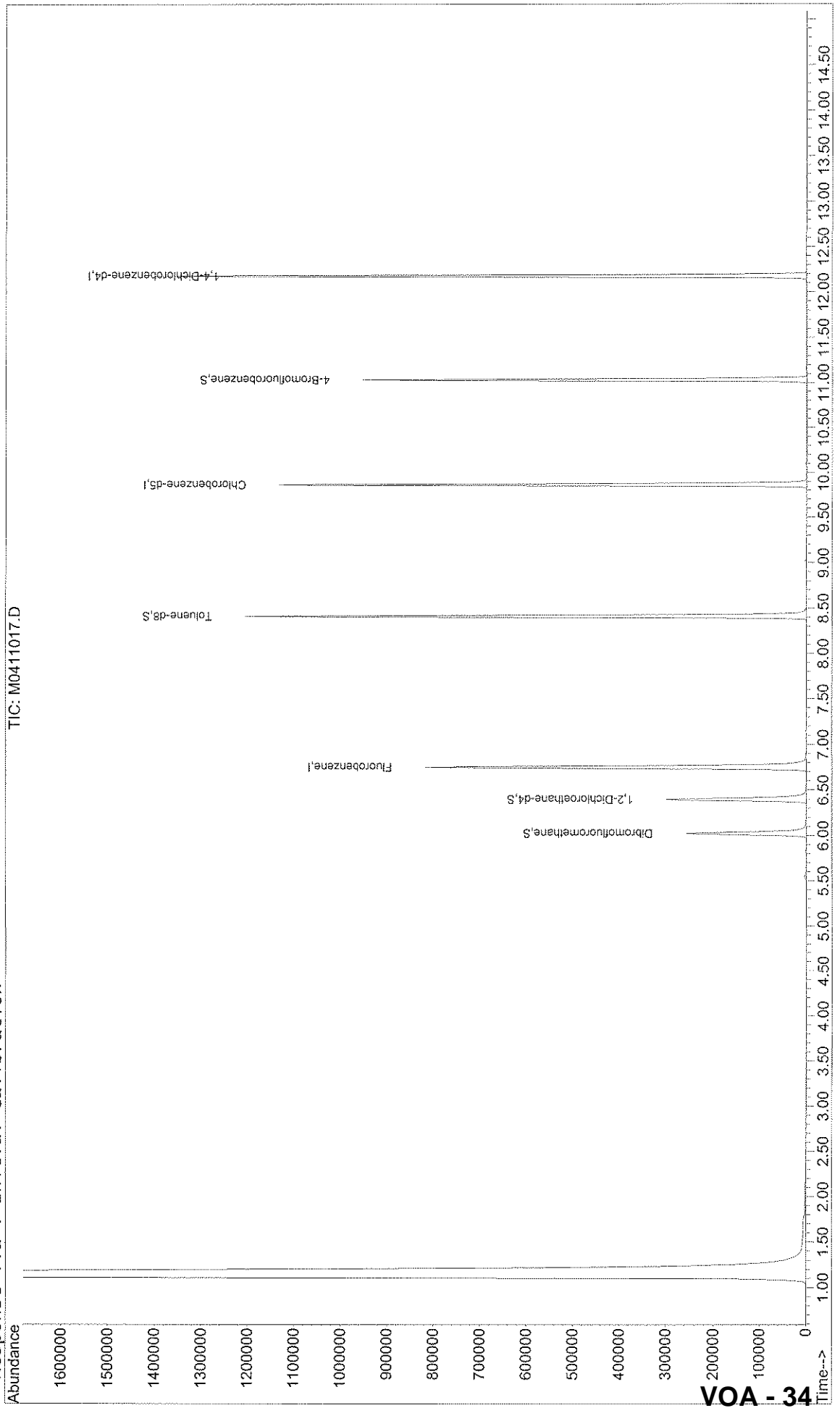
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411017.D  
Acq On : 11 Apr 2007 13:47  
Sample : JPL29-002  
Misc : #2 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:12 2007  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 34



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411017.D  
 Acq On : 11 Apr 2007 13:47  
 Sample : JPL29-002  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:12 2007

Vial: 57  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)

Title : VOA 524- 5ML Calibration 5973M

Last Update : Tue Apr 10 10:47:40 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	785115	50.00	ug/l	0.00 74.97%
50) Chlorobenzene-d5	9.86	82	329797	50.00	ug/l	0.00 72.80%
70) 1,4-Dichlorobenzene-d4	12.18	152	332435	50.00	ug/l	0.00 71.14%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	183737	49.18	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	98.36%
37) 1,2-Dichloroethane-d4	6.40	65	227922	52.89	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	105.78%
51) Toluene-d8	8.41	98	798601	52.09	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.18%
72) 4-Bromofluorobenzene	11.03	95	280392	51.57	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	103.14%

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.07	76	66	N.D.		
15) Allyl chloride	3.26	76	56	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.56	63	84	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.39	96	76	N.D.		
27) 2-Butanone	5.56	43	129	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.82	83	726	N.D.		
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	6.09	117	62	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.41	78	82	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0411017.D 524.M Thu Apr 12 09:13:29 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411017.D  
 Acq On : 11 Apr 2007 13:47  
 Sample : JPL29-002  
 Misc : #2 5ml +Is/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:12 2007

Vial: 57  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.16	130	275		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.73	83	81		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	234		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.01	166	865		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.86	91	1071		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	1517		N.D.	
65) m,p-Xylene	10.09	106	516		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	d
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	229		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	221		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	67		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.75	119	57		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	12.11	146	481		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	12.20	146	701		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	816		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	72		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	14.66	180	206		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL29-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/28/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/11/2007 14:10

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.35	J
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.53	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.3	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-003  
 Lab File ID: M0411018.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:10  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.61	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-003  
 Lab File ID: M0411018.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:10  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.25	JB

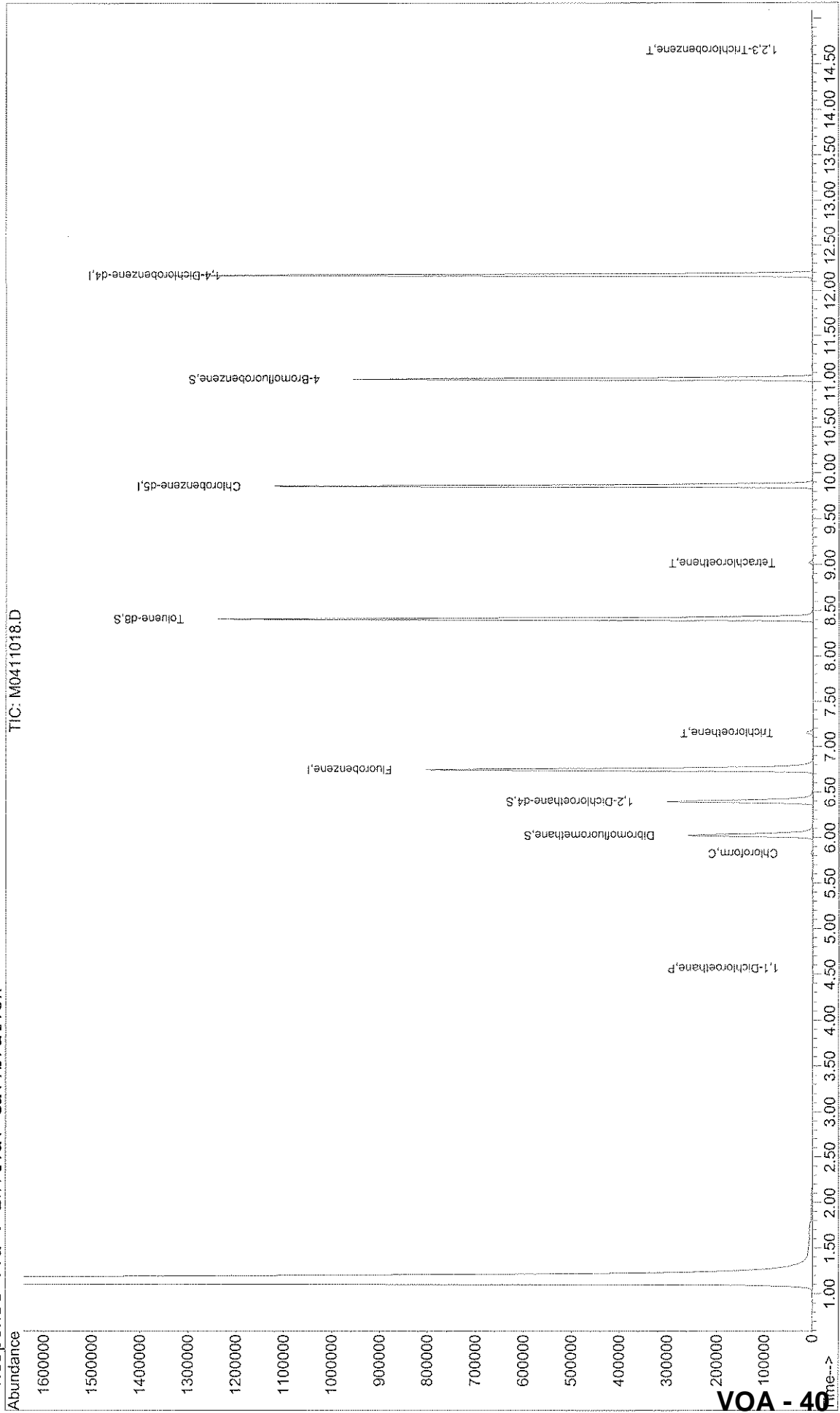
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411018.D  
Acq On : 11 Apr 2007 14:10  
Sample : JPL29-003  
Misc : #4 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:14 2007

Vial: 58  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 40

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411018.D  
 Acq On : 11 Apr 2007 14:10  
 Sample : JPL29-003  
 Misc : #4 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:14 2007

Vial: 58  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar)
1) Fluorobenzene	6.75	96	789241	50.00	ug/l	0.00	75.36%
50) Chlorobenzene-d5	9.86	82	335520	50.00	ug/l	0.00	74.06%
70) 1,4-Dichlorobenzene-d4	12.18	152	330694	50.00	ug/l	0.00	70.76%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	187700	49.98	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.96%	
37) 1,2-Dichloroethane-d4	6.40	65	232770	53.73	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	107.46%	
51) Toluene-d8	8.41	98	804023	51.55	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.10%	
72) 4-Bromofluorobenzene	11.03	95	277964	51.39	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	102.78%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.57	63	1978	0.35	ug/l	71
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.39	96	774	N.D.		
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.83	83	3317	0.53	ug/l	96
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.41	78	322	N.D.		

LU 4/12/07

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411018.D  
 Acq On : 11 Apr 2007 14:10  
 Sample : JPL29-003  
 Misc : #4 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:14 2007

Vial: 58  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

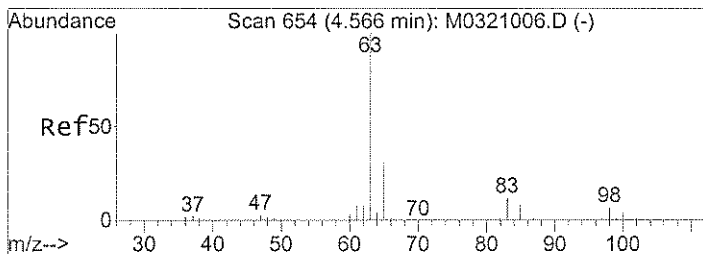
Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	4945	1.31	ug/l	95
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	7.40	63	58		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.72	83	319		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	151		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	2323	0.61	ug/l	96
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	125		N.D.	
62) 1-Chlorohexane	9.86	91	1246		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	1321		N.D.	
65) m,p-Xylene	10.10	106	667		N.D.	
66) o-xylene	10.50	106	57		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	350		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	84		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	57		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	224		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D. d	
83) 1,3-Dichlorobenzene	12.11	146	1347		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	12.11	146	1347		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	580		N.D.	
87) n-Butylbenzene	12.52	91	67		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	372		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D. d	
92) 1,2,3-Trichlorobenzene	14.66	180	933	0.25	ug/l #	93

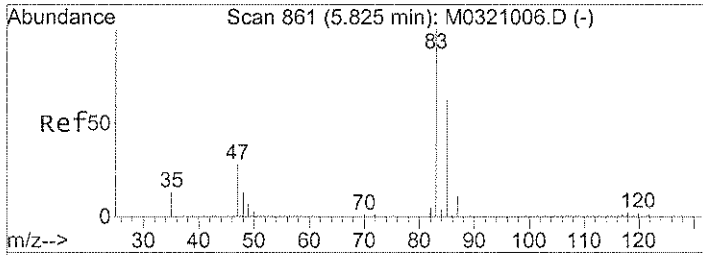
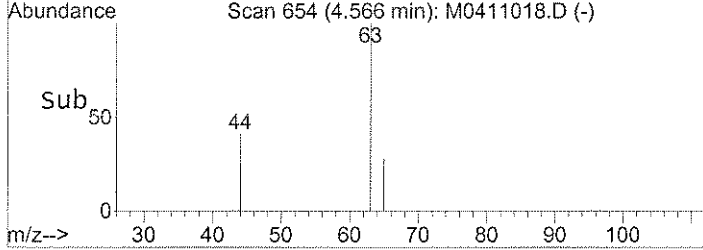
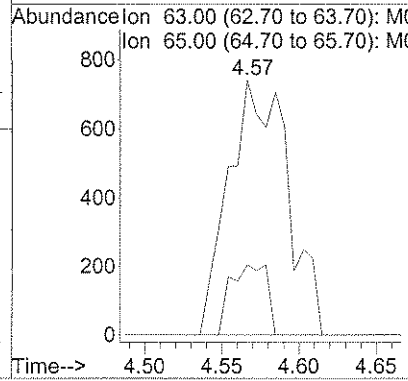
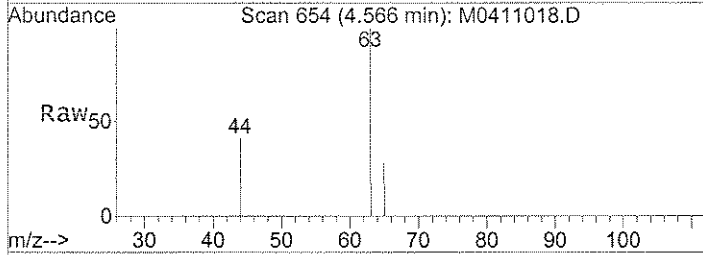
W 4/12/07





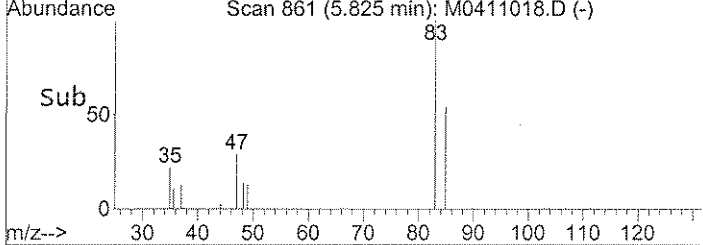
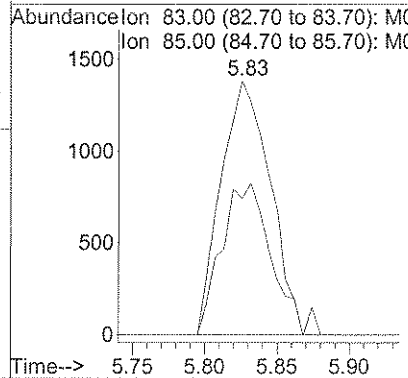
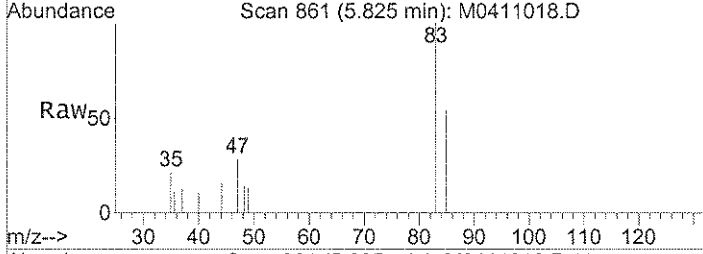
#22  
 1,1-Dichloroethane  
 Concen: 0.35 ug/l  
 RT: 4.57 min Scan# 654  
 Delta R.T. 0.01 min  
 Lab File: M0411018.D  
 Acq: 11 Apr 2007 14:10

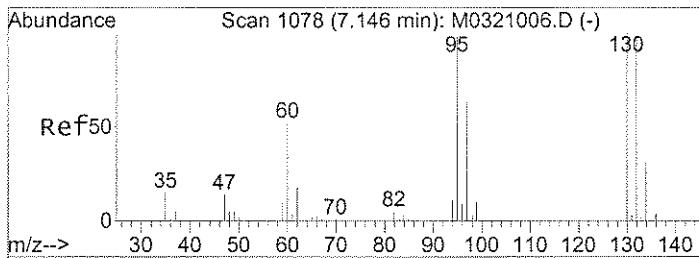
Tgt Ion: 63 Resp: 1978  
 Ion Ratio Lower Upper  
 63 100  
 65 17.0 13.7 53.7



#31  
 Chloroform  
 Concen: 0.53 ug/l  
 RT: 5.83 min Scan# 861  
 Delta R.T. -0.00 min  
 Lab File: M0411018.D  
 Acq: 11 Apr 2007 14:10

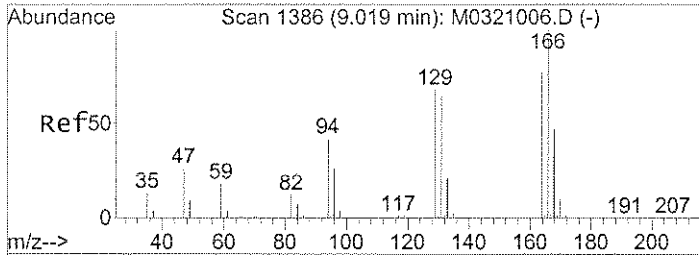
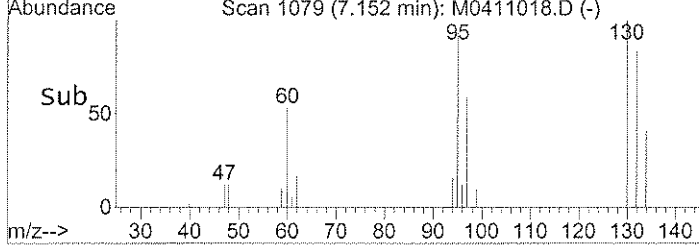
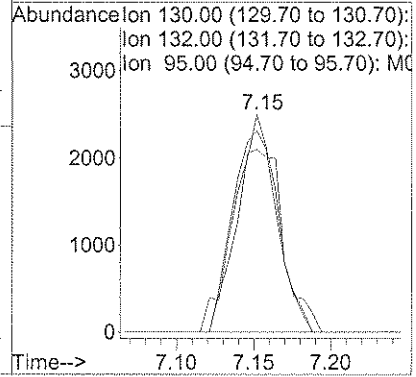
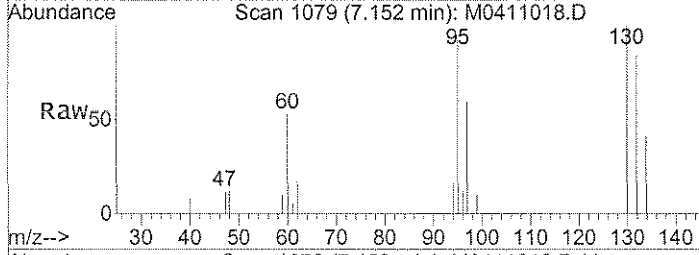
Tgt Ion: 83 Resp: 3317  
 Ion Ratio Lower Upper  
 83 100  
 85 58.2 41.2 81.2





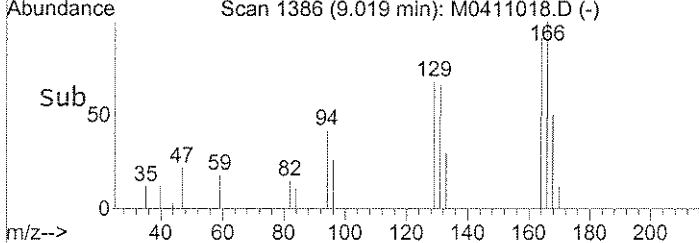
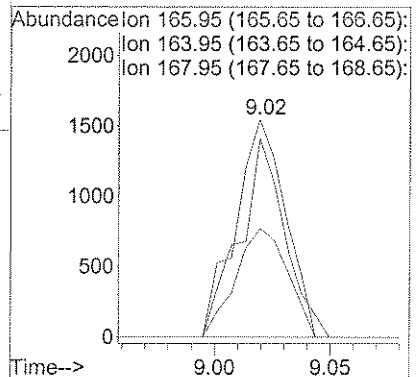
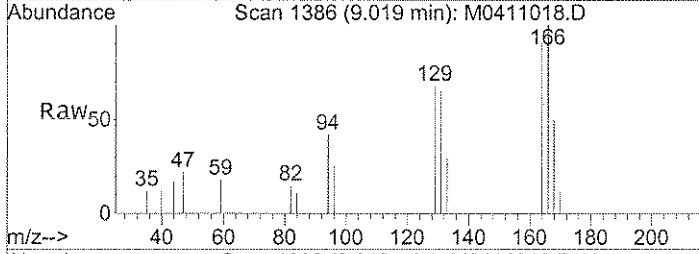
#41  
 Trichloroethene  
 Concen: 1.31 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0411018.D  
 Acq: 11 Apr 2007 14:10

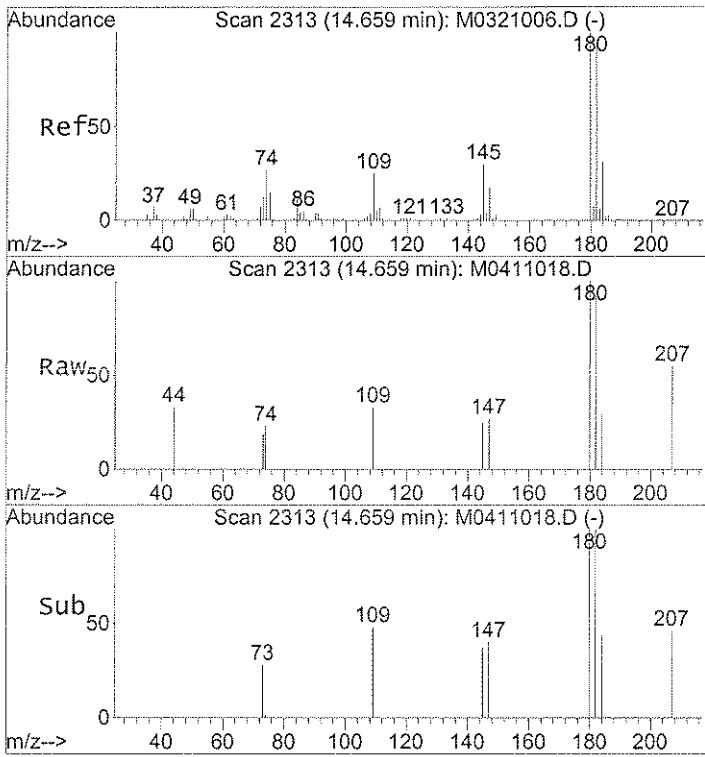
Tgt Ion	Resp	Lower	Upper
130	4945		
132	90.0	75.0	115.0
95	94.9	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 0.61 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. -0.00 min  
 Lab File: M0411018.D  
 Acq: 11 Apr 2007 14:10

Tgt Ion	Resp	Lower	Upper
166	2323		
164	83.1	63.3	94.9
168	51.7	39.6	59.4

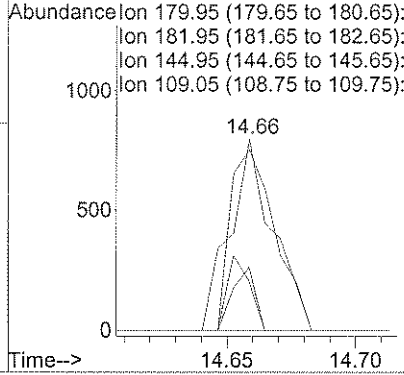




#92  
 1,2,3-Trichlorobenzene  
 Concen: 0.25 ug/l  
 RT: 14.66 min Scan# 2313  
 Delta R.T. -0.00 min  
 Lab File: M0411018.D  
 Acq: 11 Apr 2007 14:10

Tgt Ion: 180 Resp: 933

Ion	Ratio	Lower	Upper
180	100		
182	98.1	76.0	114.0
145	20.0	23.0	34.4#
109	17.1	19.0	28.4#



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-004  
 Lab File ID: M0411019.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.32	J
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.40	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	5.5	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-004  
 Lab File ID: M0411019.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.40	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-004  
 Lab File ID: M0411019.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

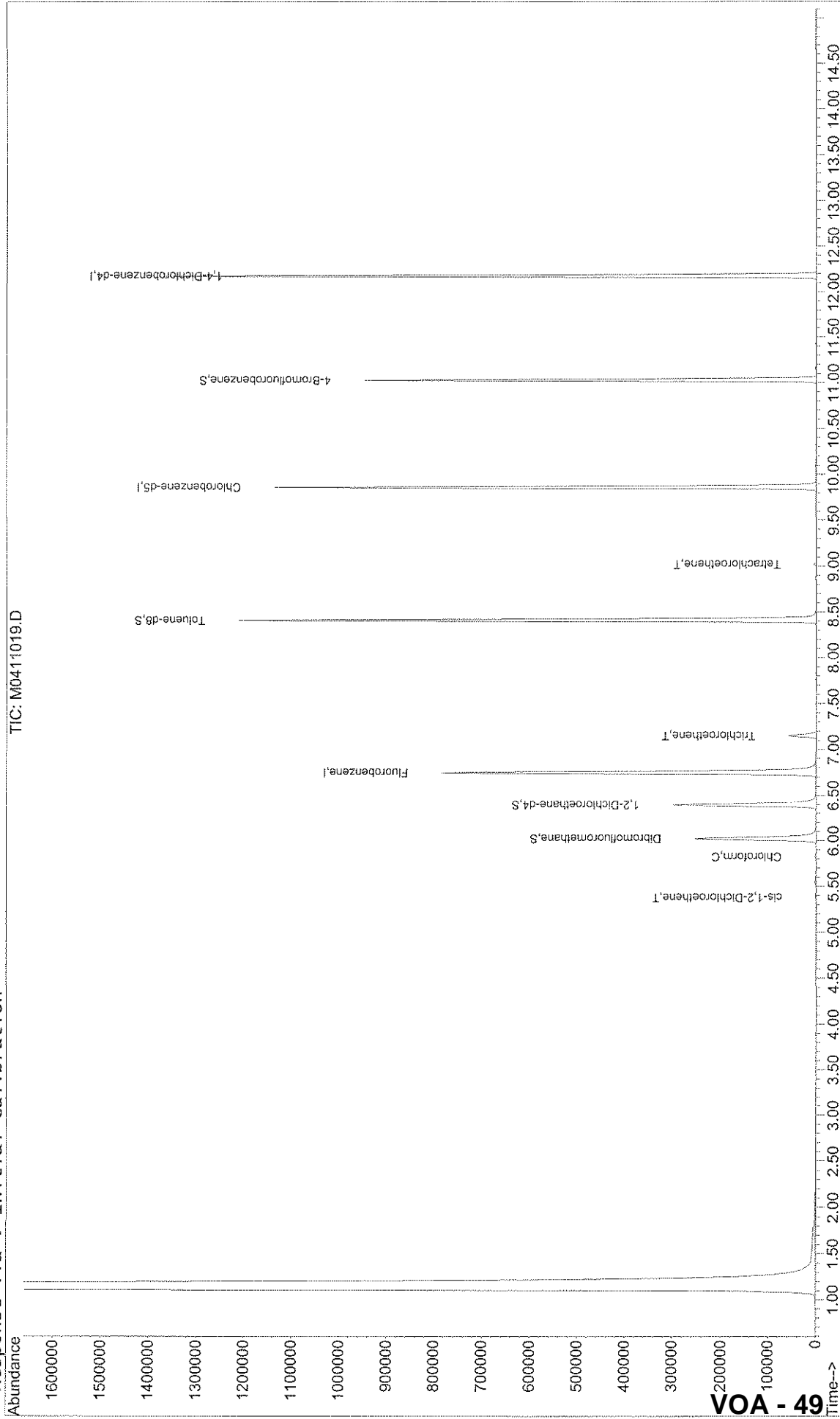
Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411019.D  
Acq On : 11 Apr 2007 14:33  
Sample : JPL29-004  
Misc : #3 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:16 2007

Vial: 59  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00

Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 49

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411019.D  
 Acq On : 11 Apr 2007 14:33  
 Sample : JPL29-004  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:16 2007

Vial: 59  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	775107	50.00	ug/l	0.00 74.01%
50) Chlorobenzene-d5	9.86	82	327264	50.00	ug/l	0.00 72.24%
70) 1,4-Dichlorobenzene-d4	12.18	152	327966	50.00	ug/l	0.00 70.18%

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) Dibromofluoromethane	6.02	111	184983	50.16	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	100.32%	
37) 1,2-Dichloroethane-d4	6.40	65	229876	54.03	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	108.06%	
51) Toluene-d8	8.41	98	792510	52.10	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	104.20%	
72) 4-Bromofluorobenzene	11.03	95	272045	50.71	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery	=	101.42%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	3.01	43	129	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	3.90	96	574	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	3.98	73	59	N.D.		
22) 1,1-Dichloroethane	4.58	63	1044	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.38	96	1079	0.32	ug/l	91
27) 2-Butanone	5.56	43	74	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.83	83	2474	0.40	ug/l	94
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.41	78	133	N.D.		

W 4/12/07



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411019.D  
 Acq On : 11 Apr 2007 14:33  
 Sample : JPL29-004  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:16 2007

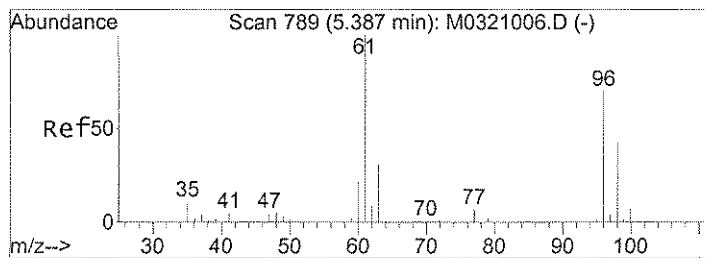
Vial: 59  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

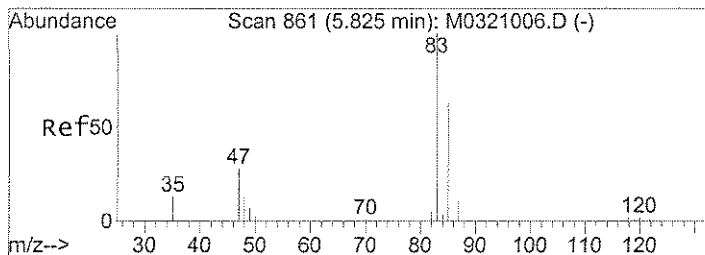
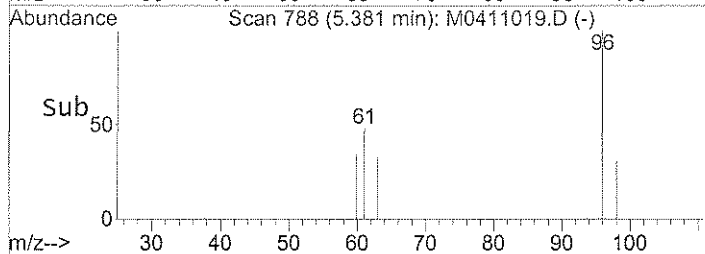
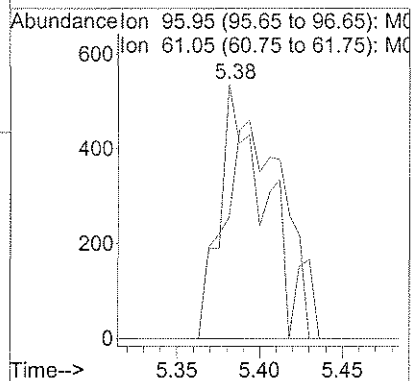
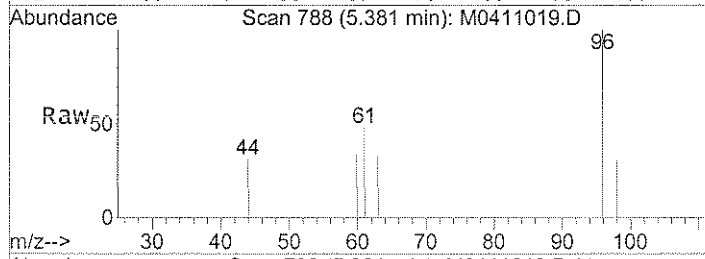
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	6.49	43	56		<del>Below Cal #</del>	22
41) Trichloroethene	7.15	130	20211	5.47	ug/l	95
42) Methylcyclohexane	0.00	83	0		N.D. d	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.72	83	173		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.47	92	85		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.01	166	1473	0.40	ug/l #	82
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.86	91	1299		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	337		N.D.	
65) m,p-Xylene	10.10	106	636		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D. d	
68) Bromoform	0.00	173	0		N.D. d	
69) Isopropylbenzene	11.03	105	193		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D. d	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	203		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.36	91	69		N.D.	
78) 4-Chlorotoluene	11.36	91	69		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	58		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D. d	
84) 4-Isopropyltoluene	0.00	119	0		N.D. d	
85) 1,4-Dichlorobenzene	12.19	146	131		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	66		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	197		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D. d	
92) 1,2,3-Trichlorobenzene	14.65	180	476		N.D.	

HW 4/12/07



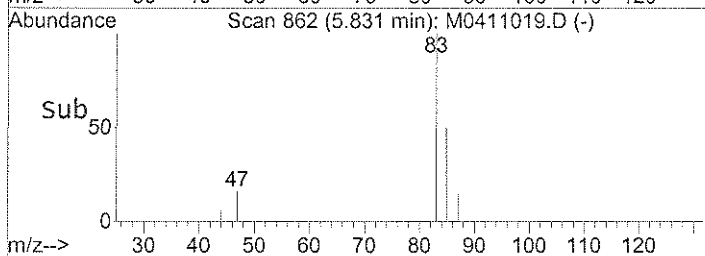
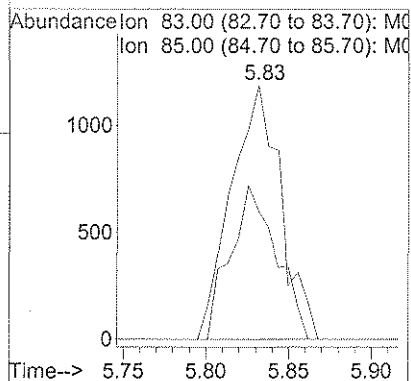
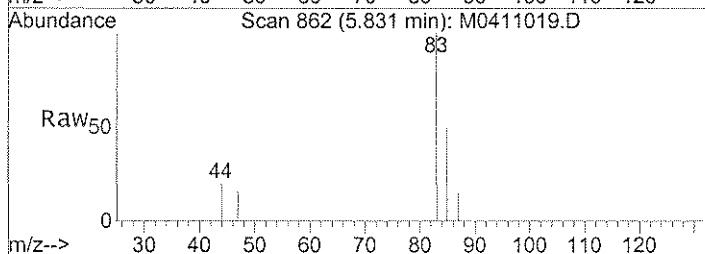
#26  
 cis-1,2-Dichloroethene  
 Concen: 0.32 ug/l  
 RT: 5.38 min Scan# 788  
 Delta R.T. -0.01 min  
 Lab File: M0411019.D  
 Acq: 11 Apr 2007 14:33

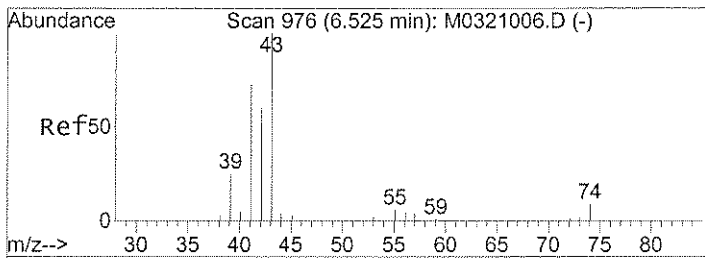
Tgt Ion: 96 Resp: 1079  
 Ion Ratio Lower Upper  
 96 100  
 61 106.9 93.4 140.2



#31  
 Chloroform  
 Concen: 0.40 ug/l  
 RT: 5.83 min Scan# 862  
 Delta R.T. 0.01 min  
 Lab File: M0411019.D  
 Acq: 11 Apr 2007 14:33

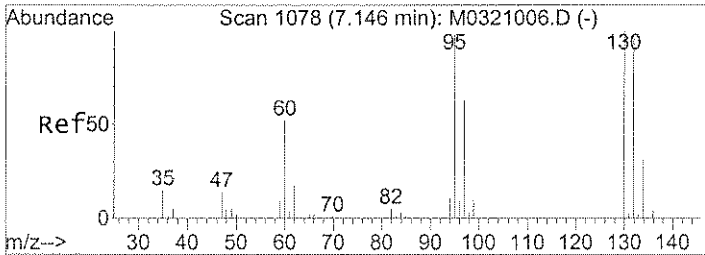
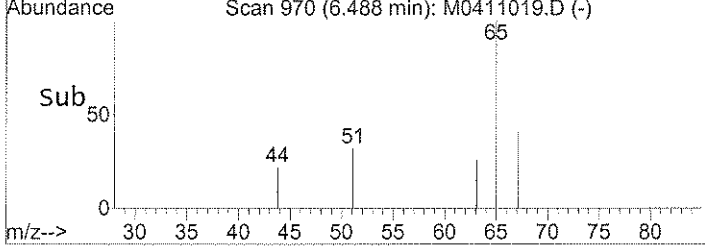
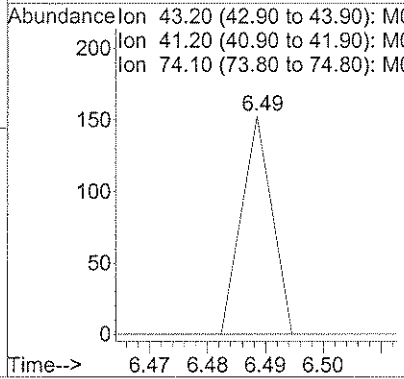
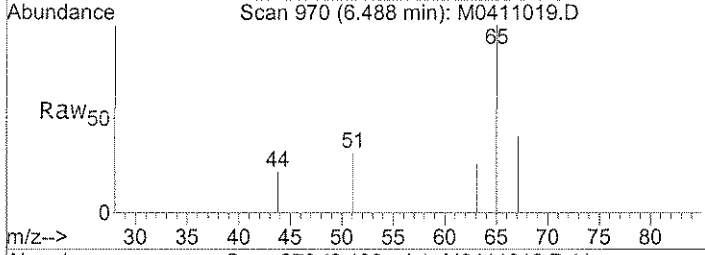
Tgt Ion: 83 Resp: 2474  
 Ion Ratio Lower Upper  
 83 100  
 85 56.6 41.2 81.2





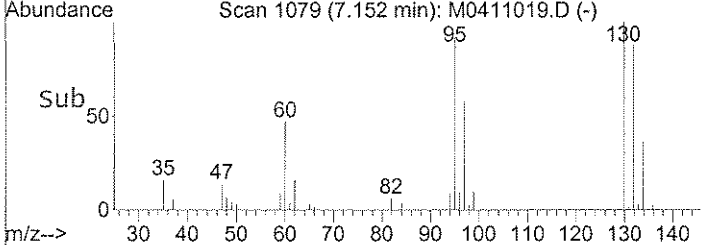
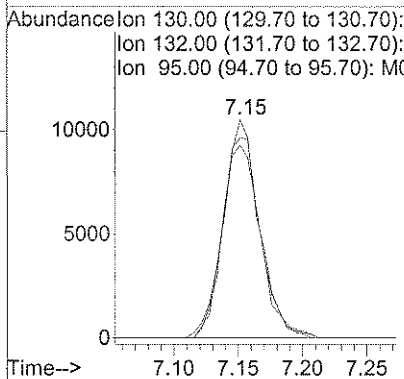
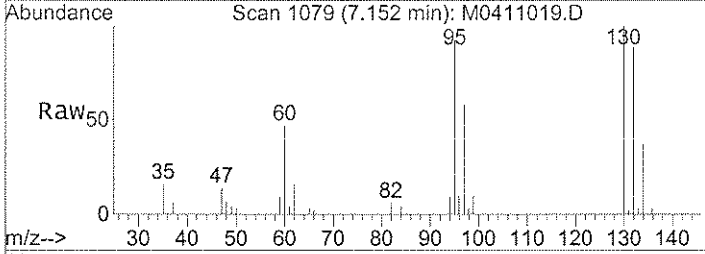
#40  
 Isobutanol  
 Concen: Below Cal  
 RT: 6.49 min Scan# 970  
 Delta R.T. -0.02 min  
 Lab File: M0411019.D  
 Acq: 11 Apr 2007 14:33

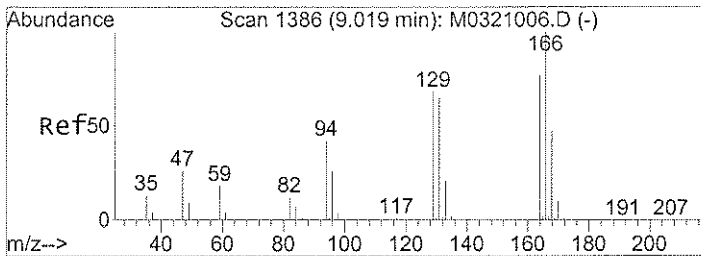
Tgt Ion	Resp	Lower	Upper
43	100		
41	0.0	55.8	83.8#
74	0.0	8.0	12.0#



#41  
 Trichloroethene  
 Concen: 5.47 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0411019.D  
 Acq: 11 Apr 2007 14:33

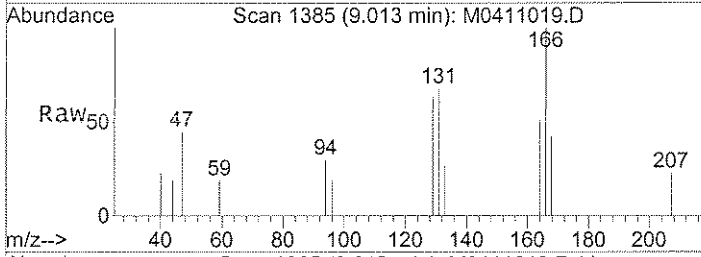
Tgt Ion	Resp	Lower	Upper
130	100		
132	95.3	75.0	115.0
95	97.8	69.4	109.4



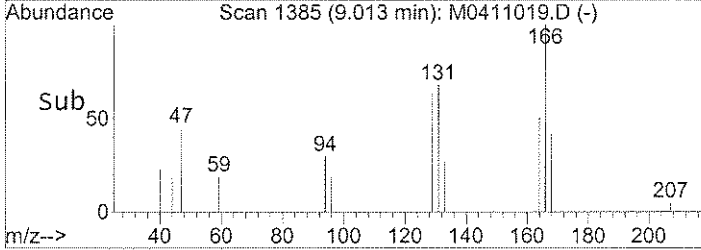
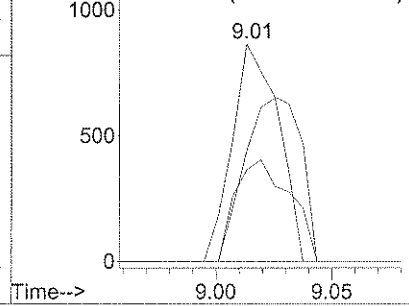


#56  
 Tetrachloroethene  
 Concen: 0.40 ug/l  
 RT: 9.01 min Scan# 1385  
 Delta R.T. -0.01 min  
 Lab File: M0411019.D  
 Acq: 11 Apr 2007 14:33

Tgt Ion	Resp	Lower	Upper
166	100		
164	56.3	63.3	94.9#
168	45.1	39.6	59.4



Abundance  
 Ion 165.95 (165.65 to 166.65):  
 Ion 163.95 (163.65 to 164.65):  
 Ion 167.95 (167.65 to 168.65):



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-005  
 Lab File ID: M0411020.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:56  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.43	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	2.5	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-005  
 Lab File ID: M0411020.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:56  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.31	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-14-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-005  
 Lab File ID: M0411020.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 14:56  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

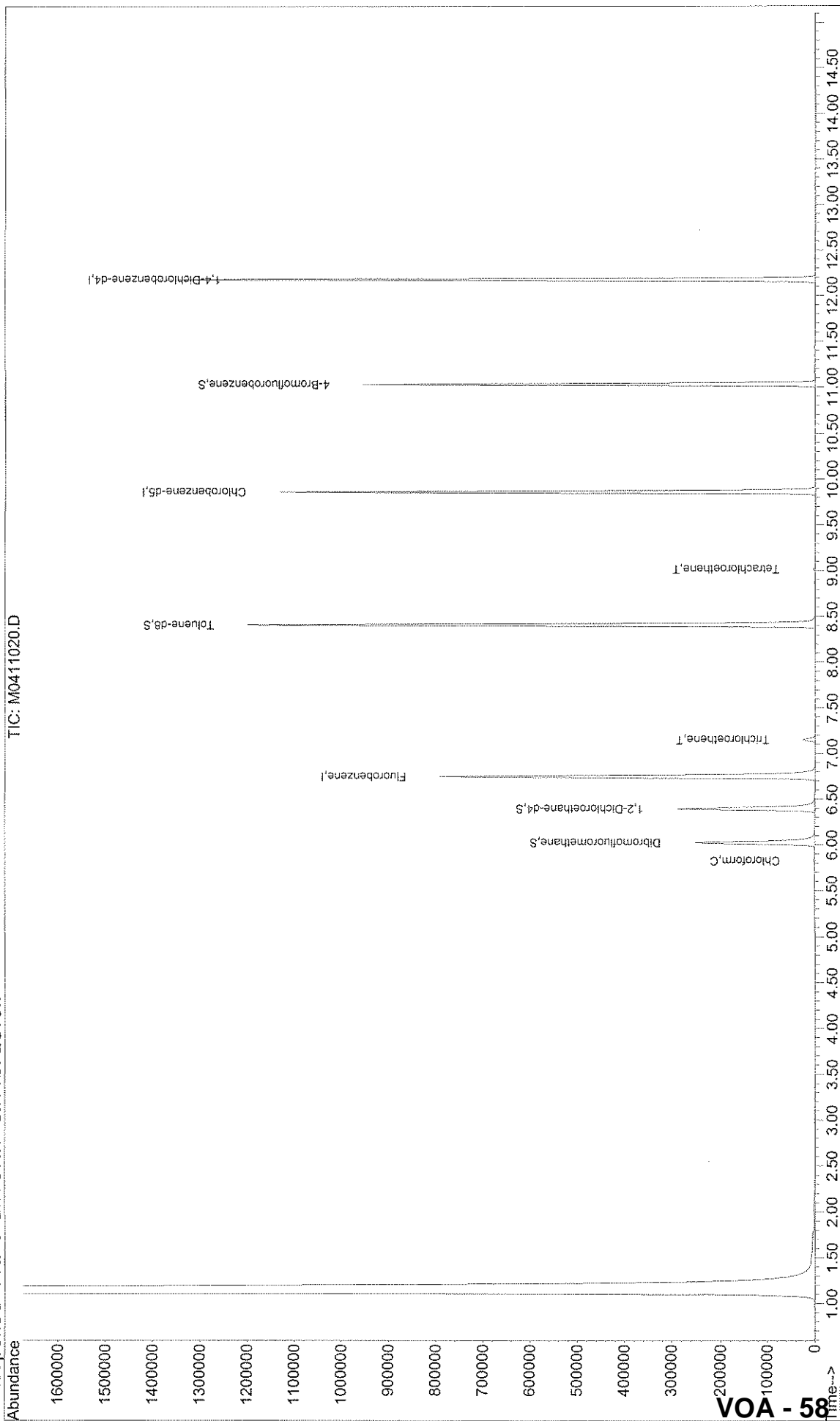
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411020.D  
Acq On : 11 Apr 2007 14:56  
Sample : JPL29-005  
Misc : #3 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:17 2007

Vial: 60  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 58



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411020.D  
 Acq On : 11 Apr 2007 14:56  
 Sample : JPL29-005  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:17 2007

Vial: 60  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.75	96	776957	50.00	ug/l	0.00 74.19%
50) Chlorobenzene-d5	9.86	82	329628	50.00	ug/l	0.00 72.76%
70) 1,4-Dichlorobenzene-d4	12.18	152	329543	50.00	ug/l	0.00 70.52%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	183240	49.57	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.14%
37) 1,2-Dichloroethane-d4	6.40	65	226103	53.02	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	106.04%
51) Toluene-d8	8.41	98	791916	51.68	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.36%
72) 4-Bromofluorobenzene	11.03	95	272124	50.48	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	100.96%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	d
11) Acetone	3.01	43	65		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	3.89	96	78		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	4.56	63	528		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.39	96	227		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.83	83	2657	0.43	ug/l	98
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	69		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0411020.D 524.M Thu Apr 12 09:17:28 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411020.D  
 Acq On : 11 Apr 2007 14:56  
 Sample : JPL29-005  
 Misc : #3 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:17 2007

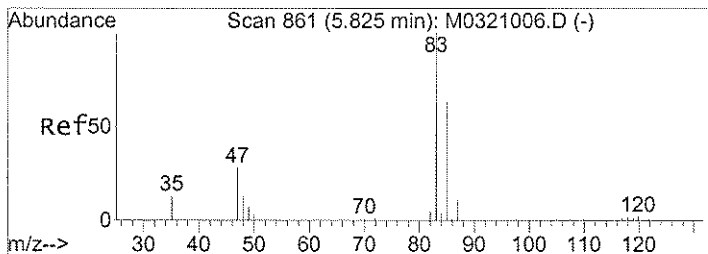
Vial: 60  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

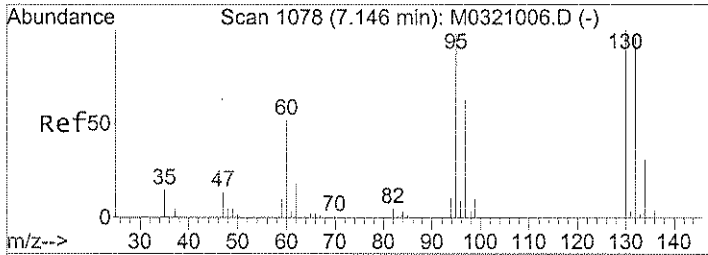
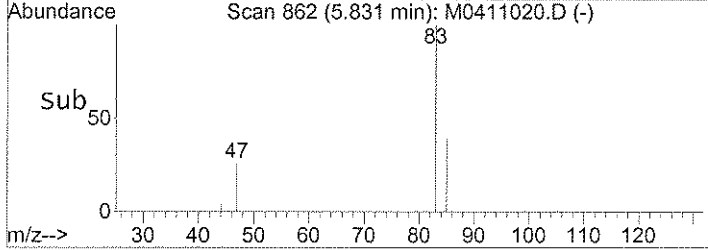
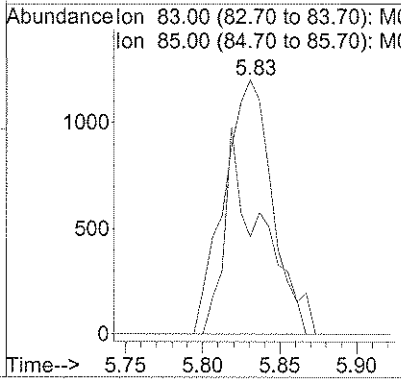
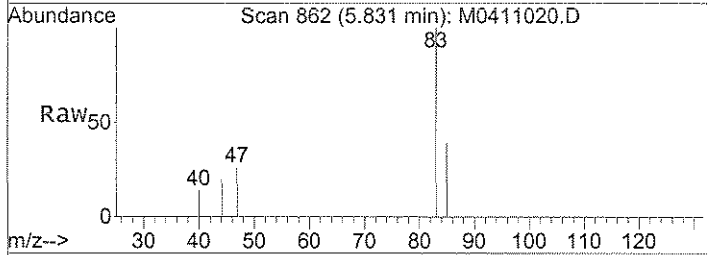
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.16	130	9416	2.54	ug/l	97
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.93	97	61		N.D.	
56) Tetrachloroethene	9.02	166	1159	0.31	ug/l	91
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.86	91	1034		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.00	91	130		N.D.	
65) m,p-Xylene	10.10	106	389		N.D.	
66) o-xylene	10.50	106	84		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	354		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	178		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	14.17	180	60		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	14.67	180	359		N.D.	

HA 4/12/07



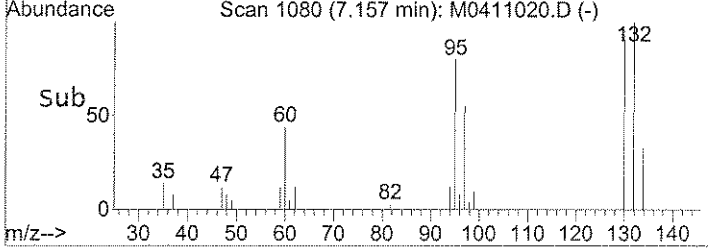
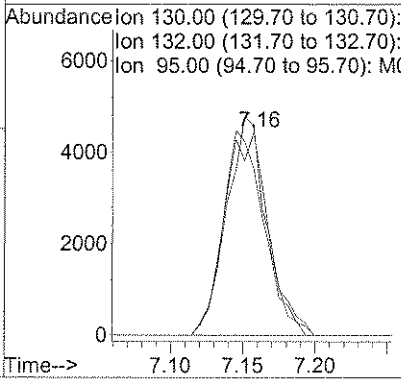
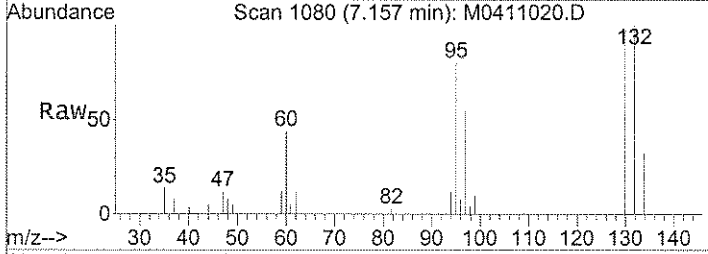
#31  
 Chloroform  
 Concen: 0.43 ug/l  
 RT: 5.83 min Scan# 862  
 Delta R.T. 0.00 min  
 Lab File: M0411020.D  
 Acq: 11 Apr 2007 14:56

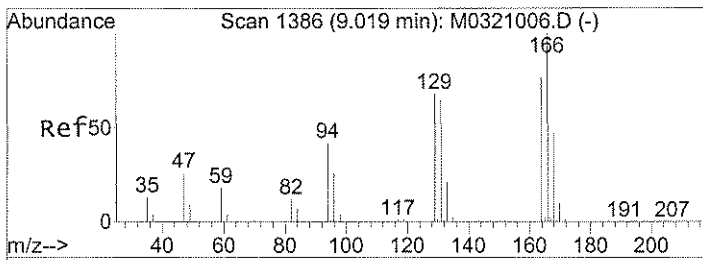
Tgt Ion	Resp	Lower	Upper
83	2657	100	
85	60.0	41.2	81.2



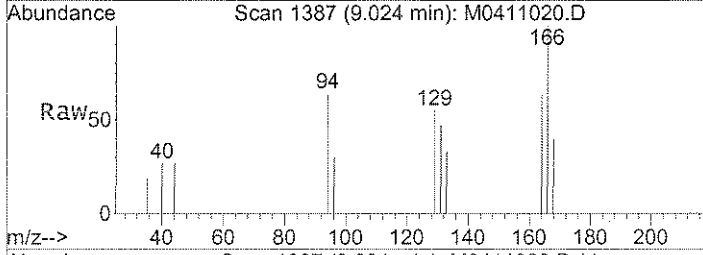
#41  
 Trichloroethene  
 Concen: 2.54 ug/l  
 RT: 7.16 min Scan# 1080  
 Delta R.T. 0.01 min  
 Lab File: M0411020.D  
 Acq: 11 Apr 2007 14:56

Tgt Ion	Resp	Lower	Upper
130	9416	100	
132	95.9	75.0	115.0
95	94.4	69.4	109.4



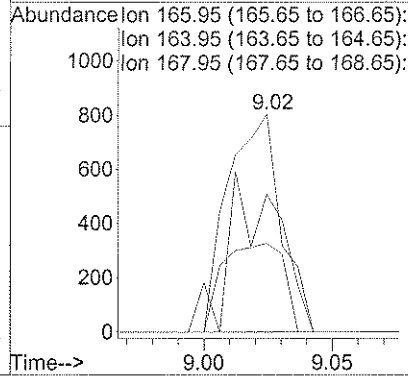
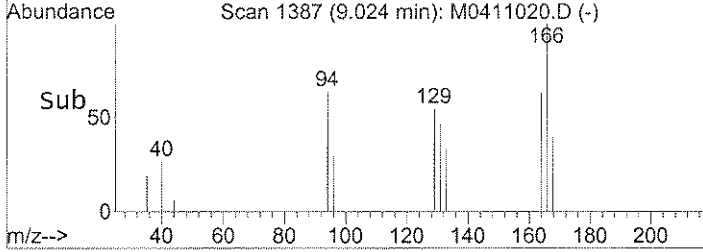


#56  
 Tetrachloroethene  
 Concen: 0.31 ug/l  
 RT: 9.02 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0411020.D  
 Acq: 11 Apr 2007 14:56



Tgt Ion: 166 Resp: 1159

Ion	Ratio	Lower	Upper
166	100		
164	68.5	63.3	94.9
168	46.2	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-006  
 Lab File ID: M0411021.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.25	JB
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-006  
 Lab File ID: M0411021.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.88	J
95-47-6	o-Xylene	0.32	J
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-006  
 Lab File ID: M0411021.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

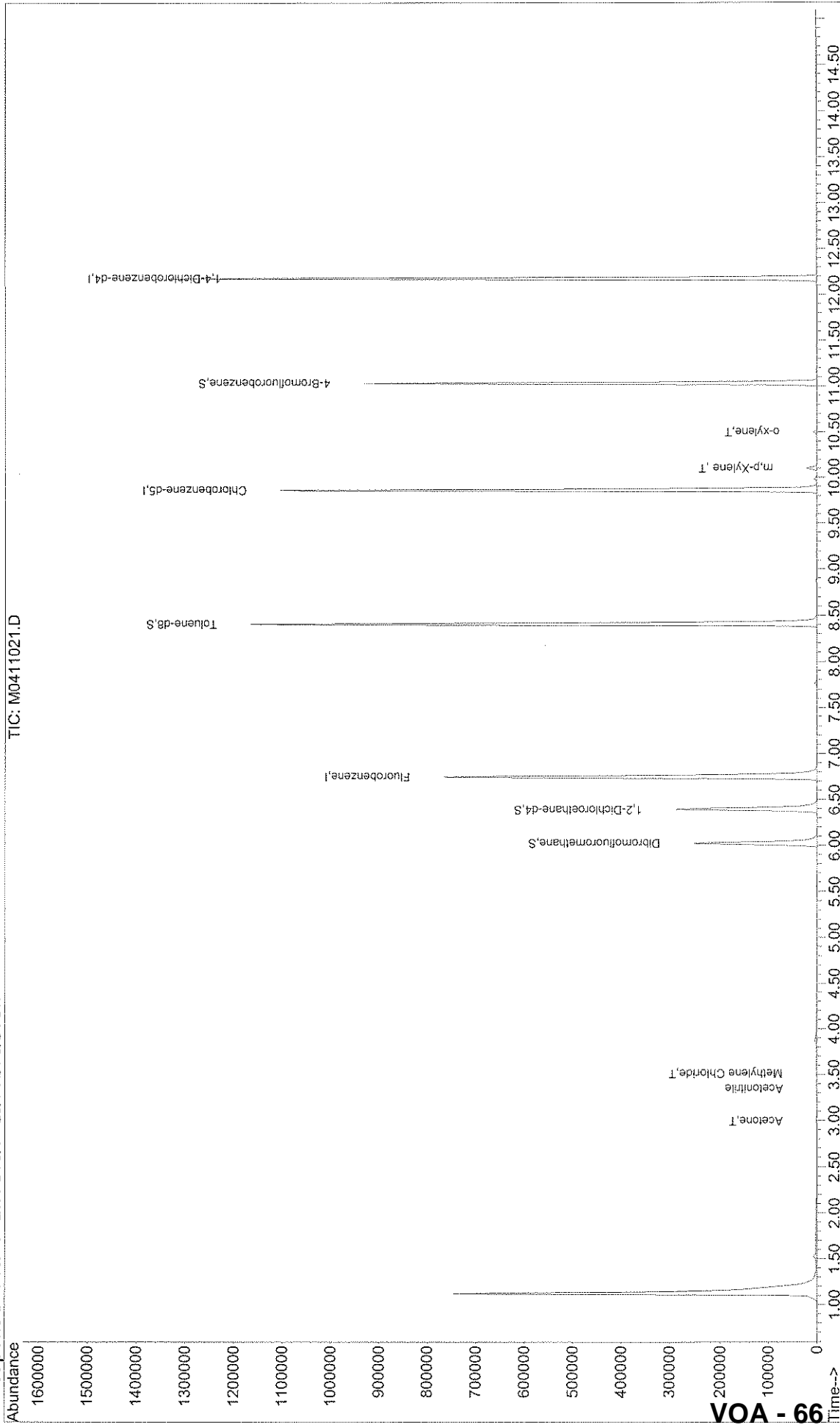
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411021.D  
Acq On : 11 Apr 2007 15:20  
Sample : JPL29-006  
Misc : #2 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:18 2007

Vial: 61  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 66



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411021.D  
 Acq On : 11 Apr 2007 15:20  
 Sample : JPL29-006  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:18 2007

Vial: 61  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)

Title : VOA 524- 5ML Calibration 5973M

Last Update : Tue Apr 10 10:47:40 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	765525	50.00	ug/l	0.00 73.10%
50) Chlorobenzene-d5	9.87	82	323532	50.00	ug/l	0.00 71.42%
70) 1,4-Dichlorobenzene-d4	12.18	152	327401	50.00	ug/l	0.00 70.06%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	181924	49.95	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	99.90%		
37) 1,2-Dichloroethane-d4	6.40	65	222201	52.88	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	105.76%		
51) Toluene-d8	8.41	98	780108	51.87	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	103.74%		
72) 4-Bromofluorobenzene	11.03	95	272824	50.95	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery =	101.90%		

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D. d	
11) Acetone	3.00	43	1586	1.95 ug/l	97
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	0.00	76	0	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	3.35	40	112	0.68 ug/l #	1
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	3.51	84	72	0.25 ug/l #	61
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	5.55	43	206	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	5.84	83	73	N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D. d	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	6.42	78	217	N.D.	

HJ 4/12/07

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411021.D  
 Acq On : 11 Apr 2007 15:20  
 Sample : JPL29-006  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:18 2007

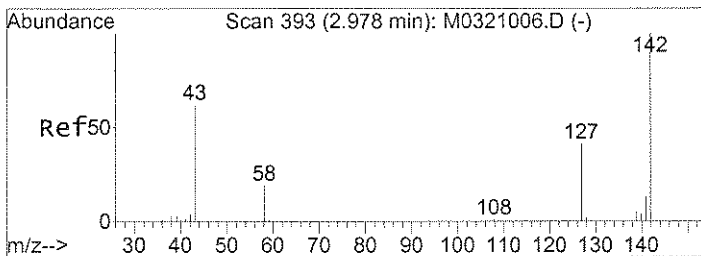
Vial: 61  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

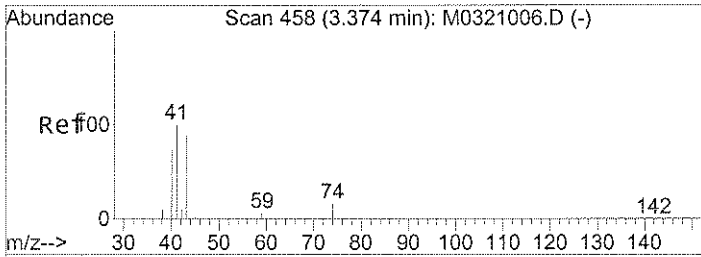
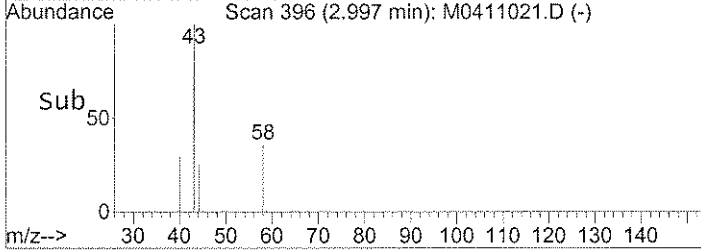
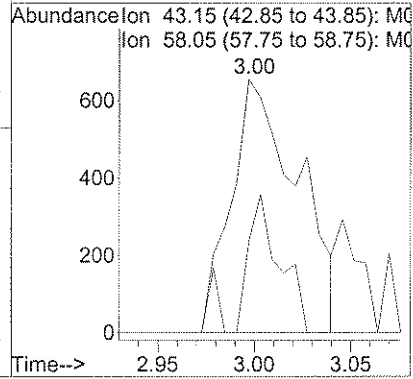
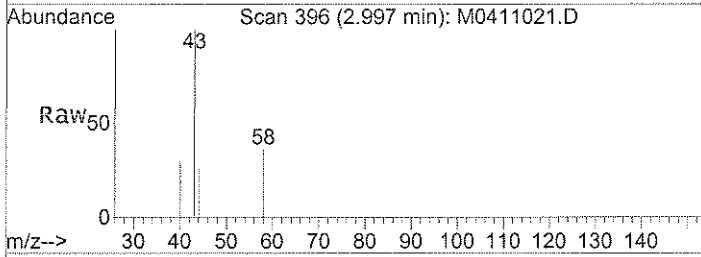
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	206		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	9.00	43	74		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.87	91	1300		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	4389		N.D.	
65) m,p-Xylene	10.10	106	6214	0.88	ug/l	99
66) o-xylene	10.49	106	2035	0.32	ug/l	89
67) Styrene	0.00	104	0		N.D.	d
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	10.85	105	61		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.37	156	56		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	58		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	77		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	d
82) sec-butylbenzene	0.00	105	0		N.D.	d
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	12.54	91	56		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	14.66	180	80		N.D.	

W 4/12/07



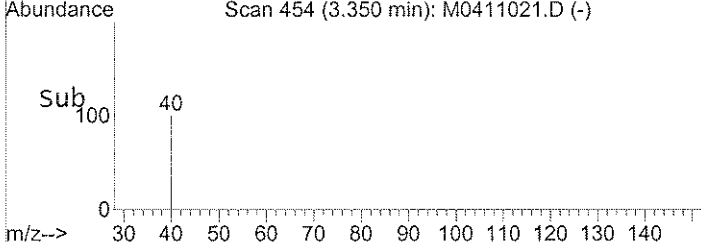
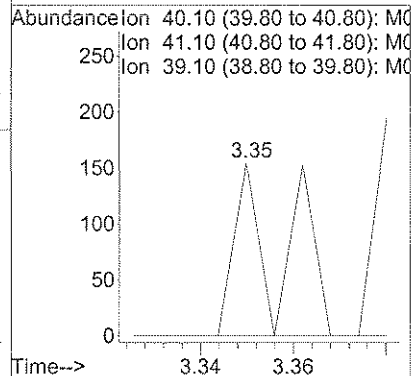
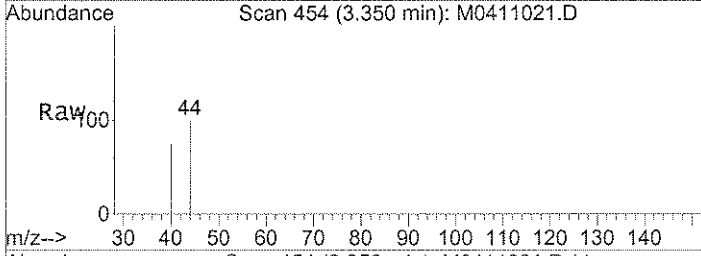
#11  
 Acetone  
 Concen: 1.95 ug/l  
 RT: 3.00 min Scan# 396  
 Delta R.T. 0.04 min  
 Lab File: M0411021.D  
 Acq: 11 Apr 2007 15:20

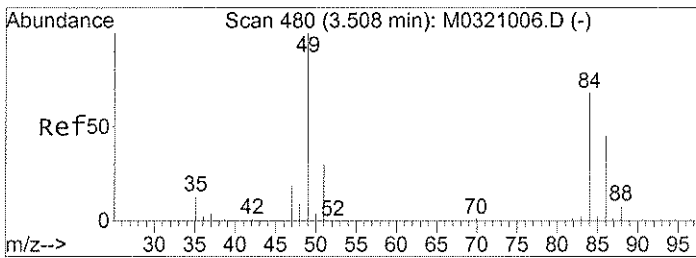
Tgt Ion	43	58	Resp	1586	Lower	Upper
Ion Ratio	100	25.7	21.9	32.9		



#16  
 Acetonitrile  
 Concen: 0.68 ug/l  
 RT: 3.35 min Scan# 454  
 Delta R.T. -0.00 min  
 Lab File: M0411021.D  
 Acq: 11 Apr 2007 15:20

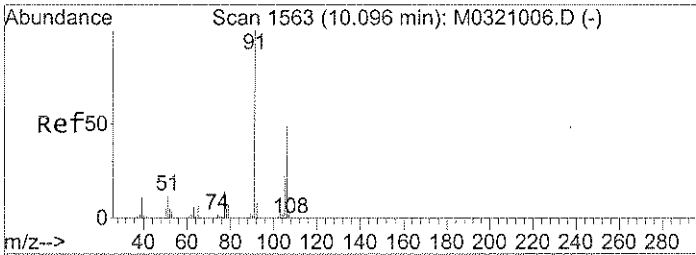
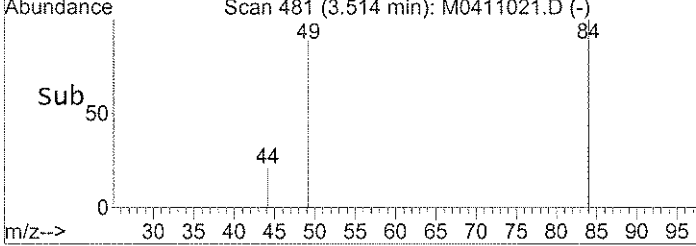
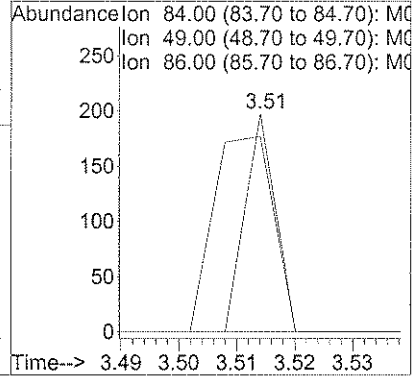
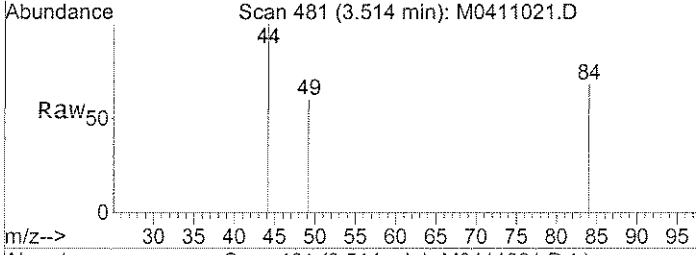
Tgt Ion	40	Resp	112	Lower	Upper
Ion Ratio	100	0.0	204.1	306.1#	
	41	0.0	76.9	115.3#	





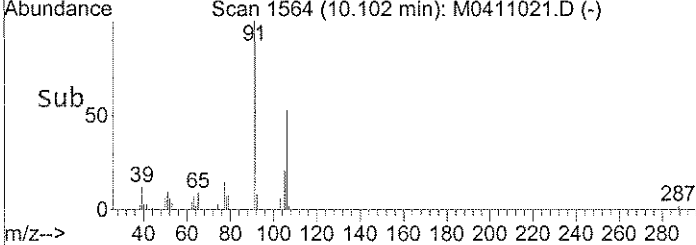
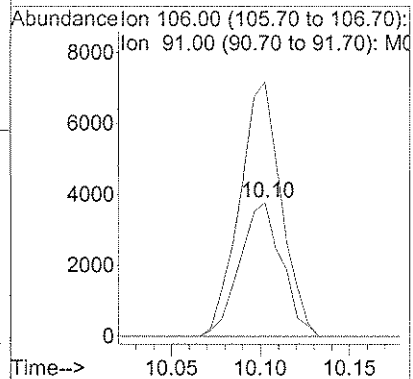
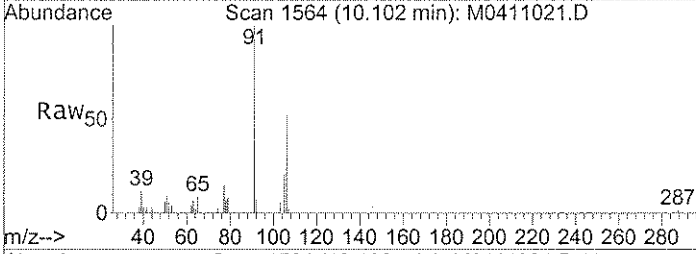
#18  
 Methylene Chloride  
 Concen: 0.25 ug/l  
 RT: 3.51 min Scan# 481  
 Delta R.T. 0.01 min  
 Lab File: M0411021.D  
 Acq: 11 Apr 2007 15:20

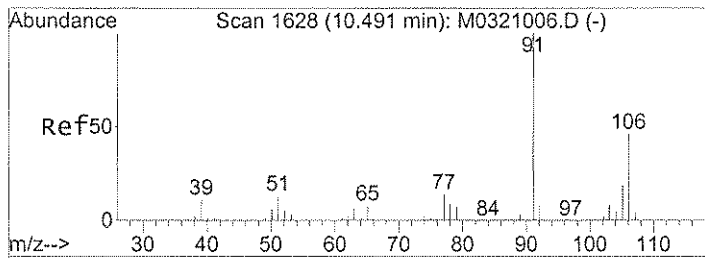
Tgt Ion	Resp	Lower	Upper
84	100		
49	176.4	133.5	173.5#
86	0.0	49.1	89.1#



#65  
 m,p-Xylene  
 Concen: 0.88 ug/l  
 RT: 10.10 min Scan# 1564  
 Delta R.T. -0.00 min  
 Lab File: M0411021.D  
 Acq: 11 Apr 2007 15:20

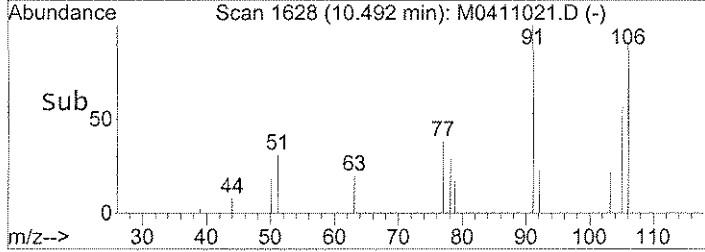
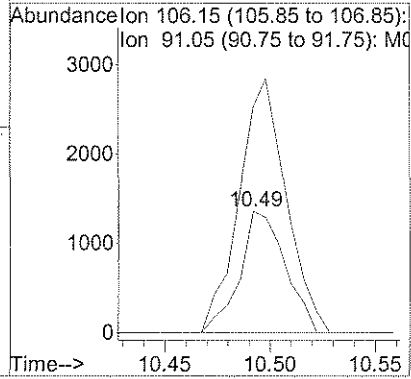
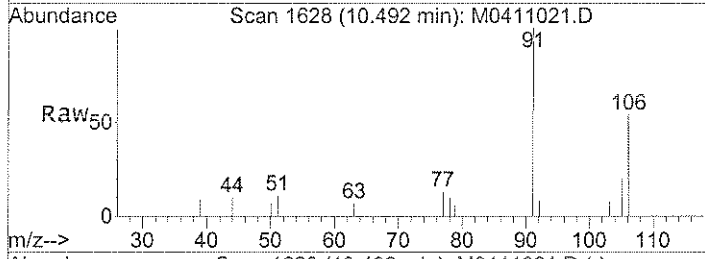
Tgt Ion	Resp	Lower	Upper
106	100		
91	187.9	169.6	209.6





#66  
 o-xylene  
 Concen: 0.32 ug/l  
 RT: 10.49 min Scan# 1628  
 Delta R.T. -0.01 min  
 Lab File: M0411021.D  
 Acq: 11 Apr 2007 15:20

Tgt Ion: 106 Resp: 2035  
 Ion Ratio Lower Upper  
 106 100  
 91 217.1 160.6 240.8



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-007  
 Lab File ID: M0411022.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.44	JB
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-007  
 Lab File ID: M0411022.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-007  
 Lab File ID: M0411022.D  
 Date Collected: 03/28/2007  
 Date/Time Analyzed: 04/11/2007 15:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

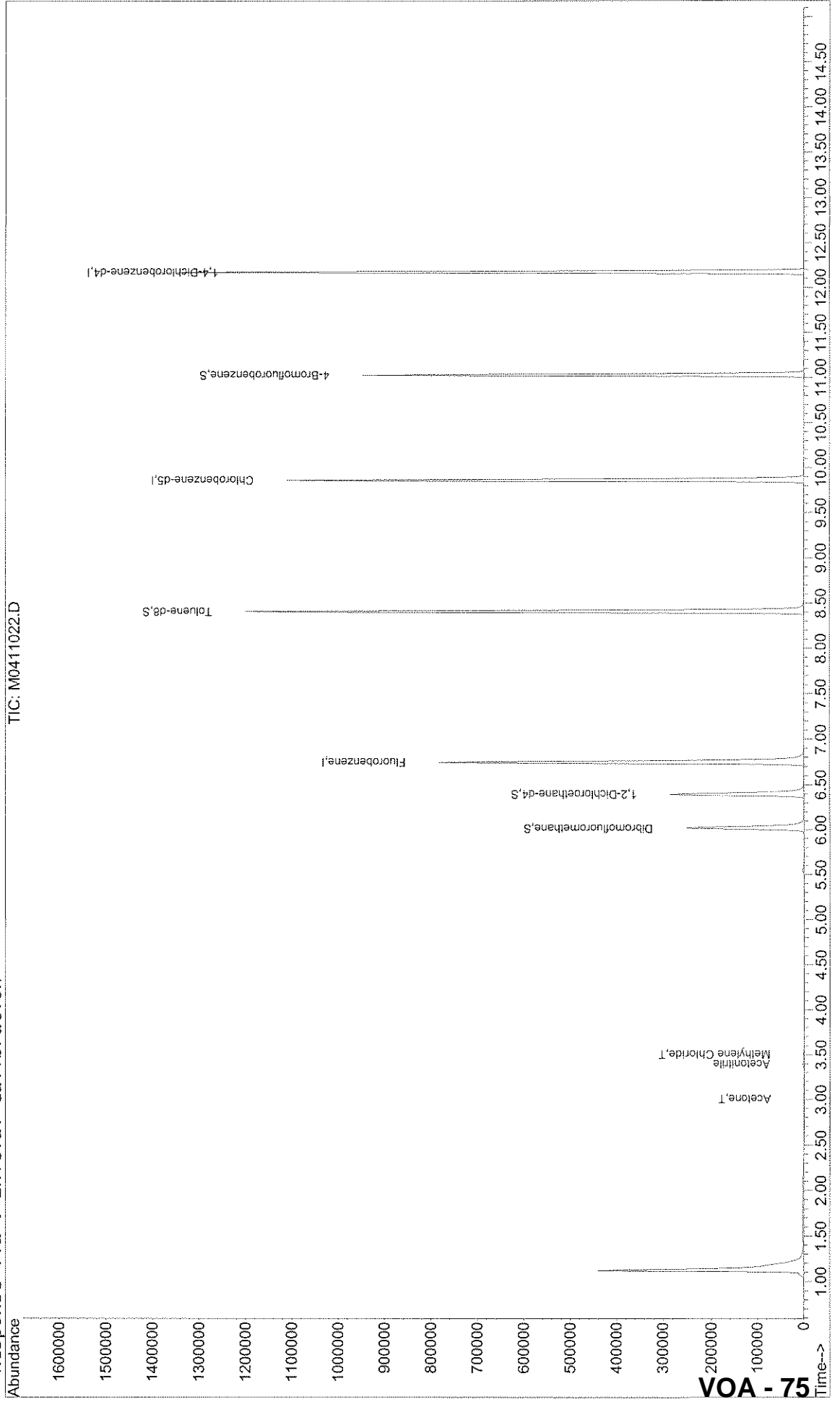
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411022.D  
Acq On : 11 Apr 2007 15:43 Vial: 62  
Sample : JPL29-007 Operator: DGA  
Misc : #2 5ml +IS/SS (524) Inst : MOBY  
MS Integration Params: rreint.p Multiplr: 1.00  
Quant Time: Apr 12 9:20 2007 Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411022.D  
 Acq On : 11 Apr 2007 15:43  
 Sample : JPL29-007  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:20 2007

Vial: 62  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar)
1) Fluorobenzene	6.75	96	769424	50.00	ug/l	0.00 73.47%
50) Chlorobenzene-d5	9.86	82	325111	50.00	ug/l	0.00 71.77%
70) 1,4-Dichlorobenzene-d4	12.18	152	327163	50.00	ug/l	0.00 70.01%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	182404	49.82	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.64%
37) 1,2-Dichloroethane-d4	6.40	65	224069	53.06	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	106.12%
51) Toluene-d8	8.41	98	783585	51.85	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	103.70%
72) 4-Bromofluorobenzene	11.03	95	270591	50.57	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	101.14%

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	3.01	43	1280	1.57	ug/l #	57
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.07	76	413	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	3.40	40	3093	18.65	ug/l #	24
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	3.51	84	676	0.44	ug/l	96
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.54	43	70	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	0.00	83	0	N.D.		
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.41	78	161	N.D.		

WS 4/12/07

(#) = qualifier out of range (m) = manual integration  
 M0411022.D 524.M Thu Apr 12 09:20:06 2007

Quantitation Report

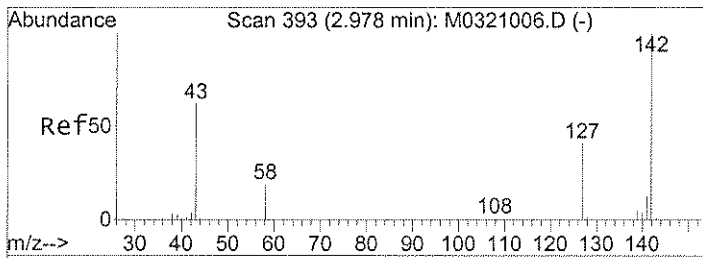
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 Acq On : 11 Apr 2007 15:43  
 Sample : JPL29-007  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:20 2007

Vial: 62  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

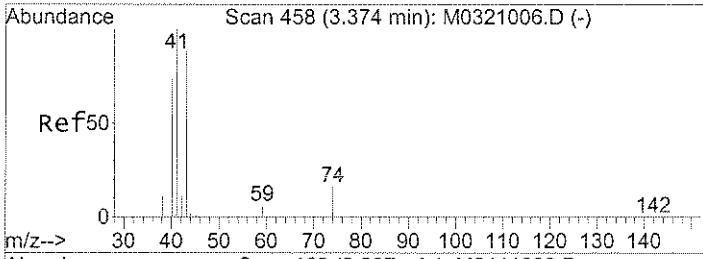
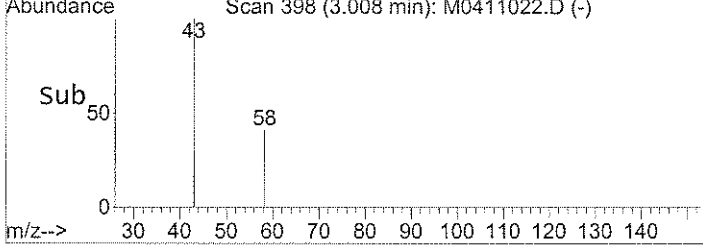
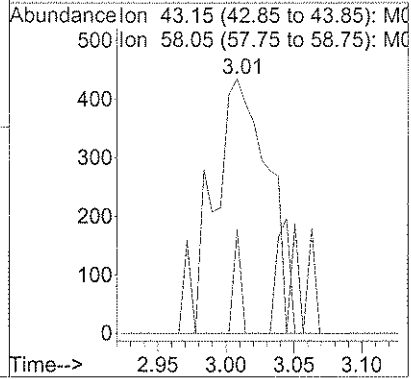
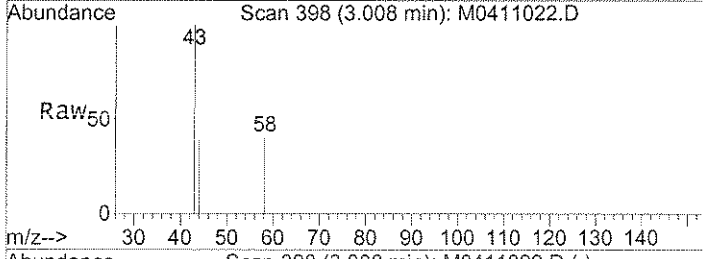
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	78		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	56		N.D.	
62) 1-Chlorohexane	9.86	91	1219		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	59		N.D.	
65) m,p-Xylene	10.10	106	64		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	335		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	55		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	143		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	139		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	d
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	12.53	91	122		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



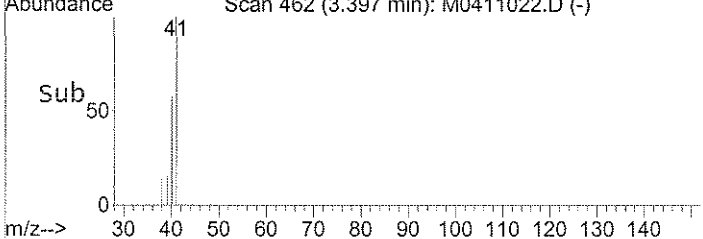
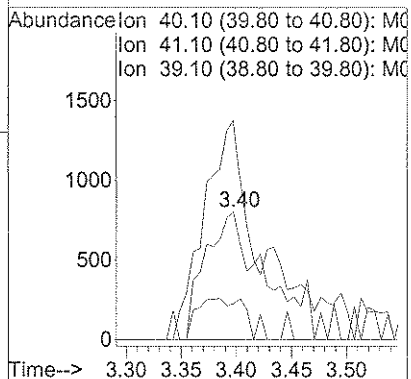
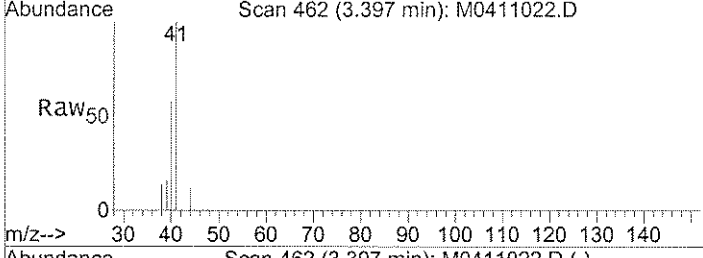
#11  
 Acetone  
 Concen: 1.57 ug/l  
 RT: 3.01 min Scan# 398  
 Delta R.T. 0.05 min  
 Lab File: M0411022.D  
 Acq: 11 Apr 2007 15:43

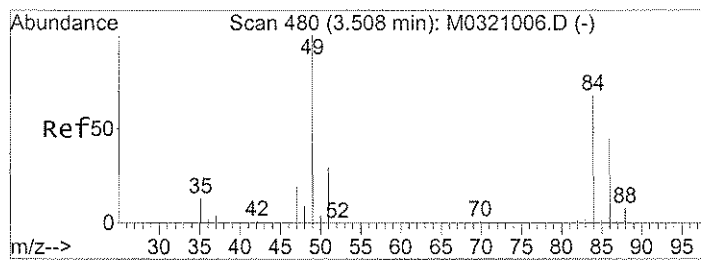
Tgt Ion: 43 Resp: 1280  
 Ion Ratio Lower Upper  
 43 100  
 58 5.1 21.9 32.9#



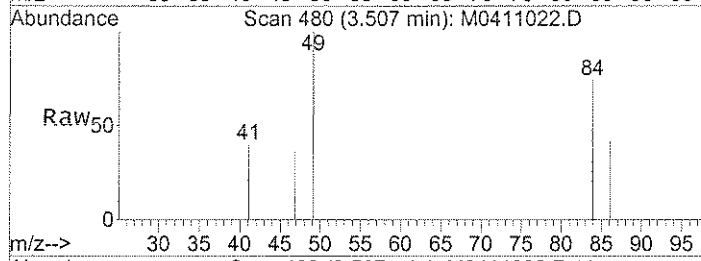
#16  
 Acetonitrile  
 Concen: 18.65 ug/l  
 RT: 3.40 min Scan# 462  
 Delta R.T. 0.05 min  
 Lab File: M0411022.D  
 Acq: 11 Apr 2007 15:43

Tgt Ion: 40 Resp: 3093  
 Ion Ratio Lower Upper  
 40 100  
 41 118.0 204.1 306.1#  
 39 25.9 76.9 115.3#



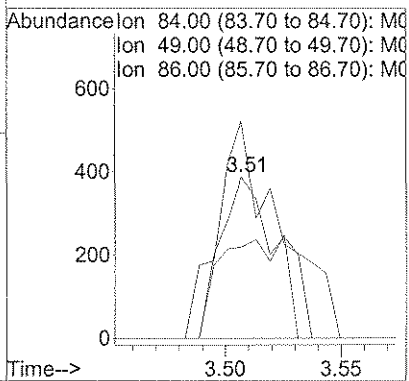
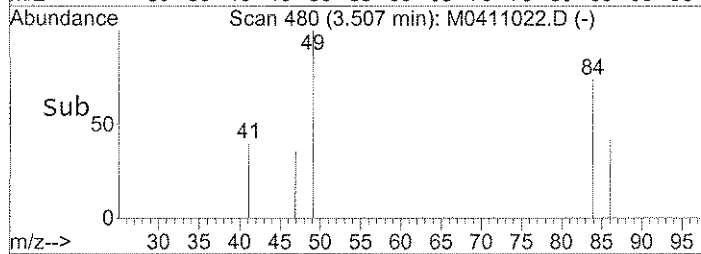


#18  
 Methylene Chloride  
 Concen: 0.44 ug/l  
 RT: 3.51 min Scan# 480  
 Delta R.T. 0.00 min  
 Lab File: M0411022.D  
 Acq: 11 Apr 2007 15:43



Tgt Ion: 84 Resp: 676

Ion	Ratio	Lower	Upper
84	100		
49	146.7	133.5	173.5
86	68.9	49.1	89.1



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-008  
 Lab File ID: M0411023.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:05  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.74	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-008  
 Lab File ID: M0411023.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:05  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-008  
 Lab File ID: M0411023.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:05  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

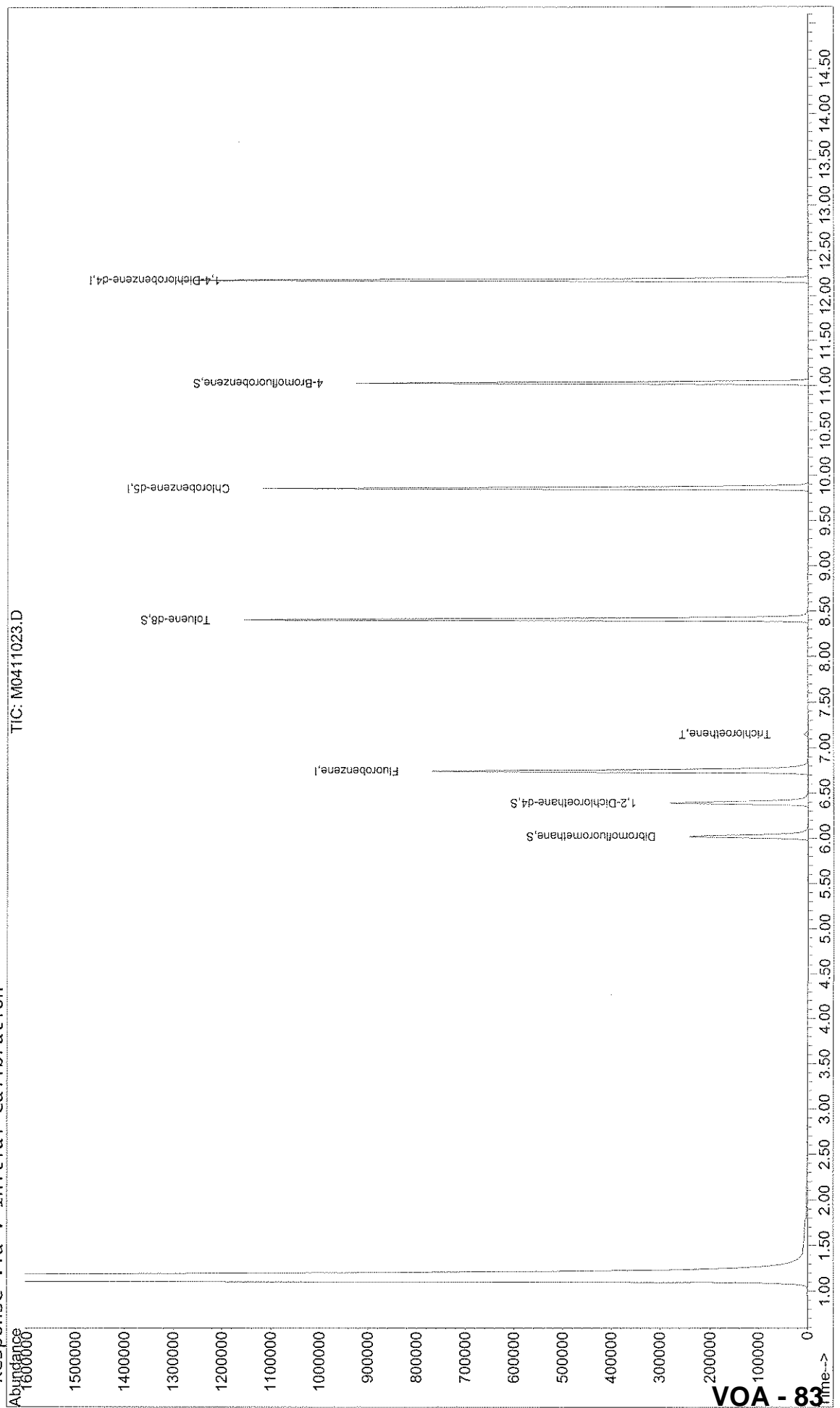


Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411023.D  
Acq On : 11 Apr 2007 16:05  
Sample : JPL29-008  
Misc : #2 5ml +IS/SS (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:27 2007

Vial: 63  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response Via : Initial Calibration



VOA - 83

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411023.D  
 Acq On : 11 Apr 2007 16:05  
 Sample : JPL29-008  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:27 2007

Vial: 63  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.75	96	747512	50.00	ug/l	0.00 71.38%
50) Chlorobenzene-d5	9.86	82	316882	50.00	ug/l	0.00 69.95%
70) 1,4-Dichlorobenzene-d4	12.18	152	317907	50.00	ug/l	0.00 68.03%

System Monitoring Compounds

33) Dibromofluoromethane	6.03	111	177758	49.98	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.96%
37) 1,2-Dichloroethane-d4	6.40	65	217911	53.11	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	106.22%
51) Toluene-d8	8.41	98	766236	52.02	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.04%
72) 4-Bromofluorobenzene	11.03	95	264905	50.94	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	101.88%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	1.44	50	63		N.D.	
4) Vinyl chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	d
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	3.95	53	75		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.84	83	935		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.42	78	332		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0411023.D 524.M Thu Apr 12 09:27:22 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411023.D  
 Acq On : 11 Apr 2007 16:05  
 Sample : JPL29-008  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:27 2007

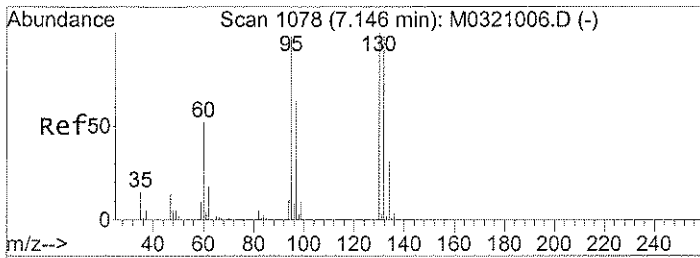
Vial: 63  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

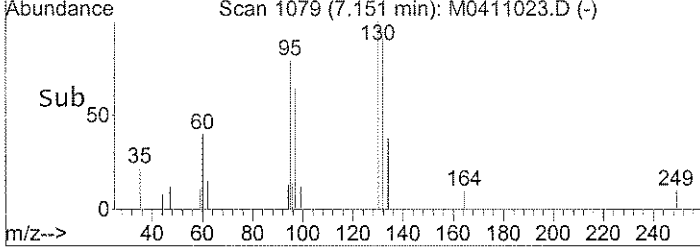
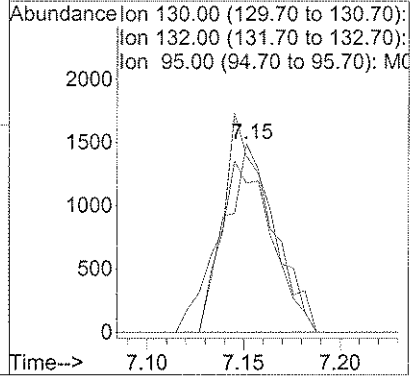
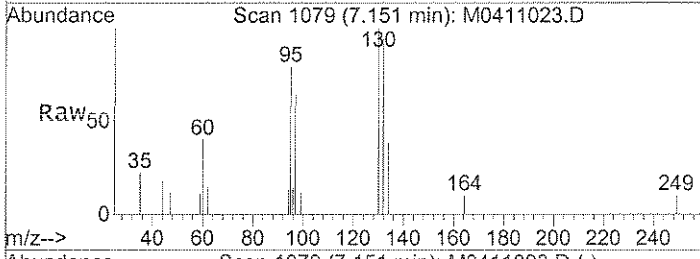
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	7.15	130	2629	0.74 ug/l	86
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d
52) Toluene	8.48	92	223	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	8.78	69	63	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	9.02	166	349	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	9.87	91	1150	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.97	91	383	N.D.	
65) m,p-Xylene	10.10	106	293	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	d
68) Bromoform	0.00	173	0	N.D.	
69) Isopropylbenzene	11.03	105	272	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	0.00	156	0	N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	222	N.D.	
75) 1,2,3-Trichloropropane	11.03	110	90	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	
77) 2-Chlorotoluene	0.00	91	0	N.D.	
78) 4-Chlorotoluene	0.00	91	0	N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0	N.D.	
80) tert-Butylbenzene	0.00	119	0	N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0	N.D.	
82) sec-butylbenzene	0.00	105	0	N.D.	d
83) 1,3-Dichlorobenzene	0.00	146	0	N.D.	d
84) 4-Isopropyltoluene	0.00	119	0	N.D.	d
85) 1,4-Dichlorobenzene	11.97	146	69	N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0	N.D.	
87) n-Butylbenzene	12.52	91	60	N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0	N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.	
90) Hexachlorobutadiene	0.00	225	0	N.D.	
91) Naphthalene	0.00	128	0	N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0	N.D.	

*MS 4/12/07*



#41  
 Trichloroethene  
 Concen: 0.74 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0411023.D  
 Acq: 11 Apr 2007 16:05

Tgt Ion	Resp	Lower	Upper
130	2629		
130	100		
132	109.0	75.0	115.0
95	102.4	69.4	109.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-009  
 Lab File ID: M0411024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:28  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.93	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	2.4	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.2	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-009  
 Lab File ID: M0411024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:28  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.31	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-009  
 Lab File ID: M0411024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/11/2007 16:28  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

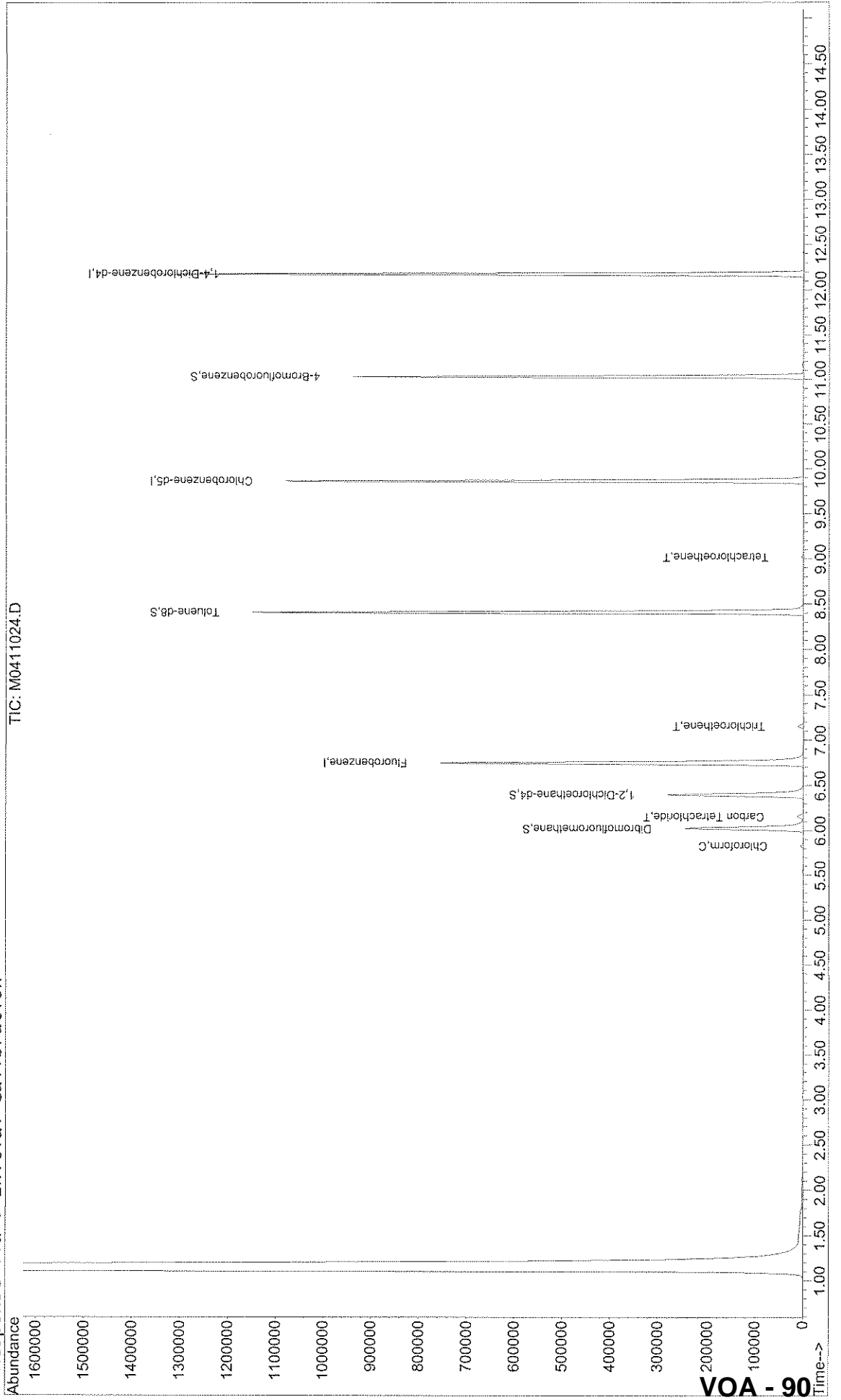
Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411024.D  
Acq On : 11 Apr 2007 16:28  
Sample : JPL29-009  
Misc : #2 5ml +IS/ss (524)  
MS Integration Params: rteint.p  
Quant Time: Apr 12 9:28 2007

Vial: 64  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 524.RES

Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Last Update : Tue Apr 10 10:47:40 2007  
Response via : Initial Calibration



VOA - 90



Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411024.D  
 Acq On : 11 Apr 2007 16:28  
 Sample : JPL29-009  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:28 2007

Vial: 64  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\040907\M0409012.D (9 Apr 2007 12:54)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar)
1) Fluorobenzene	6.75	96	757339	50.00	ug/l	0.00 72.32%
50) Chlorobenzene-d5	9.87	82	323016	50.00	ug/l	0.00 71.30%
70) 1,4-Dichlorobenzene-d4	12.18	152	317598	50.00	ug/l	0.00 67.96%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	179896	49.92	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	99.84%		
37) 1,2-Dichloroethane-d4	6.40	65	221938	53.39	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	106.78%		
51) Toluene-d8	8.41	98	769351	51.24	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	102.48%		
72) 4-Bromofluorobenzene	11.03	95	266804	51.36	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery =	102.72%		

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.98	43	56	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.57	63	827	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.54	43	57	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.83	83	5597	0.93 ug/l	✓	97
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	6.15	117	9544	2.39 ug/l	✓	91
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.42	78	160	N.D.		

*W 4/12/07*

Quantitation Report

Data File : X:\MSVOA\MOBY\041107A\M0411024.D  
 Acq On : 11 Apr 2007 16:28  
 Sample : JPL29-009  
 Misc : #2 5ml +IS/SS (524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 12 9:28 2007

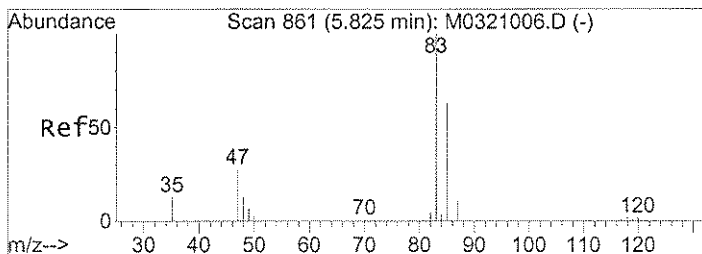
Vial: 64  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 524.RES

Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
 Title : VOA 524- 5ML Calibration 5973M  
 Last Update : Tue Apr 10 10:47:40 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

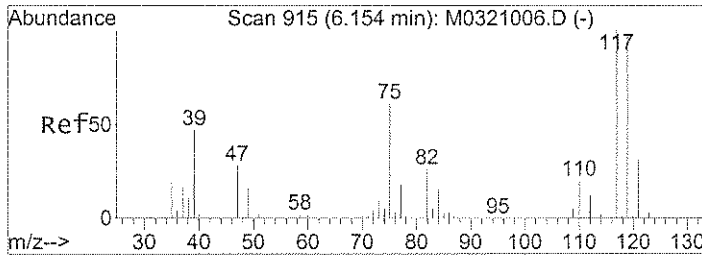
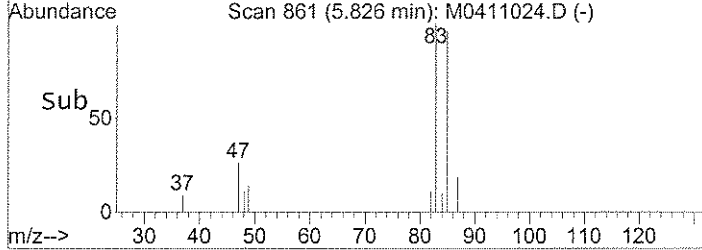
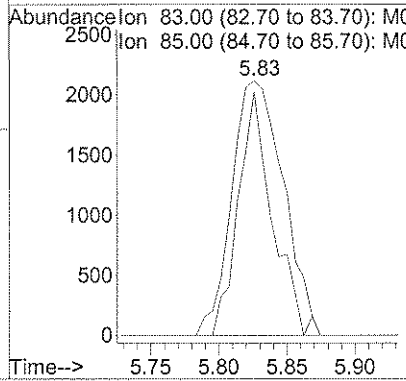
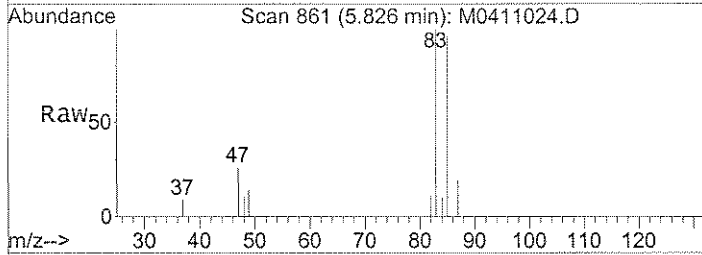
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
39) 1,2-Dichloroethane	6.49	62	205		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	4423	1.22	ug/l	96
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.72	83	213		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.47	92	390		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	1120	0.31	ug/l	89
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.98	91	298		N.D.	
65) m,p-Xylene	10.10	106	492		N.D.	
66) o-xylene	10.50	106	55		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	382		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	88		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	145		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.26	91	80		N.D.	
78) 4-Chlorotoluene	11.26	91	80		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D. d	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D. d	
84) 4-Isopropyltoluene	0.00	119	0		N.D. d	
85) 1,4-Dichlorobenzene	12.21	146	302		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	316		N.D.	
87) n-Butylbenzene	12.52	91	56		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D. d	
92) 1,2,3-Trichlorobenzene	14.65	180	380		N.D.	

*W 4/12/07*



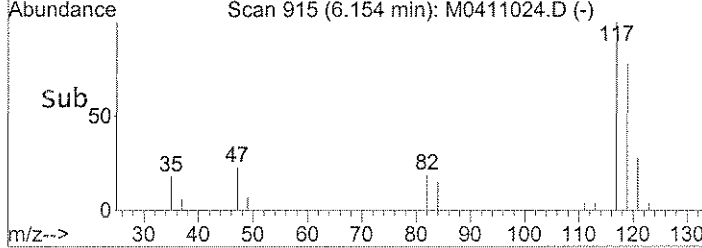
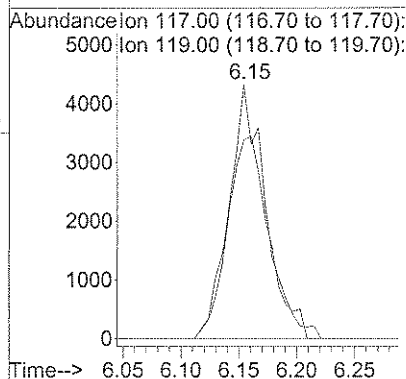
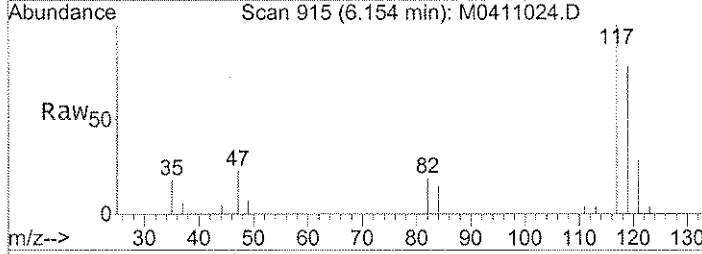
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 Chloroform  
 Concen: 0.93 ug/l  
 RT: 5.83 min Scan# 861  
 Delta R.T. -0.00 min  
 Lab File: M0411024.D  
 Acq: 11 Apr 2007 16:28

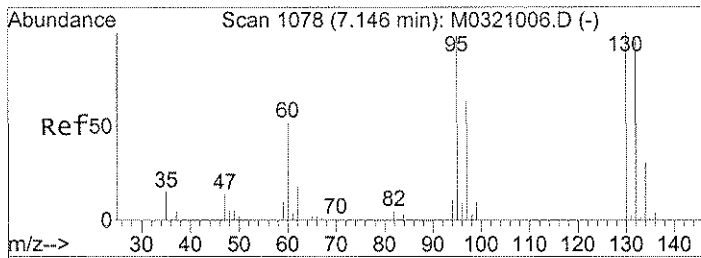
Tgt Ion: 83 Resp: 5597  
 Ion Ratio Lower Upper  
 83 100  
 85 63.2 41.2 81.2



#35  
 Carbon Tetrachloride  
 Concen: 2.39 ug/l  
 RT: 6.15 min Scan# 915  
 Delta R.T. -0.00 min  
 Lab File: M0411024.D  
 Acq: 11 Apr 2007 16:28

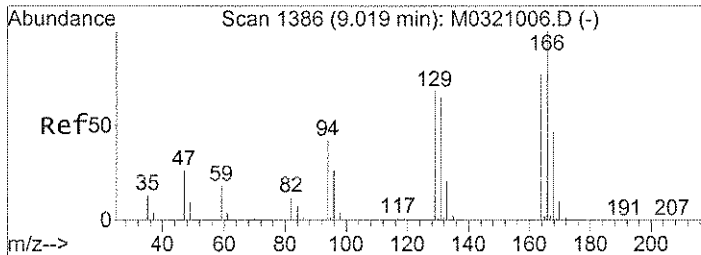
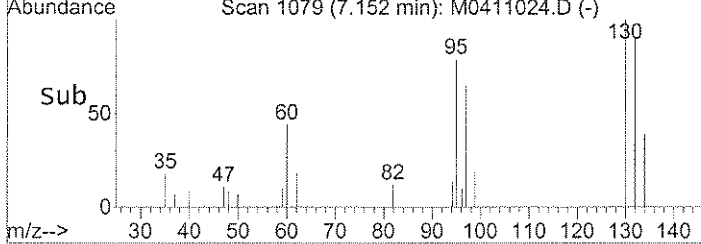
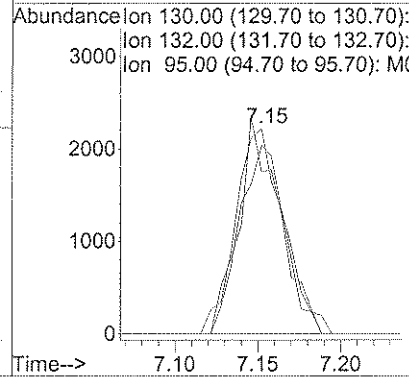
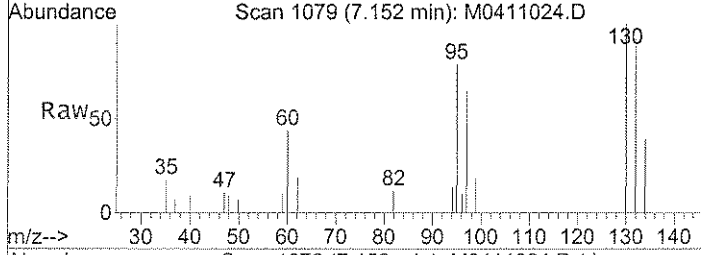
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 Ion Ratio Lower Upper  
 117 100  
 119 89.2 78.2 118.2





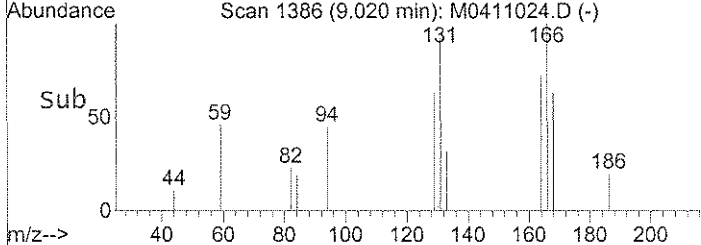
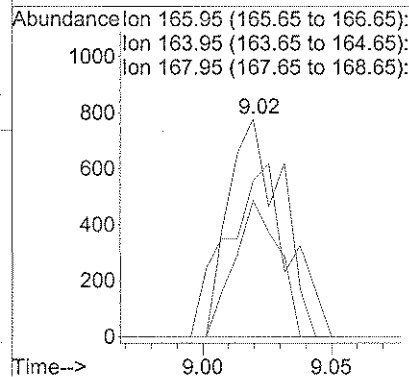
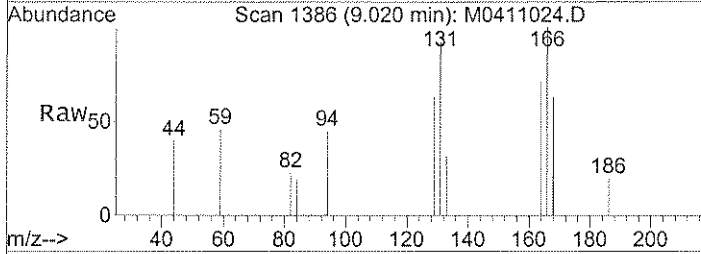
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 Trichloroethene  
 Concen: 1.22 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0411024.D  
 Acq: 11 Apr 2007 16:28

Tgt Ion	Resp	Lower	Upper
130	4423		
132	89.6	75.0	115.0
95	91.7	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 0.31 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. -0.00 min  
 Lab File: M0411024.D  
 Acq: 11 Apr 2007 16:28

Tgt Ion	Resp	Lower	Upper
166	1120		
164	92.6	63.3	94.9
168	52.2	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL29-010  
 Lab File ID: Y0413013.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/13/2007 11:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.27	J
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.60	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.3	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL29-010  
 Lab File ID: Y0413013.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/13/2007 11:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.74	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-17-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL29-010  
 Lab File ID: Y0413013.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/13/2007 11:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

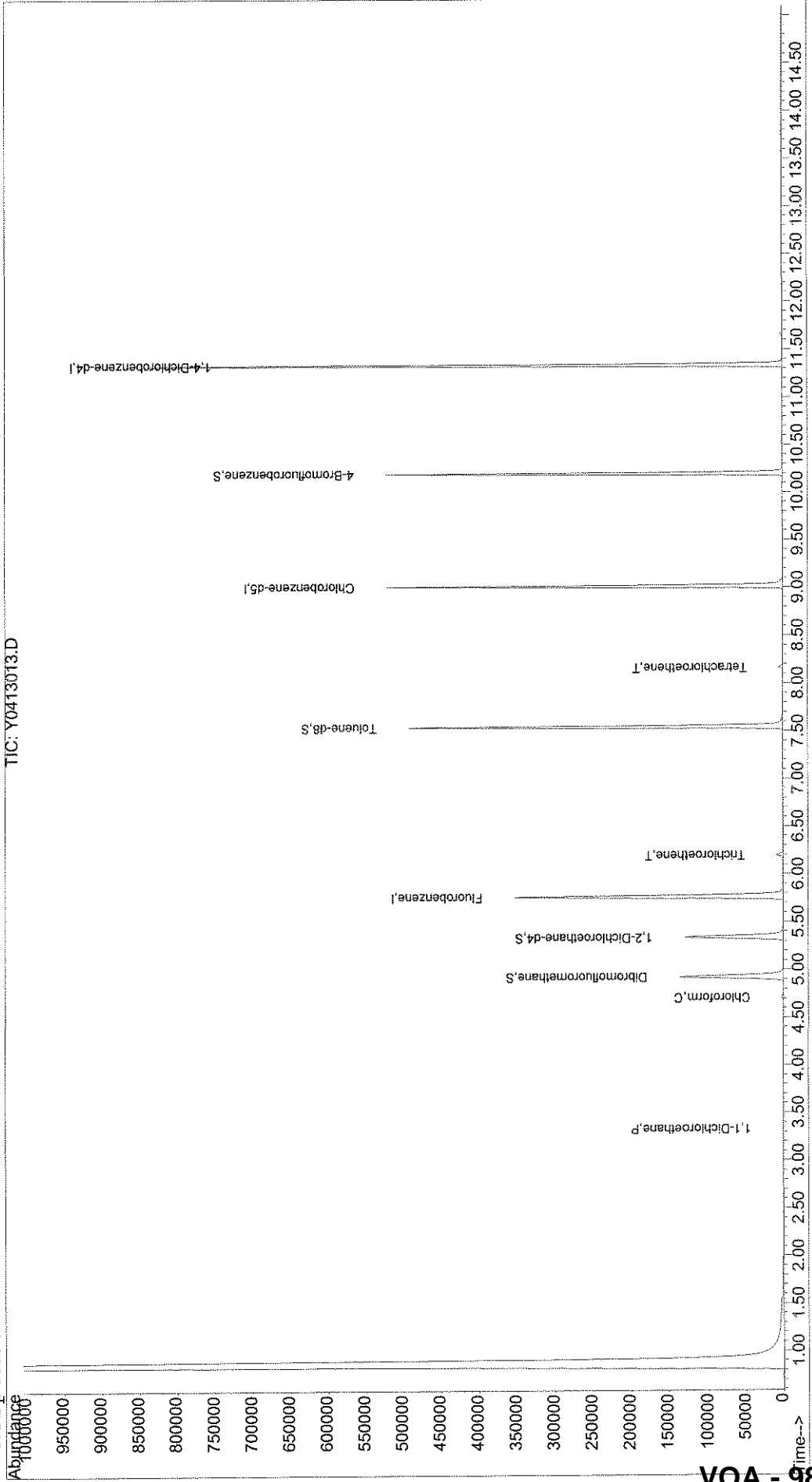
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413013.D Vial: 28  
Acq On : 13 Apr 2007 11:12 Operator: LNH  
Sample : JPL29-010 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 13 13:39 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration





Data File : X:\MSVOA\YODA\041307\Y0413013.D  
 Acq On : 13 Apr 2007 11:12  
 Sample : JPL29-010  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 13:39 2007

Vial: 28  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Fluorobenzene	5.76	96	324485	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	135585	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	183793	50.00	ug/l	0.00

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	100614	51.25	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	102.50%	
37) 1,2-Dichloroethane-d4	5.34	65	103369	49.63	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	99.26%	
51) Toluene-d8	7.54	98	294173	54.67	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	109.34%	
72) 4-Bromofluorobenzene	10.20	95	137537	51.80	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.33	63	1231	0.27	ug/l #	60
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		

LN 4/13/07

Data File : X:\MSVOA\YODA\041307\Y0413013.D

Vial: 28

Acq On : 13 Apr 2007 11:12

Operator: LNH

Sample : JPL29-010

Inst : yoda

Misc : 5mL+IS/SS #1

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 8260B.RES

Quant Time: Apr 13 13:39 2007

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

Response via : Initial Calibration

DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	4.71	83	2491	0.60	ug/l	96
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	5.11	117	710	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	5.39	78	65	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.		
41) Trichloroethene	6.20	130	2996	1.31	ug/l	98
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	0.00	41	0	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	0.00	92	0	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.		
56) Tetrachloroethene	8.16	166	1384	0.74	ug/l	93
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.		
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	0.00	112	0	N.D.		
62) 1-Chlorohexane	9.01	91	416	N.D.		
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.17	91	63	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.		
66) o-xylene	0.00	106	0	N.D.		
67) Styrene	0.00	104	0	N.D.		
68) Bromoform	0.00	173	0	N.D.		
69) Isopropylbenzene	10.19	105	54	N.D.		
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.		
73) Bromobenzene	0.00	156	0	N.D.		
74) 1,1,2,2-Tetrachloroethane	0.00	83	0	N.D.		
75) 1,2,3-Trichloropropane	0.00	110	0	N.D.		
76) n-Propylbenzene	0.00	120	0	N.D.		

(#)= qualifier out of range (m) = manual integration

Y0413013.D 8260B.M Fri Apr 13 13:39:55 2007

LU 4/13/07

Data File : X:\MSVOA\YODA\041307\Y0413013.D

Vial: 28

Acq On : 13 Apr 2007 11:12

Operator: LNH

Sample : JPL29-010

Inst : yoda

Misc : 5mL+IS/SS #1

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Time: Apr 13 13:39 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

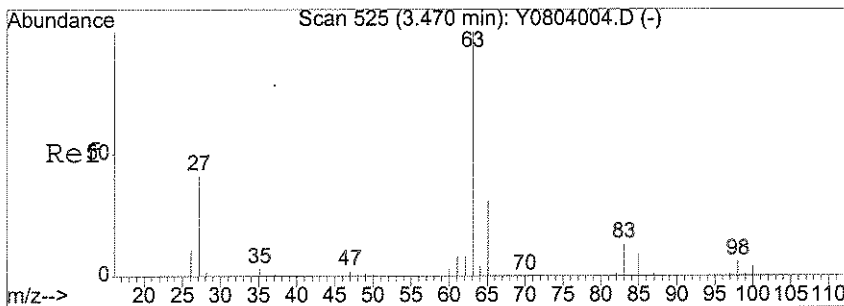
Response via : Initial Calibration

DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.47	91	233		N.D.	
78) 4-Chlorotoluene	10.64	91	62		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	60		N.D.	
82) sec-butylbenzene	11.18	105	121		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	327		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	203		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	213		N.D.	

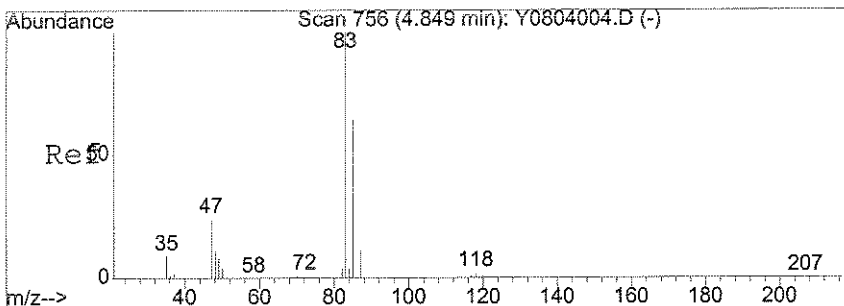
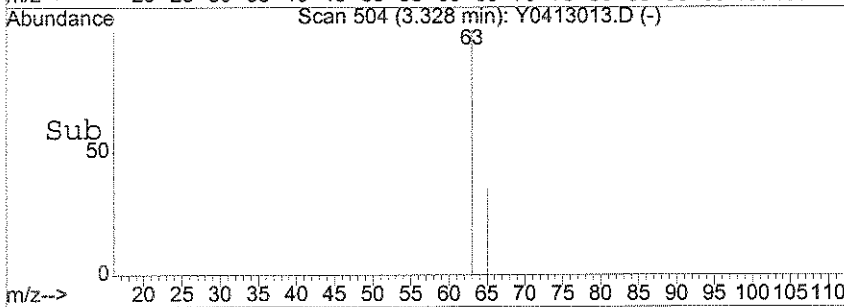
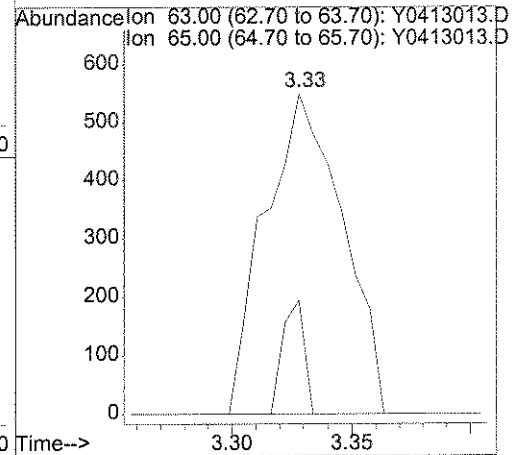
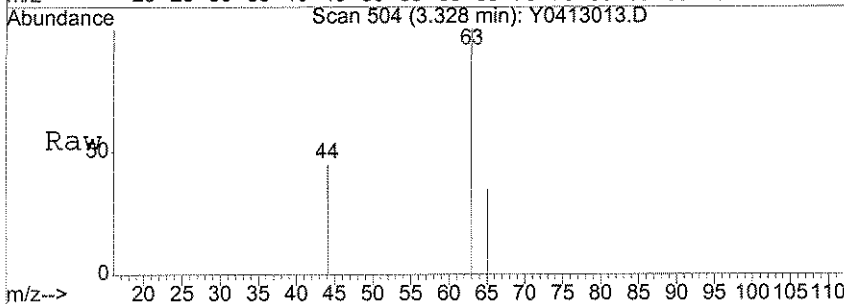
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(#) = qualifier out of range (m) = manual integration

Y0413013.D 8260B.M Fri Apr 13 13:39:55 2007



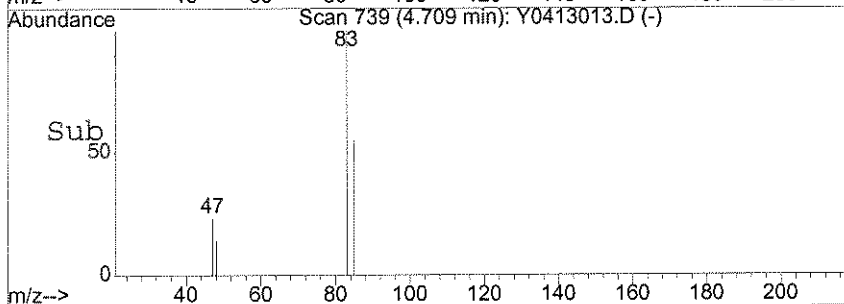
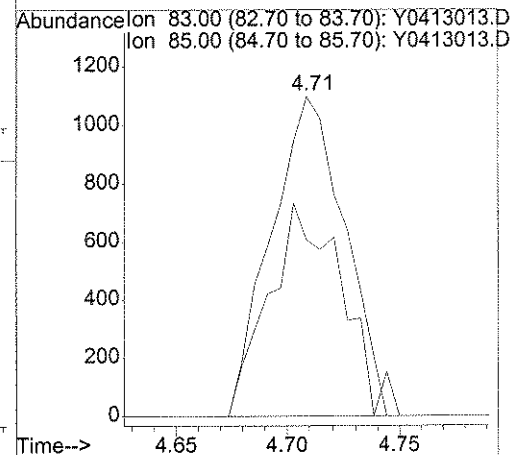
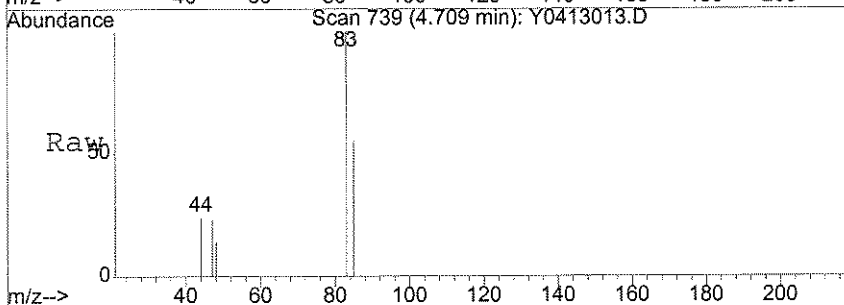
#22  
 1,1-Dichloroethane  
 Concen: 0.27 ug/l  
 RT: 3.33 min Scan# 504  
 Delta R.T. 0.01 min  
 Lab File: Y0413013.D  
 Acq: 13 Apr 2007 11:12

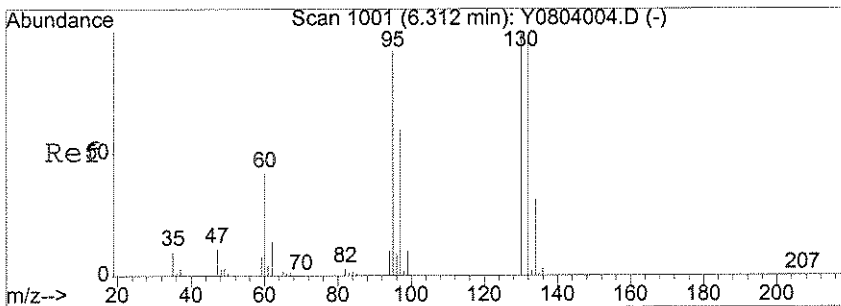
Tgt Ion	Resp	Lower	Upper
63	1231		
63	100		
65	10.1	12.3	52.3#



#31  
 Chloroform  
 Concen: 0.60 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0413013.D  
 Acq: 13 Apr 2007 11:12

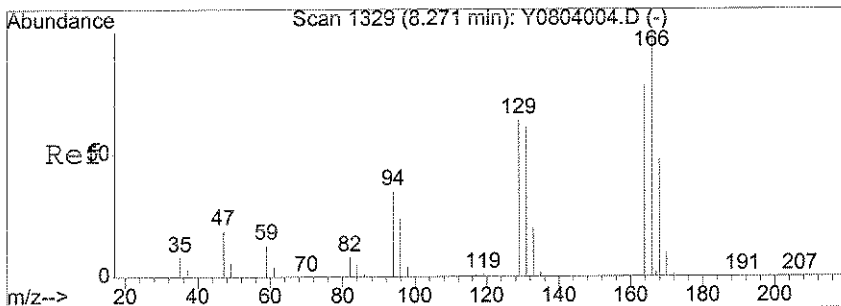
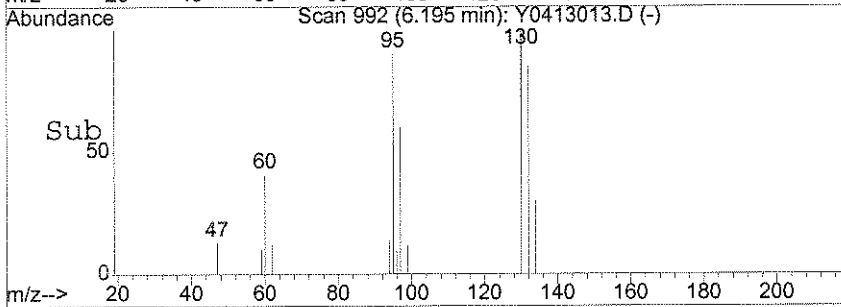
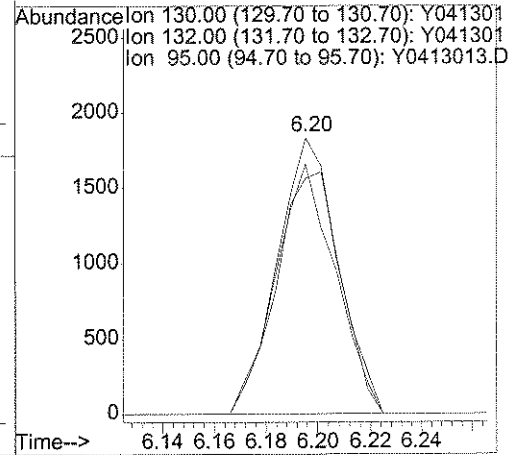
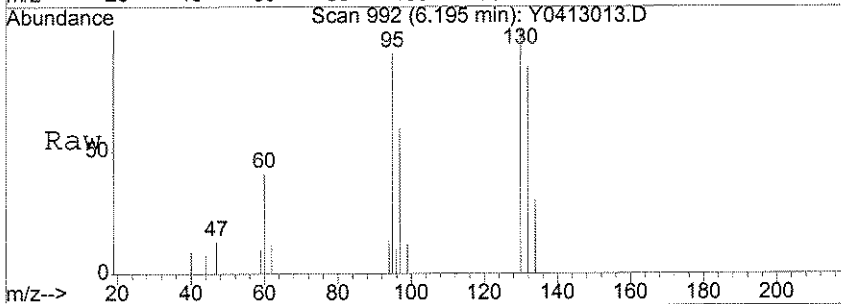
Tgt Ion	Resp	Lower	Upper
83	2491		
83	100		
85	66.2	43.3	83.3





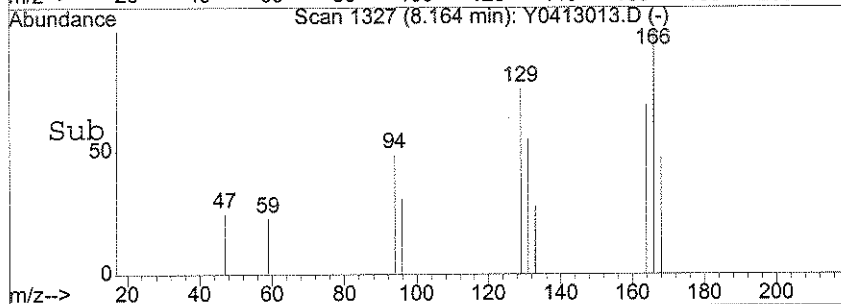
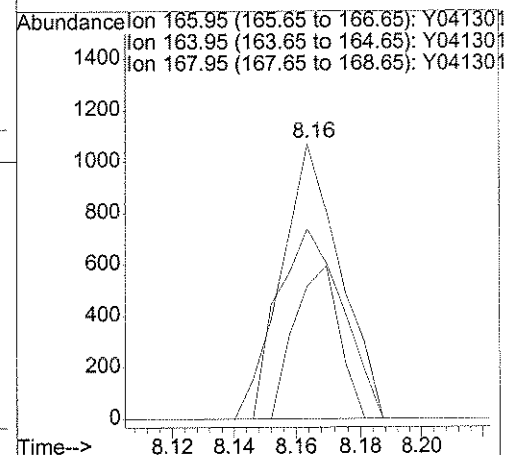
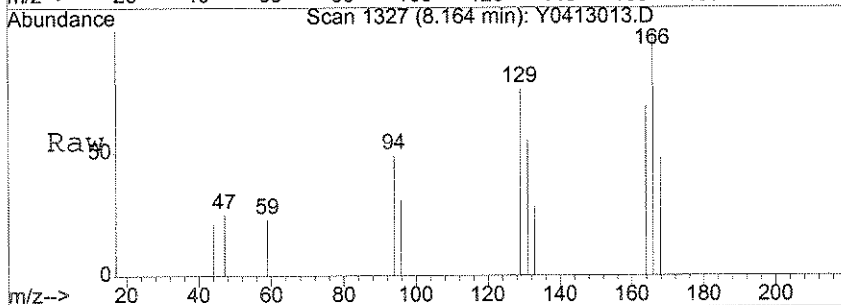
#41  
 Trichloroethene  
 Concen: 1.31 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0413013.D  
 Acq: 13 Apr 2007 11:12

Tgt Ion	Resp	Lower	Upper
130	2996		
130	100		
132	91.9	75.0	115.0
95	88.2	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 0.74 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. -0.00 min  
 Lab File: Y0413013.D  
 Acq: 13 Apr 2007 11:12

Tgt Ion	Resp	Lower	Upper
166	1384		
166	100		
164	75.4	63.3	94.9
168	42.1	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-1-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-011  
 Lab File ID: Y0412024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:29  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.59	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.4	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-1-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-011  
 Lab File ID: Y0412024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:29  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.75	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-1-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-011  
 Lab File ID: Y0412024.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:29  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

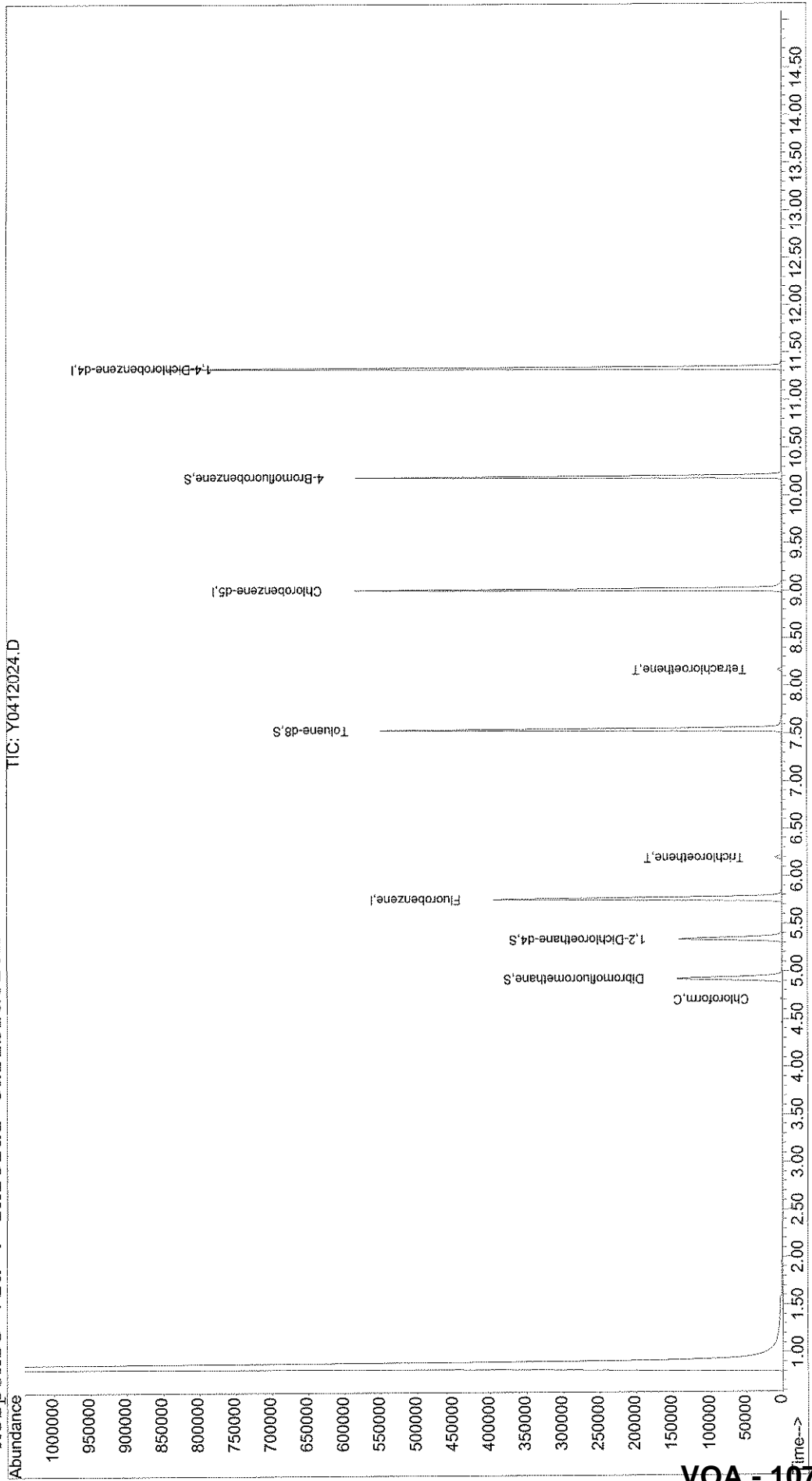
Comments:



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412024.D Vial: 32  
Acq On : 12 Apr 2007 15:29 Operator: LPM  
Sample : JPL29-011 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:04 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412024.D  
 Acq On : 12 Apr 2007 15:29  
 Sample : JPL29-011  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:04 2007

Vial: 32  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.76	96	361916	50.00	ug/l	0.00	NA%
50) Chlorobenzene-d5	9.01	82	150312	50.00	ug/l	0.00	NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	193608	50.00	ug/l	0.00	NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	108417	49.51	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.02%	
37) 1,2-Dichloroethane-d4	5.34	65	113149	48.71	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	97.42%	
51) Toluene-d8	7.54	98	333555	55.92	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	111.84%	
72) 4-Bromofluorobenzene	10.20	95	152240	54.43	ug/l	0.00	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.33	63	1177	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412024.D  
 Acq On : 12 Apr 2007 15:29  
 Sample : JPL29-011  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:04 2007

Vial: 32  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.70	83	2733	0.59 ug/l ✓	98
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	5.12	117	537	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	5.38	78	56	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	6.20	130	3476	1.36 ug/l ✓	96
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	0.00	92	0	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	8.16	166	1561	0.75 ug/l ✓	97
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	9.01	91	450	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.16	91	53	N.D.	
65) m,p-Xylene	9.30	106	220	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D.	

*Handwritten signature and date: 4/13/07*

(#) = qualifier out of range (m) = manual integration  
 Y0412024.D 8260B.M Fri Apr 13 10:07:36 2007

Quantitation Report

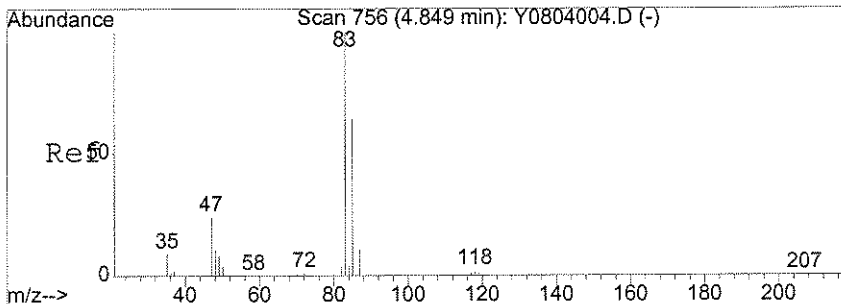
Data File : X:\MSVOA\YODA\041207\Y0412024.D  
 Acq On : 12 Apr 2007 15:29  
 Sample : JPL29-011  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:04 2007

Vial: 32  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

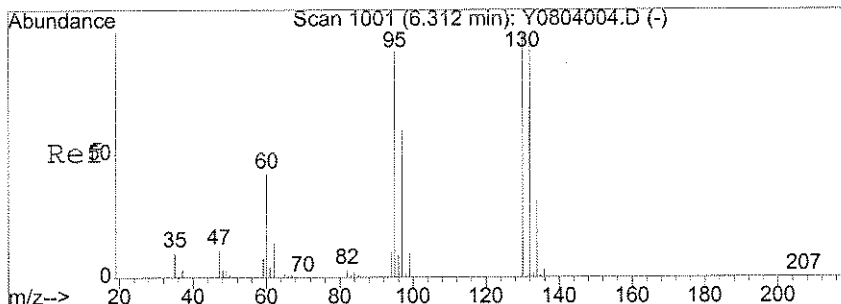
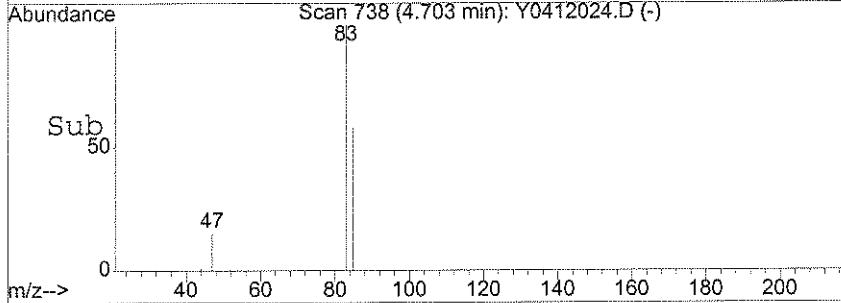
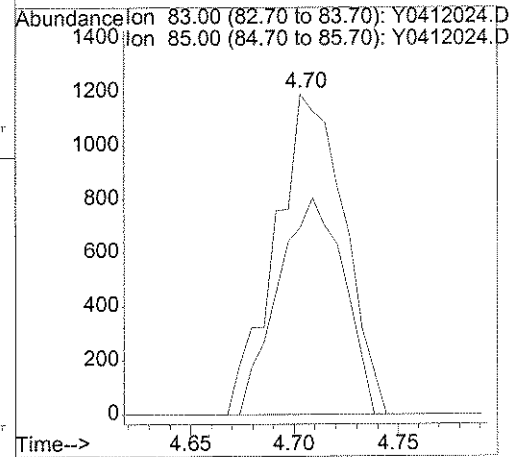
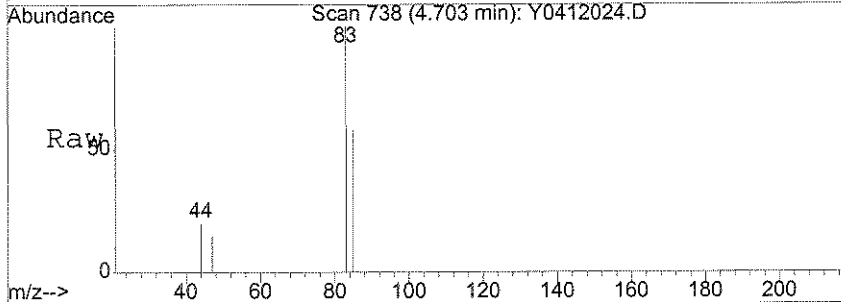
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	65		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.53	91	62		N.D.	
78) 4-Chlorotoluene	10.64	91	61		N.D.	
79) 1,3,5-Trimethylbenzene	10.66	105	192		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	137		N.D.	
82) sec-butylbenzene	11.18	105	303		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	108		N.D.	
84) 4-Isopropyltoluene	11.33	119	481		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	162		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	58		N.D.	
87) n-Butylbenzene	11.74	91	359		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.30	180	72		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	222		N.D.	



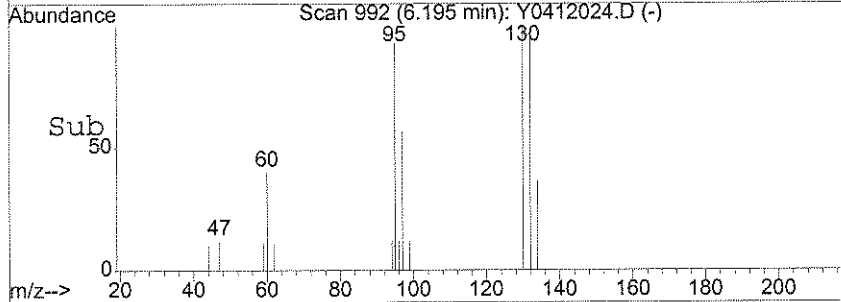
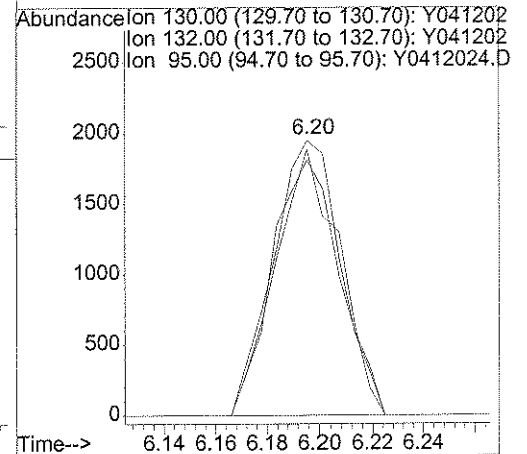
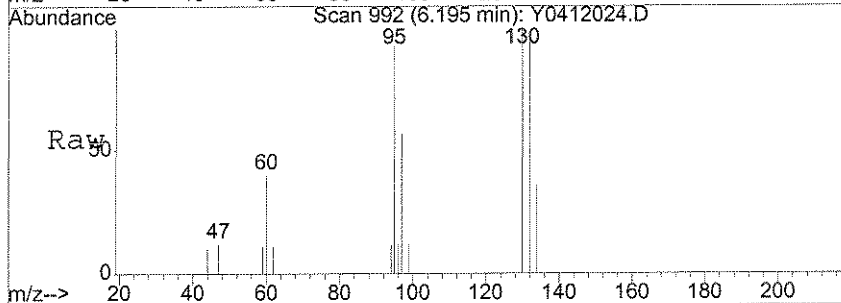
#31  
 Chloroform  
 Concen: 0.59 ug/l  
 RT: 4.70 min Scan# 738  
 Delta R.T. -0.00 min  
 Lab File: Y0412024.D  
 Acq: 12 Apr 2007 15:29

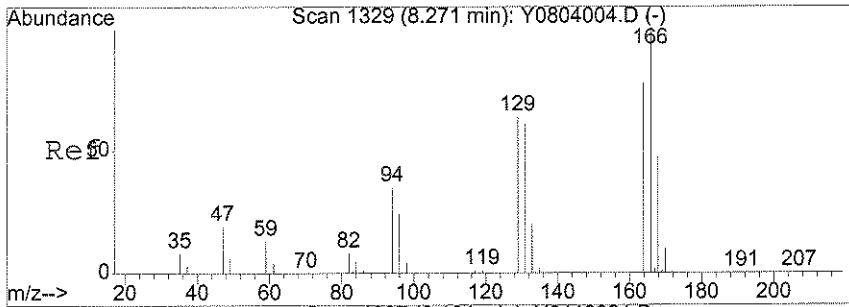
Tgt Ion:	83	Resp:	2733
Ion Ratio	Lower	Upper	
83	100		
85	64.7	43.3	83.3



#41  
 Trichloroethene  
 Concen: 1.36 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0412024.D  
 Acq: 12 Apr 2007 15:29

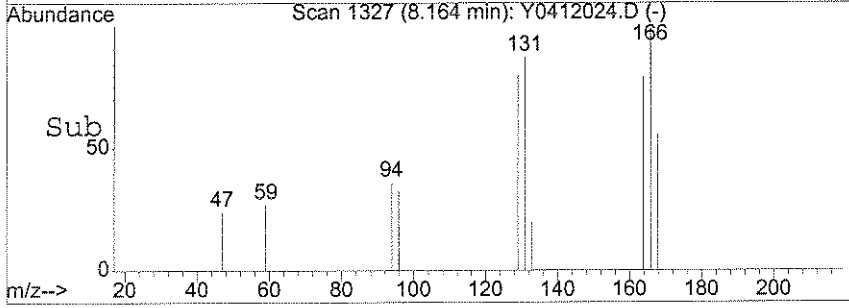
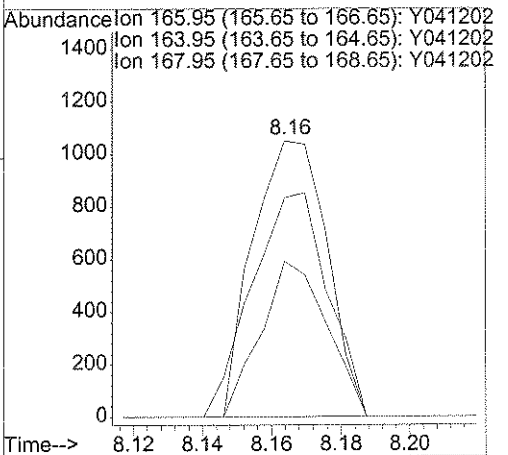
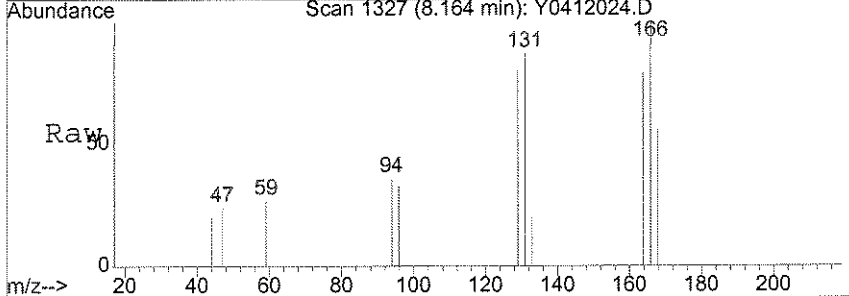
Tgt Ion:	130	Resp:	3476
Ion Ratio	Lower	Upper	
130	100		
132	91.6	75.0	115.0
95	93.3	69.4	109.4





#56  
 Tetrachloroethene  
 Concen: 0.75 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. -0.00 min  
 Lab File: Y0412024.D  
 Acq: 12 Apr 2007 15:29

Tgt Ion	Resp	Lower	Upper
166	1561		
166	100		
164	83.0	63.3	94.9
168	50.2	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-012  
 Lab File ID: Y0412025.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:53  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-012  
 Lab File ID: Y0412025.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:53  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.76	J
95-47-6	o-Xylene	0.28	J
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-012  
 Lab File ID: Y0412025.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 15:53  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

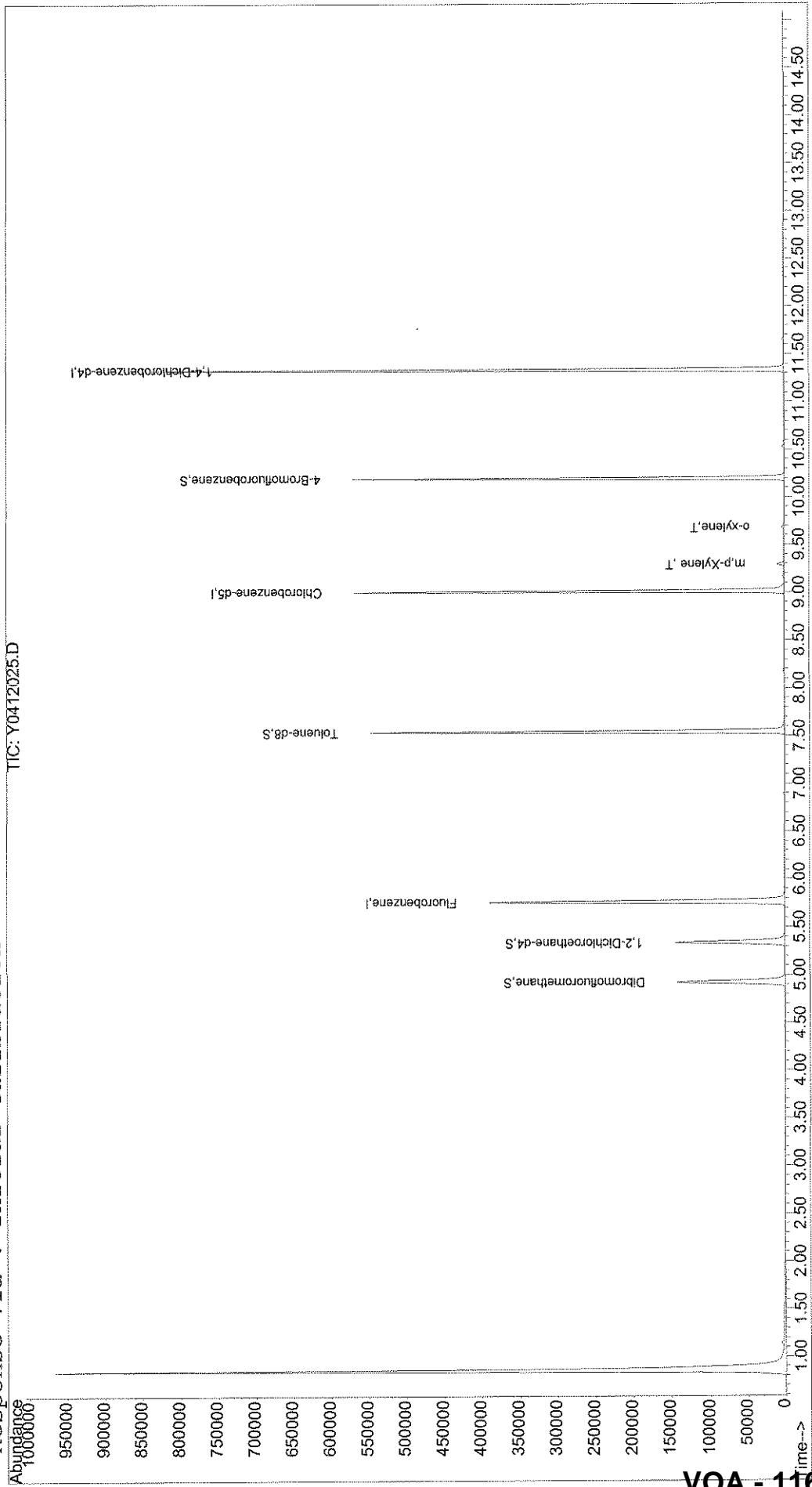
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412025.D  
Acq On : 12 Apr 2007 15:53  
Sample : JPL29-012  
Misc : 5mL+IS/SS #2  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:05 2007  
Vial: 33  
Operator: LPM  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 116

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412025.D  
 Acq On : 12 Apr 2007 15:53  
 Sample : JPL29-012  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:05 2007

Vial: 33  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	367227	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	143940	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	190698	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	107711	48.48	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	96.96%
37) 1,2-Dichloroethane-d4	5.34	65	113548	48.17	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	96.34%
51) Toluene-d8	7.54	98	331184	57.98	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	115.96%
72) 4-Bromofluorobenzene	10.20	95	145775	52.91	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0412025.D 8260B.M Fri Apr 13 10:08:33 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412025.D  
 Acq On : 12 Apr 2007 15:53  
 Sample : JPL29-012  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:05 2007

Vial: 33  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	4.34	43	66		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.72	83	55		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.40	78	59		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	423		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	1820		N.D.	
65) m,p-Xylene	9.29	106	2453	0.76	ug/l ✓	90
66) o-xylene	9.68	106	892	0.28	ug/l ✓	98
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*QNH 4/13/07*

(#) = qualifier out of range (m) = manual integration  
 Y0412025.D 8260B.M Fri Apr 13 10:08:34 2007

Quantitation Report

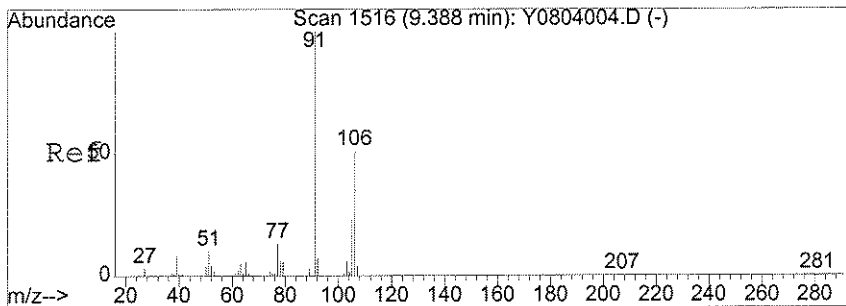
Data File : X:\MSVOA\YODA\041207\Y0412025.D  
 Acq On : 12 Apr 2007 15:53  
 Sample : JPL29-012  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:05 2007

Vial: 33  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

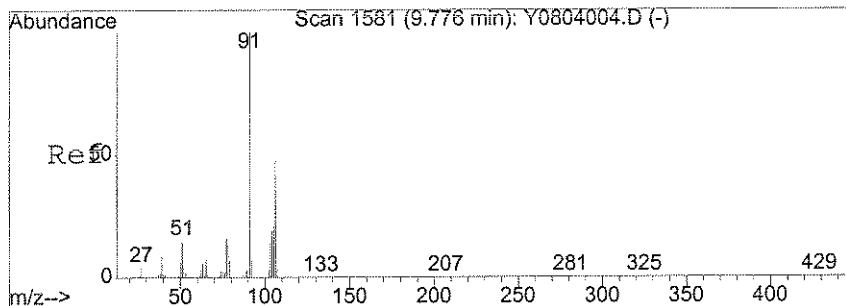
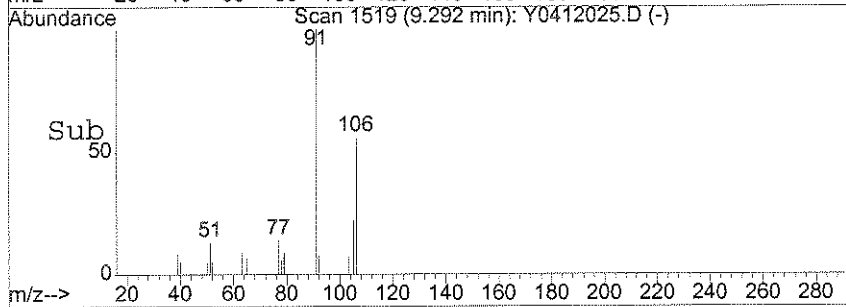
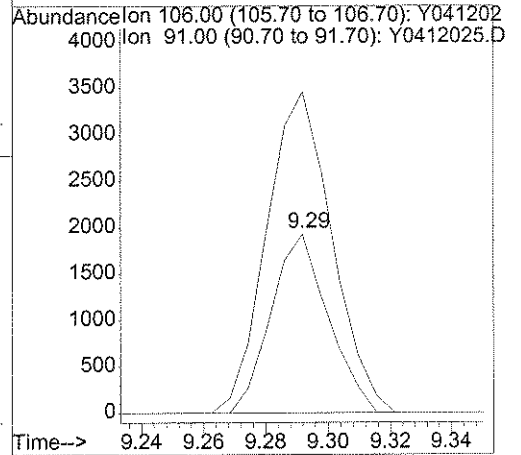
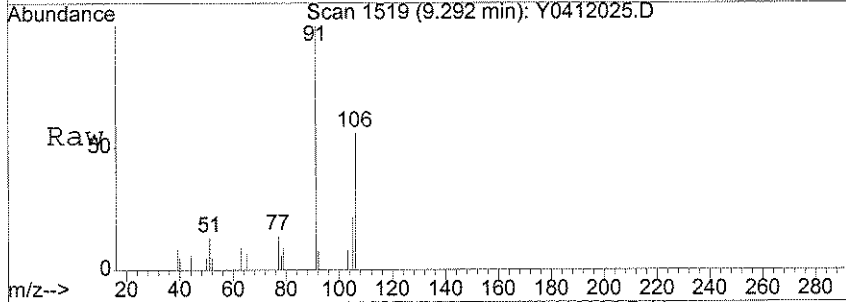
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	183		N.D.	
78) 4-Chlorotoluene	10.46	91	183		N.D.	
79) 1,3,5-Trimethylbenzene	10.66	105	59		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	248		N.D.	
82) sec-butylbenzene	11.18	105	251		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	350		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	390		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	13.55	128	336		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



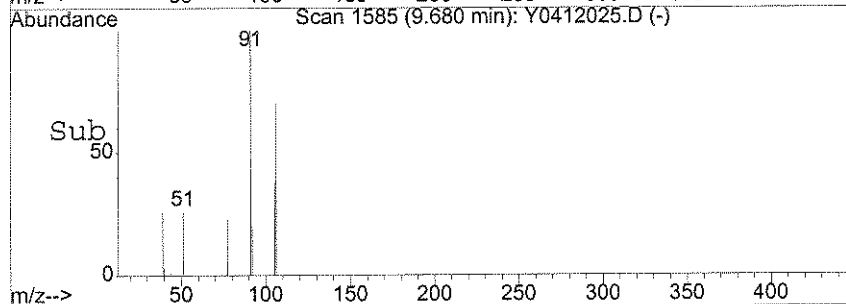
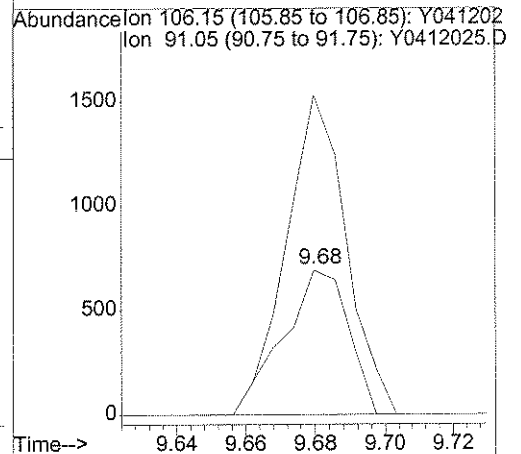
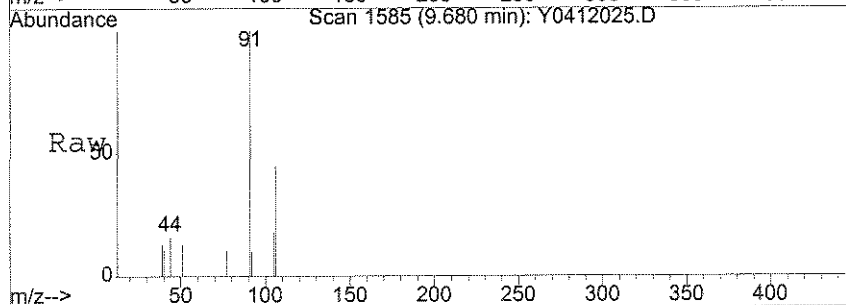
#65  
 m,p-Xylene  
 Concen: 0.76 ug/l  
 RT: 9.29 min Scan# 1519  
 Delta R.T. 0.00 min  
 Lab File: Y0412025.D  
 Acq: 12 Apr 2007 15:53

Tgt Ion	Resp	Lower	Upper
106	2453		
91	204.9	169.6	209.6



#66  
 o-xylene  
 Concen: 0.28 ug/l  
 RT: 9.68 min Scan# 1585  
 Delta R.T. 0.00 min  
 Lab File: Y0412025.D  
 Acq: 12 Apr 2007 15:53

Tgt Ion	Resp	Lower	Upper
106	892		
91	203.8	160.6	240.8



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-013  
 Lab File ID: Y0412026.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:18  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-013  
 Lab File ID: Y0412026.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:18  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-013  
 Lab File ID: Y0412026.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:18  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

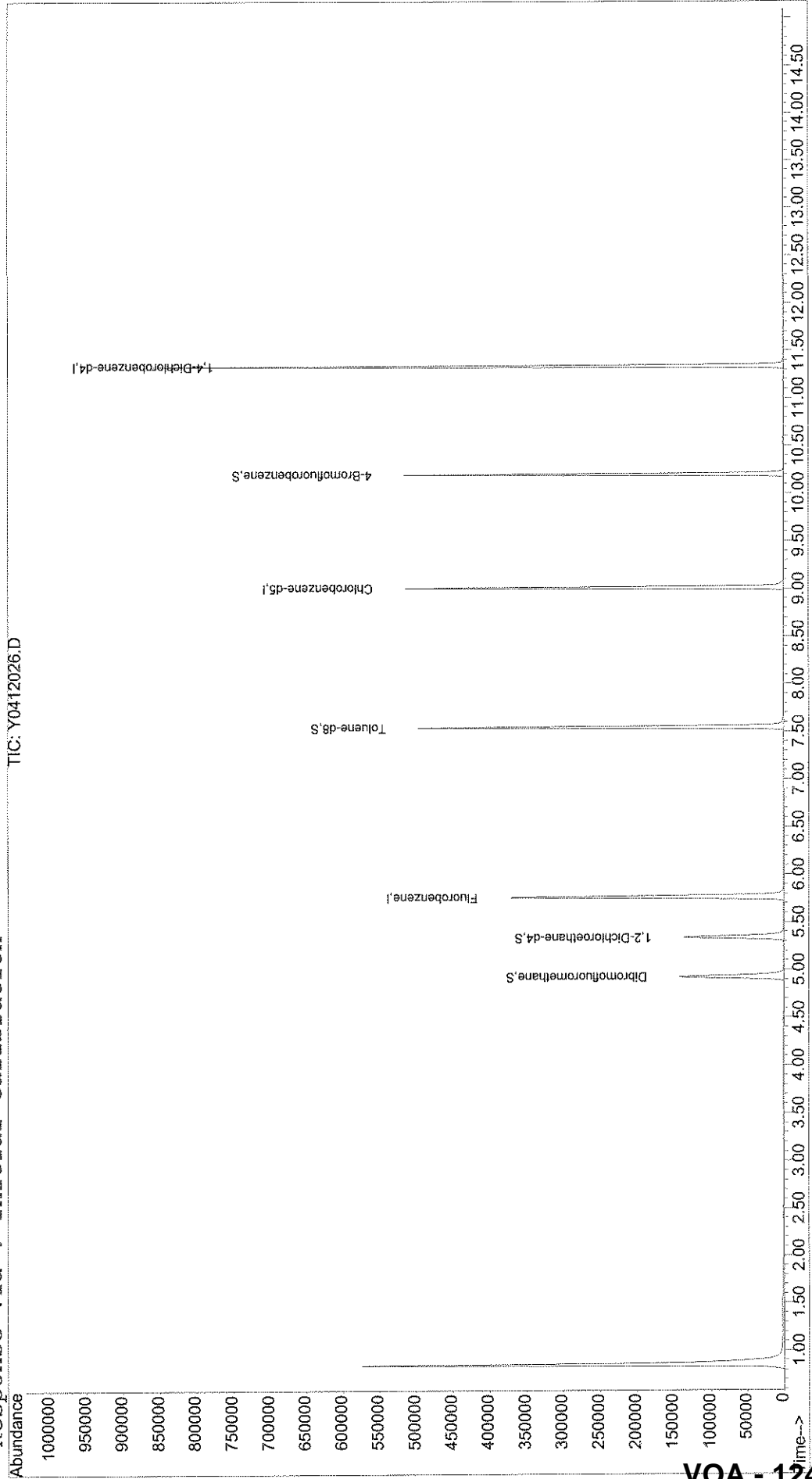
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412026.D  
Acq On : 12 Apr 2007 16:18  
Sample : JPL29-013  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:06 2007  
Vial: 34  
Operator: LPM  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 124

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412026.D  
 Acq On : 12 Apr 2007 16:18  
 Sample : JPL29-013  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Vial: 34  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	339168	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	132511	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	185589	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	104242	50.80	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.60%
37) 1,2-Dichloroethane-d4	5.34	65	106210	48.79	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	97.58%
51) Toluene-d8	7.54	98	295958	56.28	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.56%
72) 4-Bromofluorobenzene	10.20	95	136415	50.88	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	131	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	2.52	84	275	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0412026.D 8260B.M Fri Apr 13 10:08:48 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412026.D  
 Acq On : 12 Apr 2007 16:18  
 Sample : JPL29-013  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Vial: 34  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	392		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	186		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412026.D 8260B.M Fri Apr 13 10:08:49 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412026.D  
 Acq On : 12 Apr 2007 16:18  
 Sample : JPL29-013  
 Misc : 5mL+IS/SS #1

Vial: 34  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.19	105	206		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	138		N.D.	
78) 4-Chlorotoluene	10.46	91	138		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	60		N.D.	
82) sec-butylbenzene	11.18	105	72		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	333		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	389		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	88		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-014  
 Lab File ID: Y0412027.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-014  
 Lab File ID: Y0412027.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-014  
 Lab File ID: Y0412027.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 16:43  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

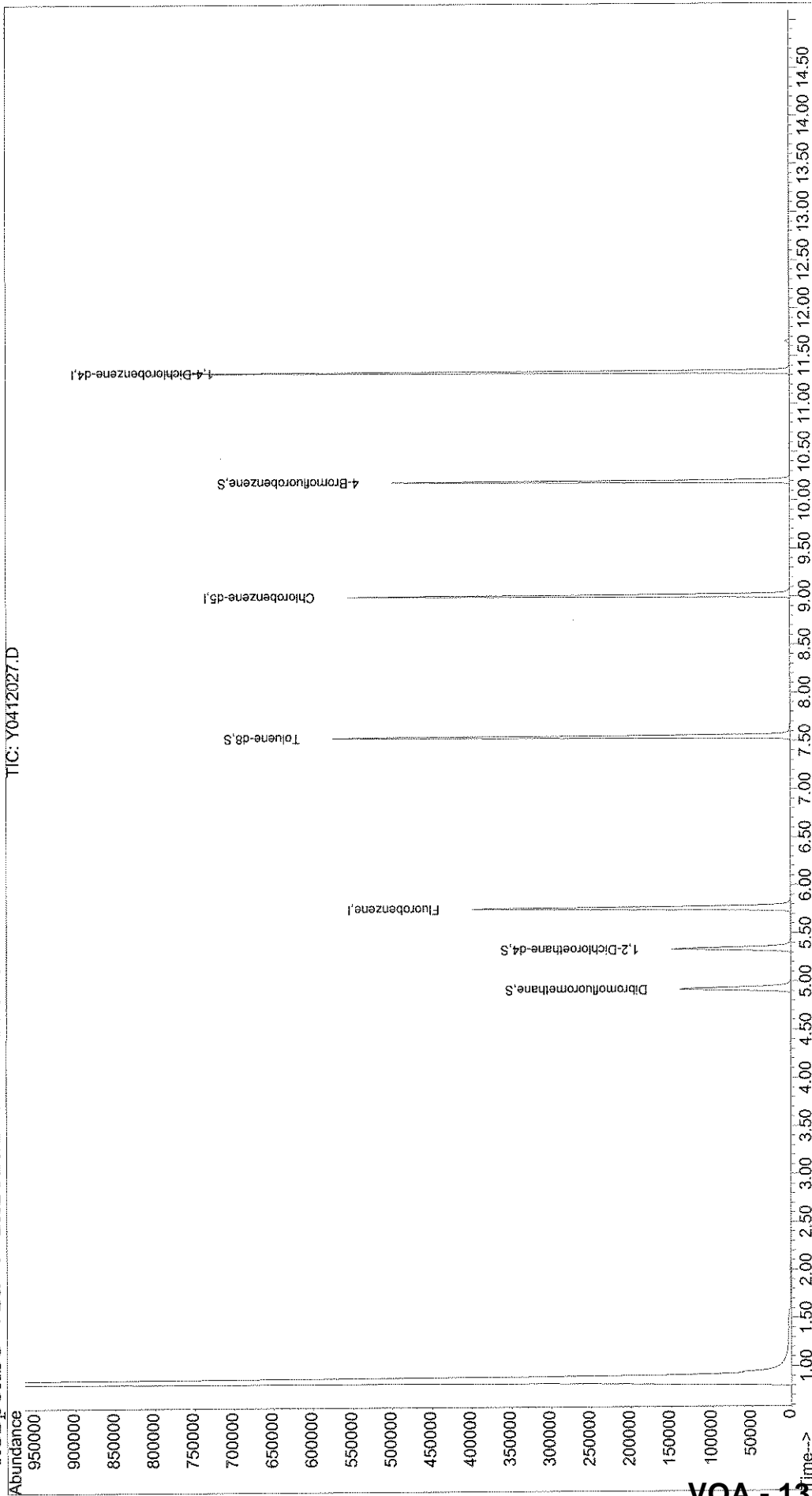
Comments:



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412027.D  
Acq On : 12 Apr 2007 16:43  
Sample : JPL29-014  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:06 2007  
Vial: 35  
Operator: LPM  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412027.D  
 Acq On : 12 Apr 2007 16:43  
 Sample : JPL29-014  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Vial: 35  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	373327	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	138672	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	176923	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	103300	45.73	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	91.46%
37) 1,2-Dichloroethane-d4	5.34	65	114784	47.90	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.80%
51) Toluene-d8	7.54	98	340974	61.96	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	123.92%#
72) 4-Bromofluorobenzene	10.20	95	130663	51.12	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	133		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	d
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412027.D 8260B.M Fri Apr 13 10:08:58 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412027.D  
 Acq On : 12 Apr 2007 16:43  
 Sample : JPL29-014  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Vial: 35  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	59		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.62	92	56		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	430		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	260		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	377		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412027.D 8260B.M Fri Apr 13 10:08:59 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412027.D  
 Acq On : 12 Apr 2007 16:43  
 Sample : JPL29-014  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:06 2007

Vial: 35  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	137		N.D.	
78) 4-Chlorotoluene	10.46	91	137		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	131		N.D.	
82) sec-butylbenzene	11.18	105	71		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	316		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	339		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-015  
 Lab File ID: Y0412028.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:08  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	1.6	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	7.1	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.1	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-015  
 Lab File ID: Y0412028.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:08  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.55	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-015  
 Lab File ID: Y0412028.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:08  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

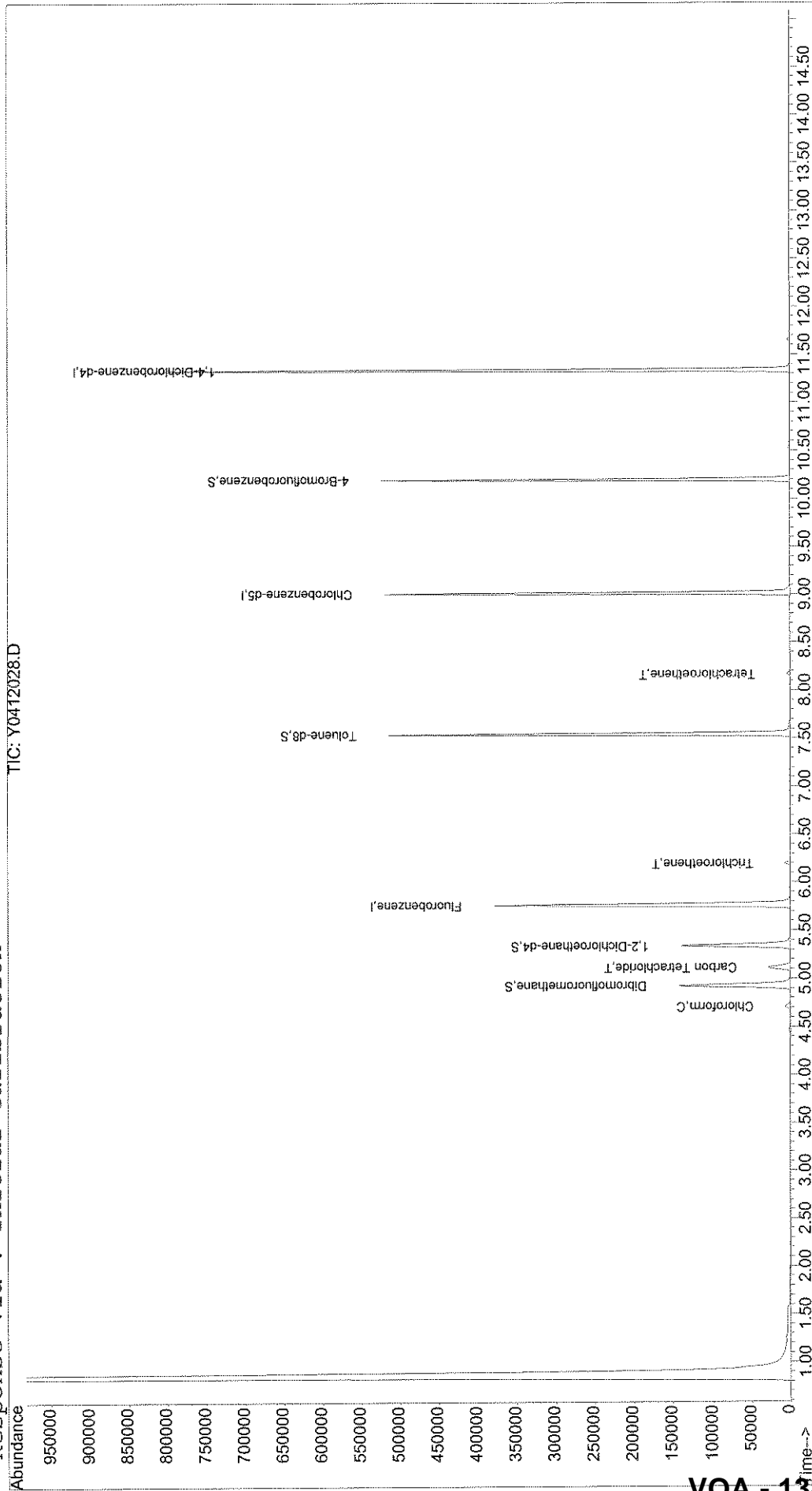
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412028.D Vial: 36  
Acq On : 12 Apr 2007 17:08 Operator: LPM  
Sample : JPL29-015 Inst : Yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:07 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration





Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412028.D  
 Acq On : 12 Apr 2007 17:08  
 Sample : JPL29-015  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:07 2007

Vial: 36  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	348438	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	132996	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	182902	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	103762	49.22	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	98.44%
37) 1,2-Dichloroethane-d4	5.34	65	108564	48.54	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	97.08%
51) Toluene-d8	7.54	98	303162	57.44	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	114.88%
72) 4-Bromofluorobenzene	10.20	95	135538	51.29	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.23	76	651	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0412028.D 8260B.M Fri Apr 13 10:09:12 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412028.D  
 Acq On : 12 Apr 2007 17:08  
 Sample : JPL29-015  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:07 2007

Vial: 36  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	6995	1.56	ug/l ✓	99
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	5.11	117	22230	7.09	ug/l ✓	100
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	2590	1.06	ug/l ✓	96
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	7.61	92	217		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	1006	0.55	ug/l ✓	95
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	372		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	218		N.D.	
65) m,p-Xylene	9.29	106	518		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*Handwritten signature and date: JPL 4/13/07*

(#) = qualifier out of range (m) = manual integration  
 Y0412028.D 8260B.M Fri Apr 13 10:09:12 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412028.D  
 Acq On : 12 Apr 2007 17:08  
 Sample : JPL29-015  
 Misc : 5mL+IS/SS #3

Vial: 36  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

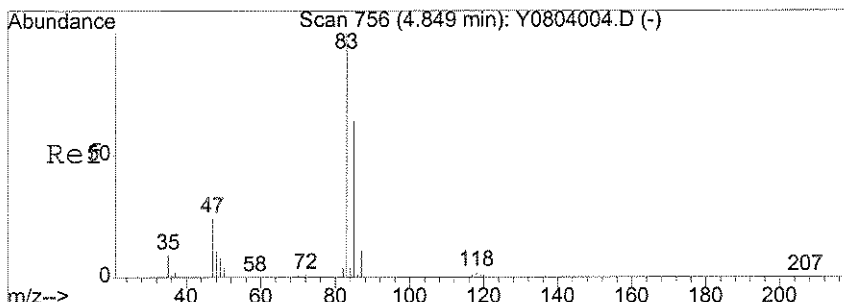
MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:07 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

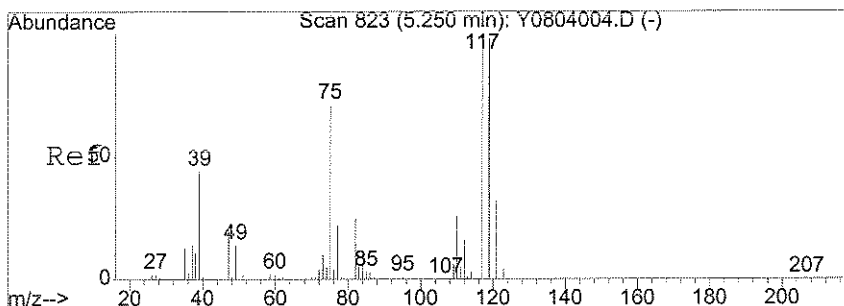
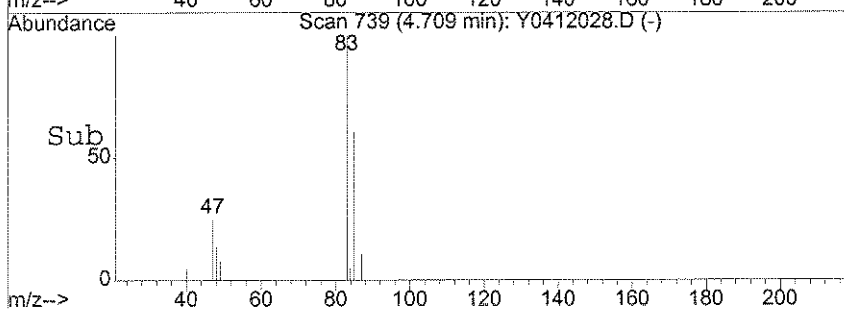
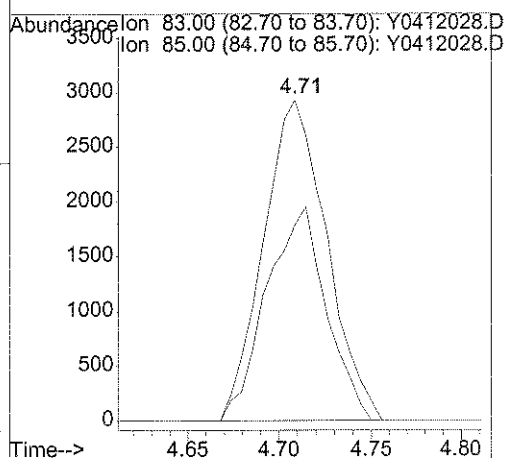
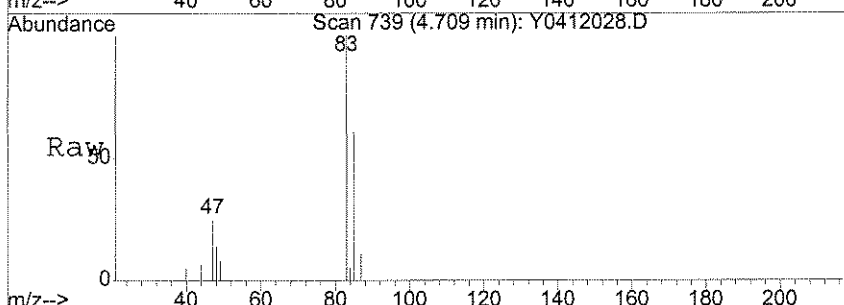
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	132		N.D.	
78) 4-Chlorotoluene	10.47	91	132		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	54		N.D.	
82) sec-butylbenzene	11.18	105	54		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	59		N.D.	
84) 4-Isopropyltoluene	11.34	119	265		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	59		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	280		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412028.D 8260B.M Fri Apr 13 10:09:13 2007



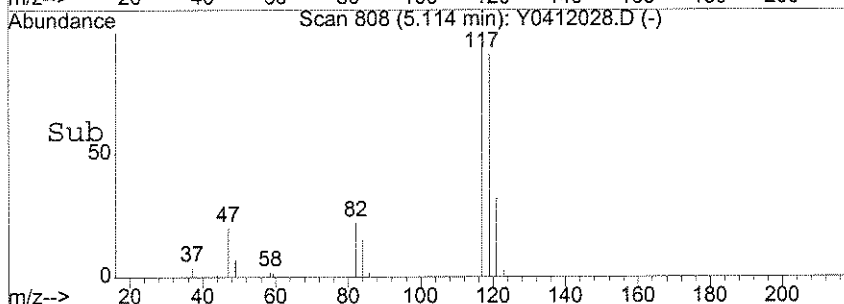
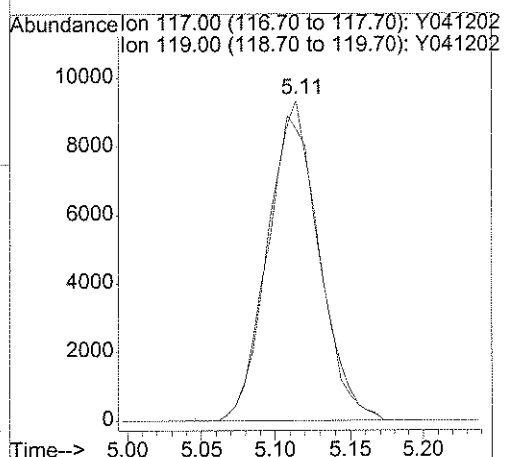
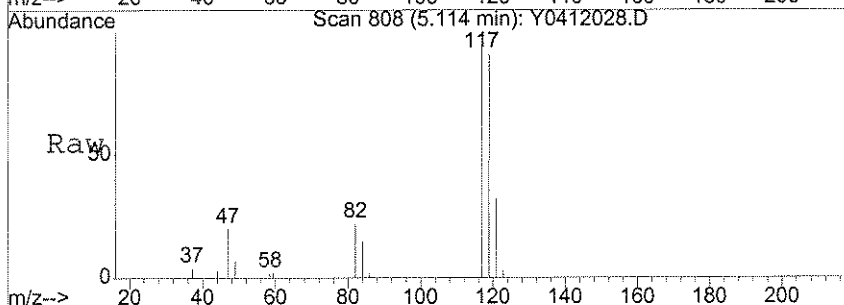
#31  
 Chloroform  
 Concen: 1.56 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0412028.D  
 Acq: 12 Apr 2007 17:08

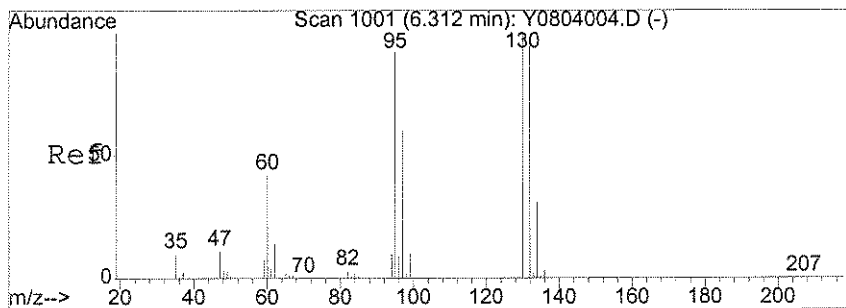
Tgt Ion	Resp	Lower	Upper
83	6995		
83	100		
85	62.7	43.3	83.3



#35  
 Carbon Tetrachloride  
 Concen: 7.09 ug/l  
 RT: 5.11 min Scan# 808  
 Delta R.T. 0.01 min  
 Lab File: Y0412028.D  
 Acq: 12 Apr 2007 17:08

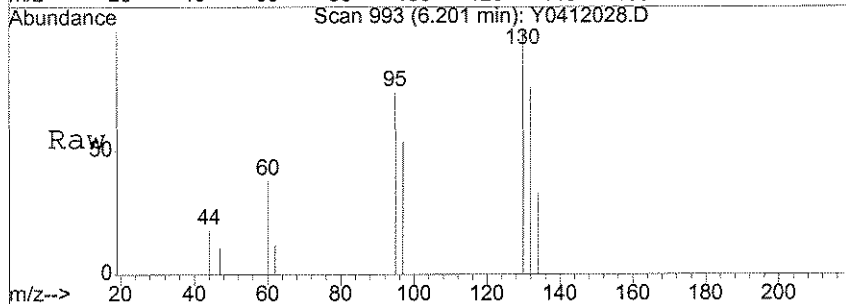
Tgt Ion	Resp	Lower	Upper
117	22230		
117	100		
119	98.0	78.2	118.2



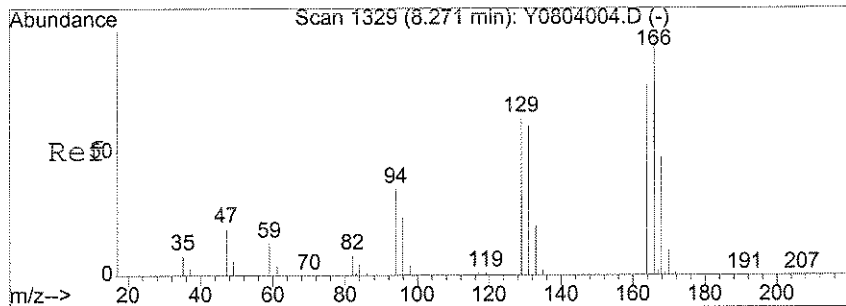
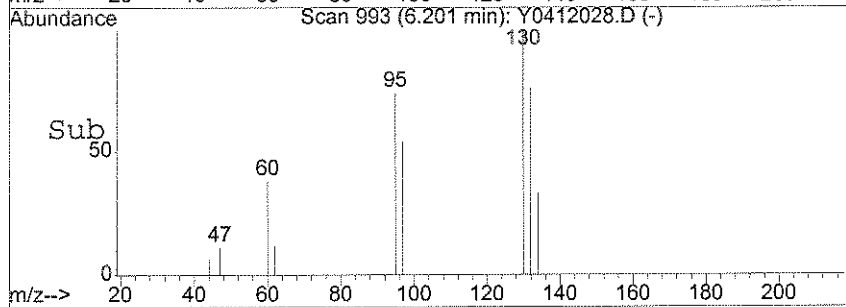
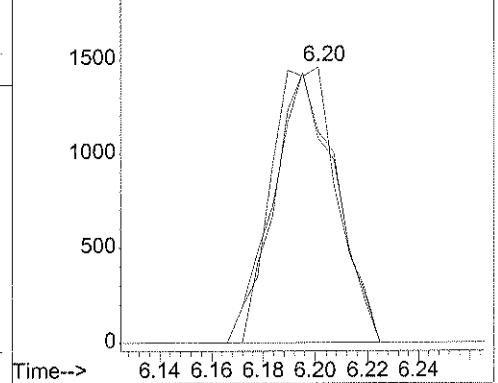


#41  
 Trichloroethene  
 Concen: 1.06 ug/l  
 RT: 6.20 min Scan# 993  
 Delta R.T. 0.01 min  
 Lab File: Y0412028.D  
 Acq: 12 Apr 2007 17:08

Tgt Ion	Resp	Lower	Upper
130	2590		
132	90.1	75.0	115.0
95	91.9	69.4	109.4

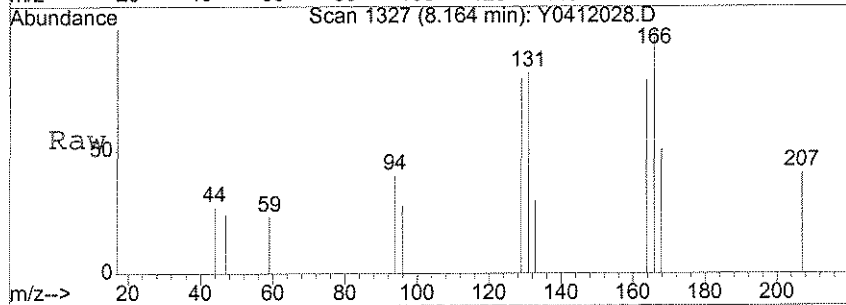


Abundance Ion 130.00 (129.70 to 130.70): Y041202  
 Ion 132.00 (131.70 to 132.70): Y041202  
 Ion 95.00 (94.70 to 95.70): Y0412028.D

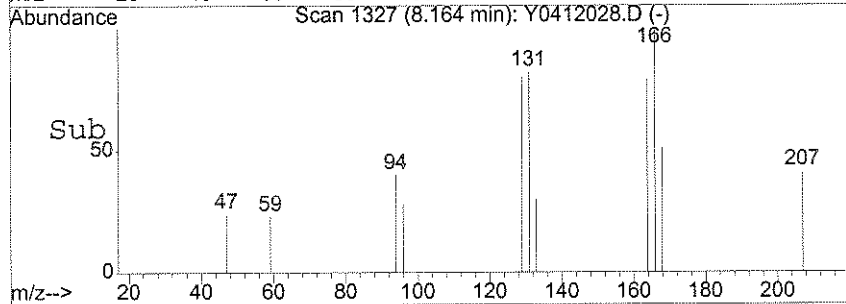
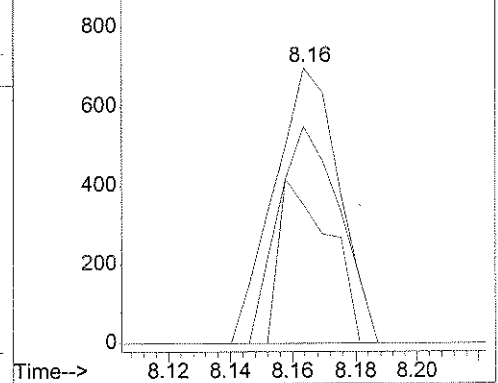


#56  
 Tetrachloroethene  
 Concen: 0.55 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. -0.00 min  
 Lab File: Y0412028.D  
 Acq: 12 Apr 2007 17:08

Tgt Ion	Resp	Lower	Upper
166	1006		
164	75.1	63.3	94.9
168	46.1	39.6	59.4



Abundance Ion 165.95 (165.65 to 166.65): Y041202  
 Ion 163.95 (163.65 to 164.65): Y041202  
 Ion 167.95 (167.65 to 168.65): Y041202



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-016  
 Lab File ID: Y0412029.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:31  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	1.1	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	6.5	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.59	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-016  
 Lab File ID: Y0412029.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:31  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-016  
 Lab File ID: Y0412029.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:31  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

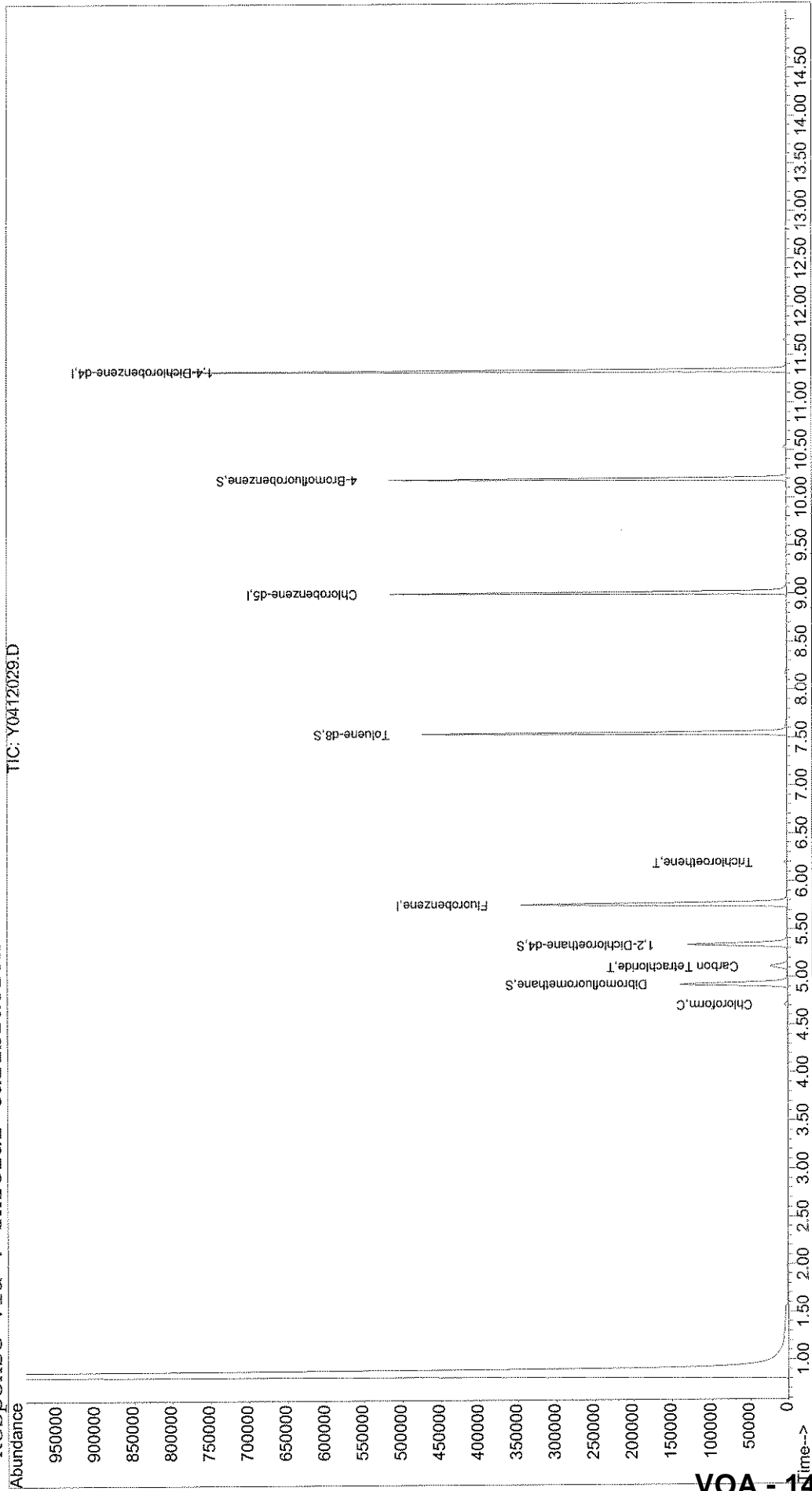


Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412029.D  
Acq On : 12 Apr 2007 17:31  
Sample : JPL29-016  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:10 2007

Vial: 37  
Operator: LPM  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412029.D  
 Acq On : 12 Apr 2007 17:31  
 Sample : JPL29-016  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:10 2007

Vial: 37  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	316113	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	129623	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	184588	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	102196	53.43	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	106.86%
37) 1,2-Dichloroethane-d4	5.34	65	101730	50.14	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.28%
51) Toluene-d8	7.54	98	279503	54.34	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	108.68%
72) 4-Bromofluorobenzene	10.20	95	134974	50.61	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.09	101	143	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	417	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412029.D  
 Acq On : 12 Apr 2007 17:31  
 Sample : JPL29-016  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:10 2007

Vial: 37  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	4.71	83	4478	1.10	ug/l ✓	98
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	5.11	117	18488	6.50	ug/l ✓	99
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	0.00	78	0	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.		
41) Trichloroethene	6.20	130	1319	0.59	ug/l ✓	97
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	0.00	41	0	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	0.00	92	0	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.		
56) Tetrachloroethene	8.17	166	430	N.D.		
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.		
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	0.00	112	0	N.D.		
62) 1-Chlorohexane	9.01	91	366	N.D.		
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.17	91	67	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.		
66) o-xylene	0.00	106	0	N.D.		
67) Styrene	0.00	104	0	N.D.		
68) Bromoform	0.00	173	0	N.D.		

*Handwritten signature and date: JNH 4/13/07*

(#) = qualifier out of range (m) = manual integration  
 Y0412029.D 8260B.M Fri Apr 13 10:09:25 2007

Quantitation Report

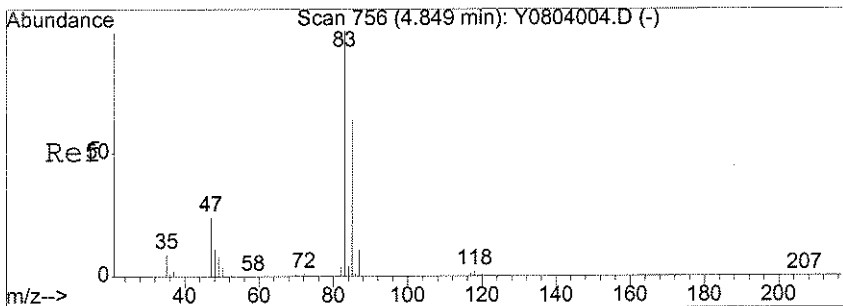
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 Acq On : 12 Apr 2007 17:31  
 Sample : JPL29-016  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:10 2007

Vial: 37  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

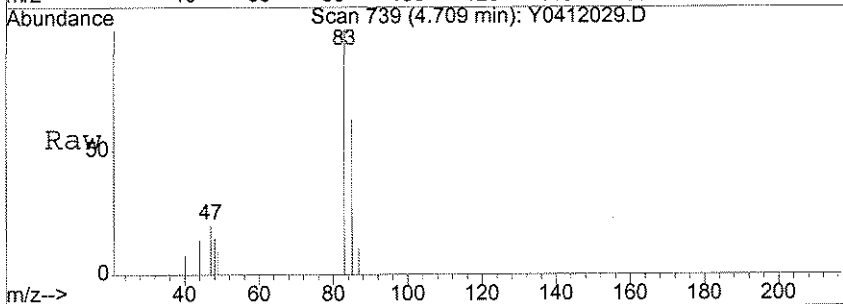
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.18	105	85		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	60		N.D.	
78) 4-Chlorotoluene	10.47	91	60		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	167		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	254		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

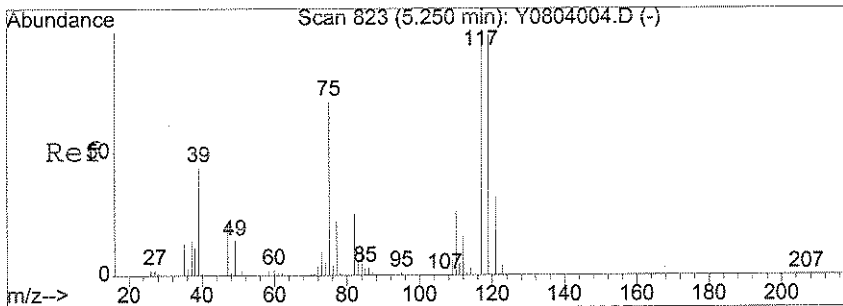
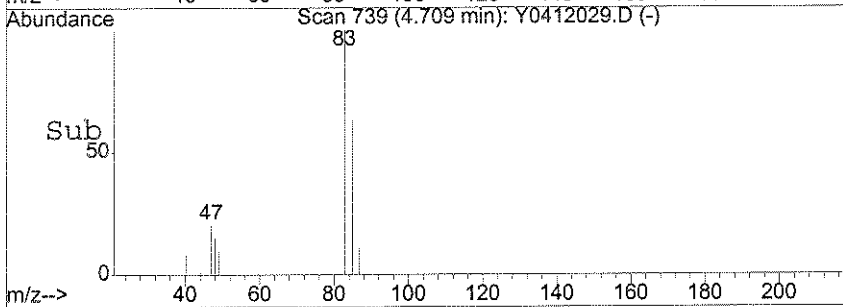
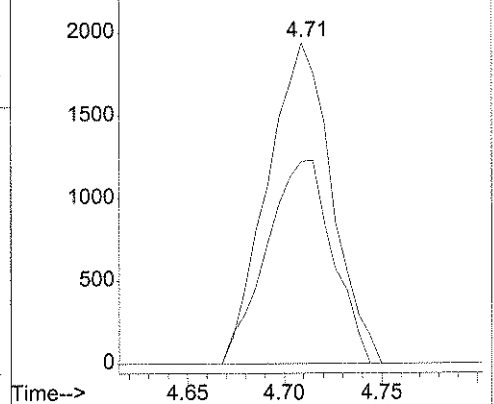


#31  
 Chloroform  
 Concen: 1.10 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0412029.D  
 Acq: 12 Apr 2007 17:31

Tgt Ion:	83	Resp:	4478
Ion Ratio	Lower	Upper	
83	100		
85	65.1	43.3	83.3

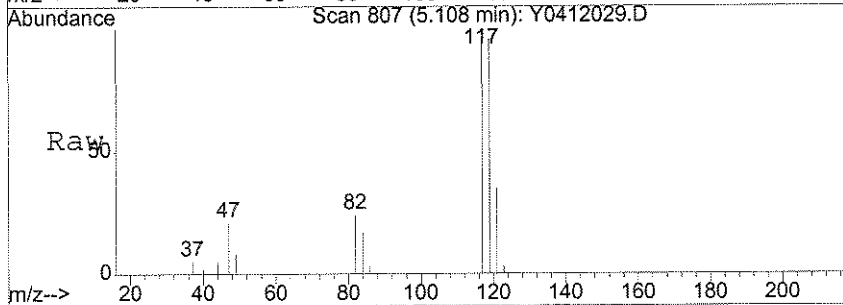


Abundance Ion 83.00 (82.70 to 83.70): Y0412029.D  
 Ion 85.00 (84.70 to 85.70): Y0412029.D

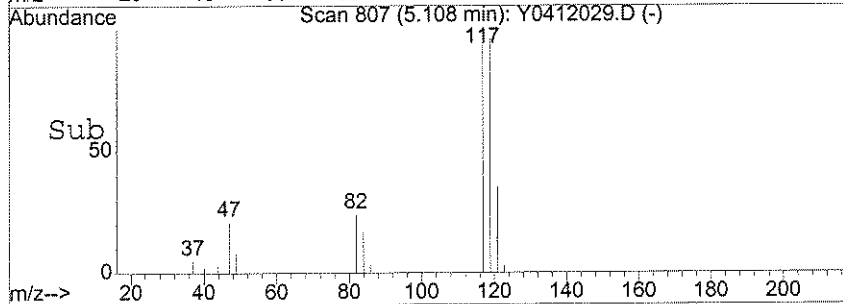
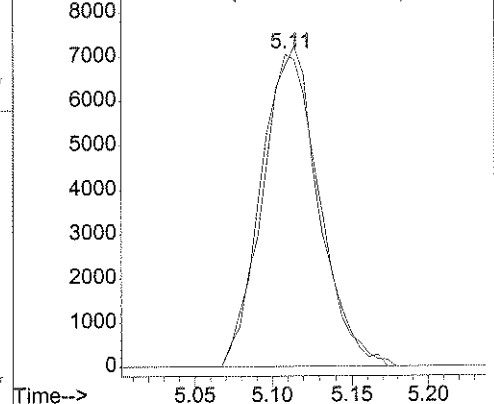


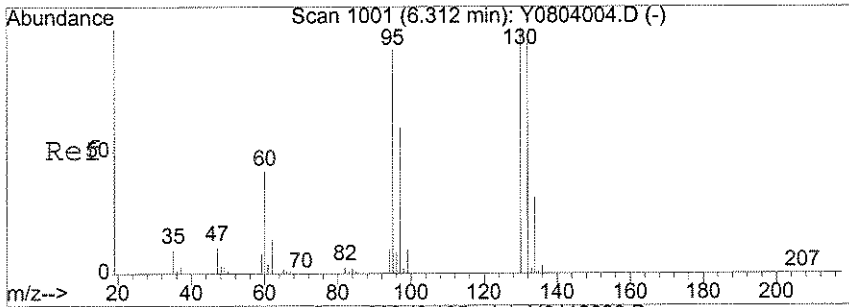
#35  
 Carbon Tetrachloride  
 Concen: 6.50 ug/l  
 RT: 5.11 min Scan# 807  
 Delta R.T. -0.00 min  
 Lab File: Y0412029.D  
 Acq: 12 Apr 2007 17:31

Tgt Ion:	117	Resp:	18488
Ion Ratio	Lower	Upper	
117	100		
119	96.8	78.2	118.2



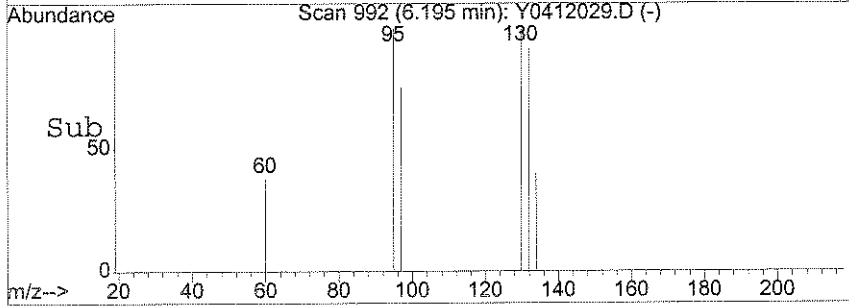
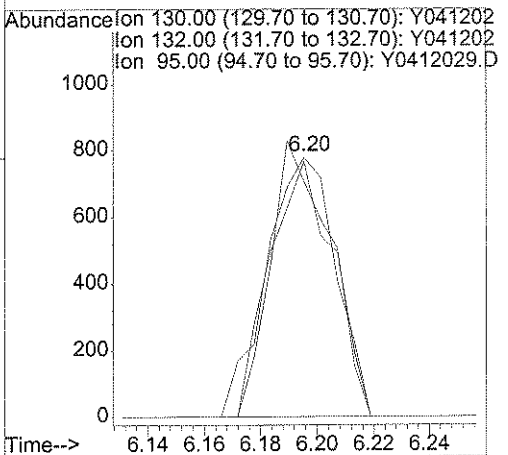
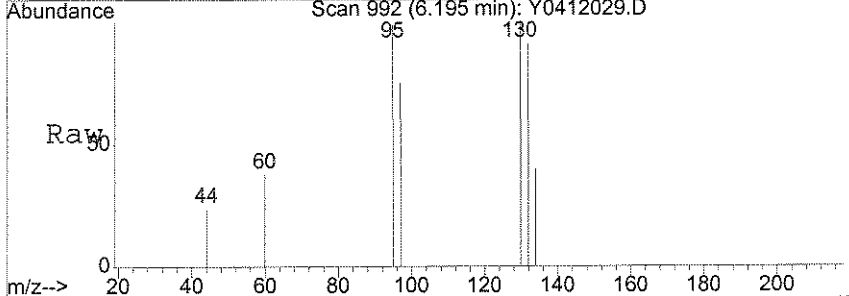
Abundance Ion 117.00 (116.70 to 117.70): Y0412029.D  
 Ion 119.00 (118.70 to 119.70): Y0412029.D





#41  
 Trichloroethene  
 Concen: 0.59 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0412029.D  
 Acq: 12 Apr 2007 17:31

Tgt Ion	Resp	Lower	Upper
130	1319		
130	100		
132	91.7	75.0	115.0
95	91.1	69.4	109.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-017  
 Lab File ID: Y0412030.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:56  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-017  
 Lab File ID: Y0412030.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 17:56  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-18-2

Lab Name: Laucks Testing Laboratories, Inc.

SDG No.: JPL29

Matrix: (SOIL/SED/WATER) Water

Sample wt/vol: 5.00 (g/mL) mL

Level: (LOW/MED) \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

GC Column: DB-624 20m ID: 0.18 (mm)

Soil Extract Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin

Run Sequence: R016785

Lab Sample ID: JPL29-017

Lab File ID: Y0412030.D

Date Collected: 03/29/2007

Date/Time Analyzed: 04/12/2007 17:56

Dilution Factor: 1.0

Soil Aliquot Volume: \_\_\_\_\_ (uL)

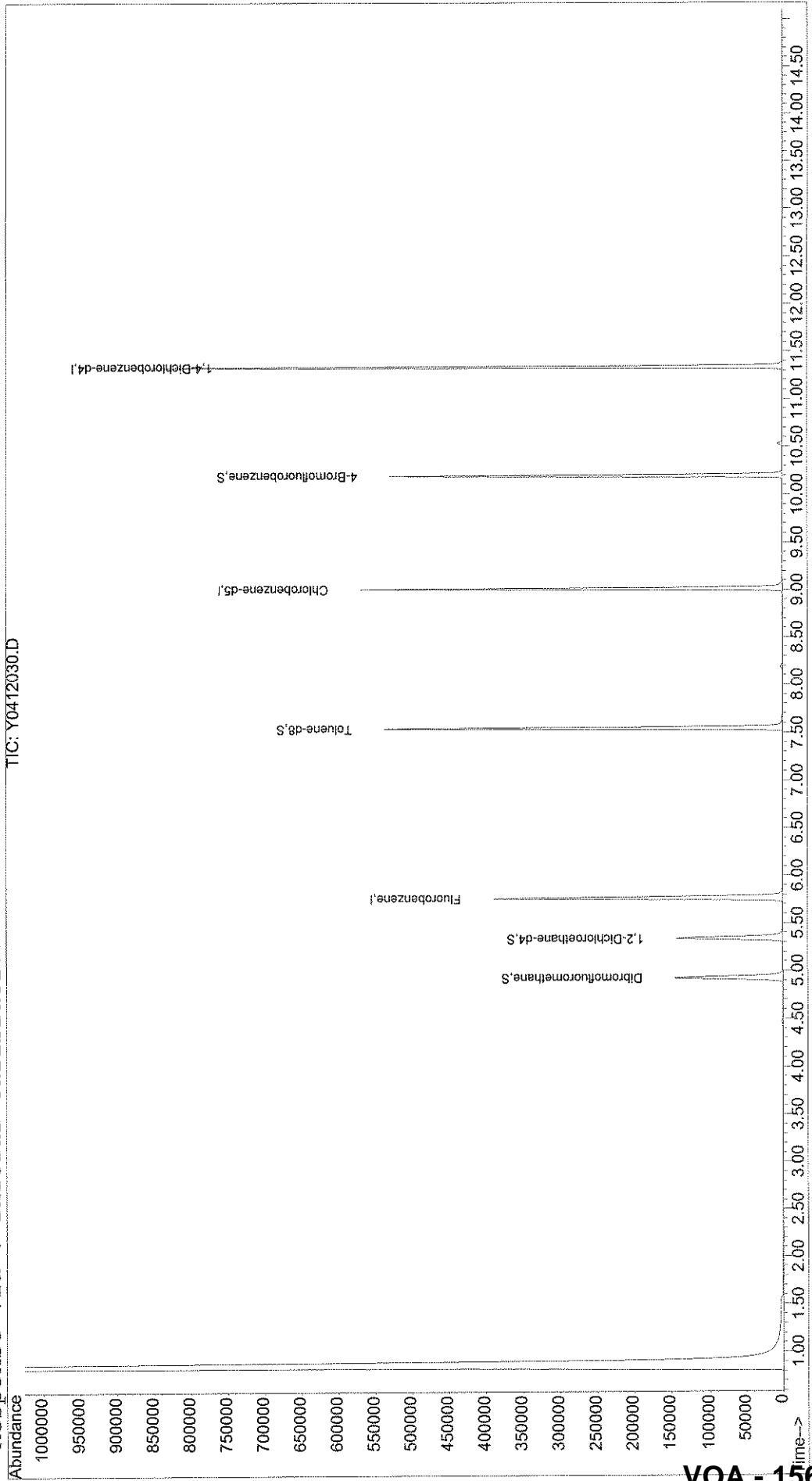
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412030.D Vial: 38  
Acq On : 12 Apr 2007 17:56 Operator: LPM  
Sample : JPL29-017 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:11 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412030.D  
 Acq On : 12 Apr 2007 17:56  
 Sample : JPL29-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:11 2007

Vial: 38  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	360155	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	143091	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	187046	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	106776	49.00	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	98.00%
37) 1,2-Dichloroethane-d4	5.34	65	113999	49.31	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.62%
51) Toluene-d8	7.55	98	320384	56.42	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.84%
72) 4-Bromofluorobenzene	10.20	95	142121	52.59	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	0.00	76	0	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412030.D 8260B.M Fri Apr 13 10:09:36 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412030.D  
 Acq On : 12 Apr 2007 17:56  
 Sample : JPL29-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:11 2007

Vial: 38  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.02	91	447		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	347		N.D.	
65) m,p-Xylene	9.30	106	144		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412030.D 8260B.M Fri Apr 13 10:09:37 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412030.D  
 Acq On : 12 Apr 2007 17:56  
 Sample : JPL29-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:11 2007

Vial: 38  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	113		N.D.	
78) 4-Chlorotoluene	10.46	91	113		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	57		N.D.	
82) sec-butylbenzene	11.18	105	54		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.33	119	260		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	190		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	13.54	128	56		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-2-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-018  
 Lab File ID: Y0412031.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 18:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-2-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-018  
 Lab File ID: Y0412031.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 18:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-2-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL29  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-018  
 Lab File ID: Y0412031.D  
 Date Collected: 03/29/2007  
 Date/Time Analyzed: 04/12/2007 18:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

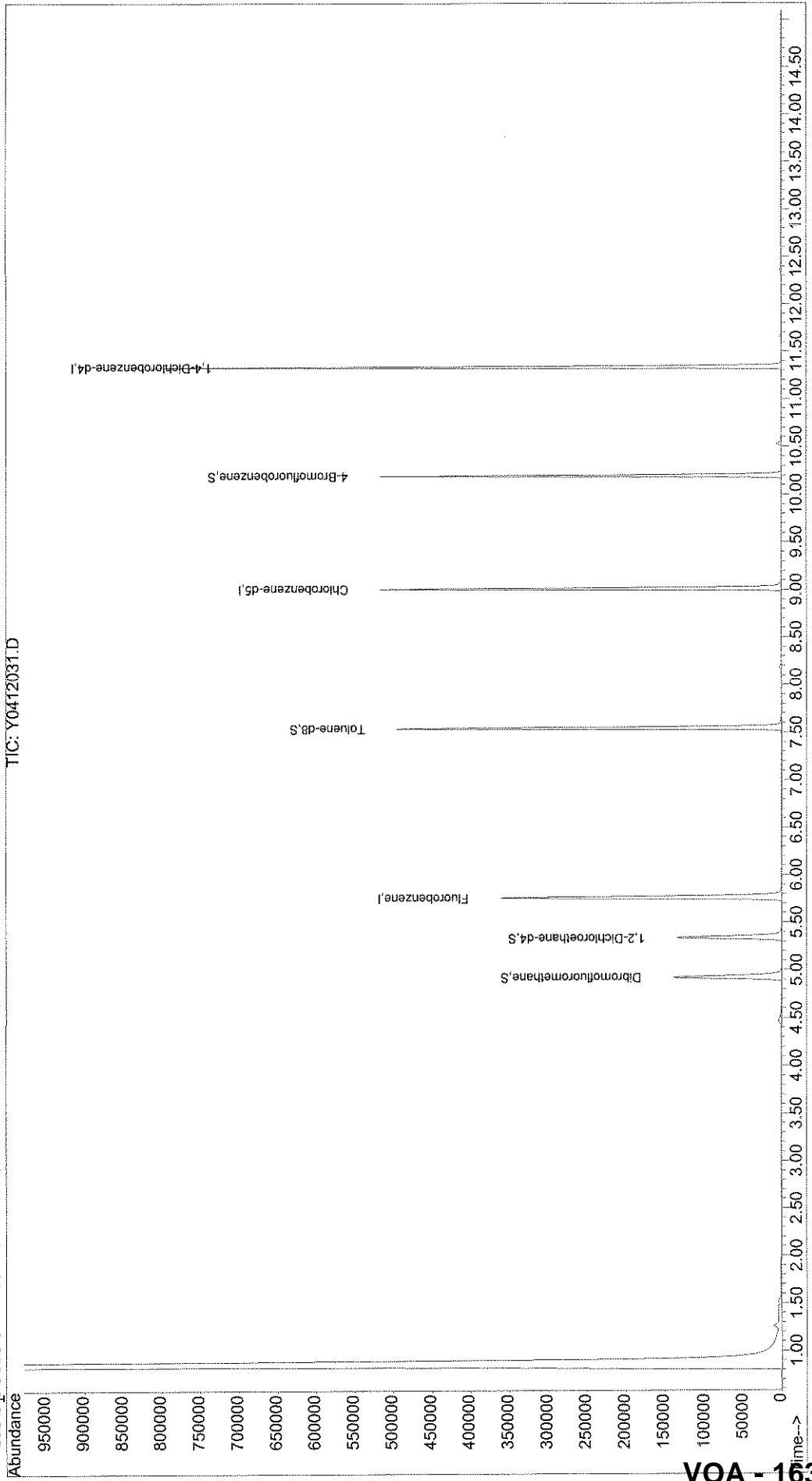
Comments:



Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412031.D Vial: 39  
Acq On : 12 Apr 2007 18:20 Operator: LPM  
Sample : JPL29-018 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 13 8:12 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 163

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412031.D  
 Acq On : 12 Apr 2007 18:20  
 Sample : JPL29-018  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:12 2007

Vial: 39  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	342211	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	133319	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	175947	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	103979	50.22	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.44%
37) 1,2-Dichloroethane-d4	5.34	65	108934	49.59	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.18%
51) Toluene-d8	7.54	98	296965	56.13	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.26%
72) 4-Bromofluorobenzene	10.20	95	133830	52.65	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412031.D 8260B.M Fri Apr 13 10:09:54 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412031.D  
 Acq On : 12 Apr 2007 18:20  
 Sample : JPL29-018  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:12 2007

Vial: 39  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	391		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	421		N.D.	
65) m,p-Xylene	9.29	106	54		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0412031.D 8260B.M Fri Apr 13 10:09:55 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041207\Y0412031.D  
 Acq On : 12 Apr 2007 18:20  
 Sample : JPL29-018  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 13 8:12 2007

Vial: 39  
 Operator: LPM  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	68		N.D.	
78) 4-Chlorotoluene	10.47	91	68		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	325		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	193		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

**TICS SUMMARY**

SDG JPL29

VOLATILES ANALYSIS

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041107MVOWM3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: B041107MVOWM3  
 Lab File ID: M0411009.D  
 Date Collected: \_\_\_\_\_  
 Date Analyzed: 04/11/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411009.D                   Vial: 56  
Acq On    : 11 Apr 2007 10:13                   Operator: DGA  
Sample    : B041107MVOWM3                   Inst    : MOBY  
Misc      : 5ml PFW+IS/SS(MV8-38-11)       Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411009.D 524.M Thu Apr 12 09:10:37 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041207MVOWY1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016785

Matrix: (SOIL/WATER) Water

Lab Sample ID: B041207MVOWY1

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0412014.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/12/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:



Library Search Compound Report

Data File : X:\MSVOA\YODA\041207\Y0412014.D Vial: 22  
Acq On : 12 Apr 2007 11:15 Operator: LPM  
Sample : B041207MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-1) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0412014.D 8260B.M Fri Apr 13 10:33:17 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041307MVOWY1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: B041307MVOWY1

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413007.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413007.D Vial: 22  
Acq On : 13 Apr 2007 8:44 Operator: LNH  
Sample : B041307MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-1) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413007.D 8260B.M Fri Apr 13 14:06:21 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-14-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-001  
 Lab File ID: M0411016.D  
 Date Collected: 03/29/2007  
 Date Analyzed: 04/11/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411016.D                   Vial: 56  
Acq On    : 11 Apr 2007 13:25                   Operator: DGA  
Sample    : JPL29-001                         Inst     : MOBY  
Misc      : #3 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411016.D 524.M Thu Apr 12 09:12:09 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-14-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-002  
 Lab File ID: M0411017.D  
 Date Collected: 03/29/2007  
 Date Analyzed: 04/11/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411017.D Vial: 57  
Acq On : 11 Apr 2007 13:47 Operator: DGA  
Sample : JPL29-002 Inst : MOBY  
Misc : #2 5ml +IS/SS (524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411017.D 524.M Thu Apr 12 09:13:36 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-14-3

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/29/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/11/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411018.D                   Vial: 58  
Acq On    : 11 Apr 2007  14:10                   Operator: DGA  
Sample    : JPL29-003                           Inst     : MOBY  
Misc      : #4 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411018.D  524.M   Thu Apr 12 09:14:54 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-14-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/29/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/11/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411019.D                   Vial: 59  
Acq On    : 11 Apr 2007 14:33                   Operator: DGA  
Sample    : JPL29-004                         Inst     : MOBY  
Misc      : #3 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411019.D 524.M Thu Apr 12 09:16:22 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-14-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-005  
 Lab File ID: M0411020.D  
 Date Collected: 03/29/2007  
 Date Analyzed: 04/11/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411020.D Vial: 60  
Acq On : 11 Apr 2007 14:56 Operator: DGA  
Sample : JPL29-005 Inst : MOBY  
Misc : #3 5ml +IS/SS (524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title : VOA 524- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411020.D 524.M Thu Apr 12 09:17:35 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/29/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/11/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411021.D                   Vial: 61  
Acq On    : 11 Apr 2007 15:20                   Operator: DGA  
Sample    : JPL29-006                         Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411021.D 524.M Thu Apr 12 09:19:11 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-2-3/28/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/29/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/11/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
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Comments:



Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411022.D                    Vial: 62  
Acq On    : 11 Apr 2007 15:43                                    Operator: DGA  
Sample    : JPL29-007    Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)                                Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411022.D 524.M Thu Apr 12 09:20:11 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-17-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016814  
 Lab Sample ID: JPL29-008  
 Lab File ID: M0411023.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/11/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411023.D                   Vial: 63  
Acq On    : 11 Apr 2007 16:05                   Operator: DGA  
Sample    : JPL29-008                         Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library     : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411023.D 524.M Thu Apr 12 09:27:27 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-17-3

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016814

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0411024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/11/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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04					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041107A\M0411024.D                   Vial: 64  
Acq On    : 11 Apr 2007 16:28                   Operator: DGA  
Sample    : JPL29-009                           Inst     : MOBY  
Misc      : #2 5ml +IS/SS (524)               Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\524.M (RTE Integrator)  
Title      : VOA 524- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0411024.D 524.M    Thu Apr 12 09:28:32 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-17-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413013.D Vial: 28  
Acq On : 13 Apr 2007 11:12 Operator: LNH  
Sample : JPL29-010 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413013.D 8260B.M Fri Apr 13 14:19:33 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-1-1Q07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-011  
 Lab File ID: Y0412024.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/12/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:



Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 15:29  
Data File: X:\MSVOA\YODA\041207\Y0412024.D  
Name: JPL29-011  
Misc: 5mL+IS/SS #1  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Y0412024.D 8260B.M			Fri Apr 13 10:08:27 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016785

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0412025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/12/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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04				
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 15:53  
Data File: X:\MSVOA\YODA\041207\Y0412025.D  
Name: JPL29-012  
Misc: 5mL+IS/SS #2  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Y0412025.D 8260B.M			Fri Apr 13 10:08:40 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-3-3/29/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-013  
 Lab File ID: Y0412026.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/12/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 16:18  
Data File: X:\MSVOA\YODA\041207\Y0412026.D  
Name: JPL29-013  
Misc: 5mL+IS/SS #1  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
-----								
Y0412026.D 8260B.M			Fri Apr 13 10:08:53 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-18-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL)      mL  
 Level: (LOW/MED)       
 % Moisture: not dec.       
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume:      (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-014  
 Lab File ID: Y0412027.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/12/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume:      (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 16:43  
Data File: X:\MSVOA\YODA\041207\Y0412027.D  
Name: JPL29-014  
Misc: 5mL+IS/SS #3  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Y0412027.D 8260B.M			Fri Apr 13 10:09:04 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-18-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-015  
 Lab File ID: Y0412028.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/12/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:



Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 17:08  
Data File: X:\MSVOA\YODA\041207\Y0412028.D  
Name: JPL29-015  
Misc: 5mL+IS/SS #3  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Y0412028.D 8260B.M			Fri Apr 13 10:09:18 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-18-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL29  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL)      mL  
 Level: (LOW/MED)       
 % Moisture: not dec.       
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume:      (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016785  
 Lab Sample ID: JPL29-016  
 Lab File ID: Y0412029.D  
 Date Collected: 03/30/2007  
 Date Analyzed: 04/12/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume:      (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 17:31  
Data File: X:\MSVOA\YODA\041207\Y0412029.D  
Name: JPL29-016  
Misc: 5mL+IS/SS #3  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
Y0412029.D 8260B.M			Fri Apr 13 10:09:30 2007					

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-18-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016785

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0412030.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/12/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 17:56  
Data File: X:\MSVOA\YODA\041207\Y0412030.D  
Name: JPL29-017  
Misc: 5mL+IS/SS #2  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC	Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
-----									
Y0412030.D	8260B.M			Fri Apr 13 10:09:42 2007					

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-2-1Q07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL29

Run Sequence: R016785

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL29-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0412031.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/12/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Tentatively Identified Compound (LSC) summary

Operator ID: LPM      Date Acquired: 12 Apr 2007 18:20  
Data File: X:\MSVOA\YODA\041207\Y0412031.D  
Name: JPL29-018  
Misc: 5mL+IS/SS #1  
Method: X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title: VOA 8260- 5ML Calibration 5973Y  
Library Searched: D:\DATABASE\NIST129K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc
-----								
Y0412031.D 8260B.M			Fri Apr 13 10:10:00 2007					

**Metals Data**

**JPL29**



COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories  
 Lab Code: LAUCKS  
 SOW No.: \_\_\_\_\_

Contract: JPL Groundwater Monitorin  
 SDG No.: JPL29

Sample No.	Lab Sample ID
MW-14-3	JPL29-003
MW-14-3MS	JPL29-003MS
MW-14-3MSD	JPL29-003MSD
MW-14-2	JPL29-004
MW-14-1	JPL29-005
EB-2-3/28/07	JPL29-006
MW-17-4	JPL29-008
MW-17-3	JPL29-009
MW-17-2	JPL29-010
DUPE-1-1Q07	JPL29-011
DUPE-1-1Q07MS	JPL29-011MS
DUPE-1-1Q07MSD	JPL29-011MSD
EB-3-3/29/07	JPL29-012
MW-18-4	JPL29-015
MW-18-3	JPL29-016
MW-18-2	JPL29-017
DUPE-2-1Q07	JPL29-018

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

---

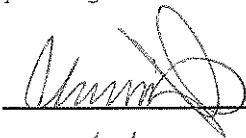


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I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature:  Name: Cheronna Oreiro

Date: 01/10/2007 Title: metals Lead

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-14-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-003

Level (low/med): LOW

Date Received: 03/29/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-14-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-004

Level (low/med): LOW

Date Received: 03/29/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.83			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-14-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-005

Level (low/med): LOW

Date Received: 03/29/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-2-3/28/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-006

Level (low/med): LOW

Date Received: 03/29/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.53			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-17-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-008

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-17-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-009

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.34			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-17-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-010

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.69			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-1-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-011

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.79			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

SW-846

-1-

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-3-3/29/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-012

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.41			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-18-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-015

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.73			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-18-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-016

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	4.90			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-18-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-017

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

SW-846

-1-

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-2-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL29

Matrix (soil/water): Water

Lab Sample ID: JPL29-018

Level (low/med): LOW

Date Received: 03/30/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL29**



**COVER PAGE - INORGANIC ANALYSES DATA PACKAGE**

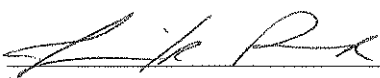
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL29

Client Identification	Lab Sample Work Order Number
MW-14-5	JPL29-001DL
MW-14-4	JPL29-002DL
MW-14-3	JPL29-003DL
MW-14-3MS	JPL29-003MS
MW-14-3MSD	JPL29-003MSD
MW-14-2	JPL29-004DL
MW-14-1	JPL29-005DL
EB-2-3/28/07	JPL29-006
MW-17-4	JPL29-008
MW-17-3	JPL29-009DL
MW-17-2	JPL29-010DL
DUPE-1-1Q07	JPL29-011DL
DUPE-1-1Q07MS	JPL29-011MS
DUPE-1-1Q07MSD	JPL29-011SDM
EB-3-3/29/07	JPL29-012
MW-18-5	JPL29-014
MW-18-4	JPL29-015DL
MW-18-3	JPL29-016DL
MW-18-2	JPL29-017DL
DUPE-2-1Q07	JPL29-018DL

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Pennor

Date: 4-16-07

Title: Inorganics Lead

## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-14-5 **Date/Time Collected:** 03/28/2007 08:13  
**Lab Sample ID:** JPL29-001 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/11/2007	04/12/2007	R016614

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-14-4 **Date/Time Collected:** 03/28/2007 08:40  
**Lab Sample ID:** JPL29-002 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	2.5	U	2.5	0.35	04/11/2007	04/12/2007	R016614

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-14-3 **Date/Time Collected:** 03/28/2007 09:10  
**Lab Sample ID:** JPL29-003 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/11/2007	04/12/2007	R016614

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-14-2 **Date/Time Collected:** 03/28/2007 10:03  
**Lab Sample ID:** JPL29-004 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-14-1 **Date/Time Collected:** 03/28/2007 10:34  
**Lab Sample ID:** JPL29-005 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** EB-2-3/28/07 **Date/Time Collected:** 03/28/2007 10:22  
**Lab Sample ID:** JPL29-006 **Date/Time Received:** 03/29/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/09/2007	04/10/2007	R016670



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-17-4 **Date/Time Collected:** 03/29/2007 07:49  
**Lab Sample ID:** JPL29-008 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-17-3 **Date/Time Collected:** 03/29/2007 08:21  
**Lab Sample ID:** JPL29-009 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	47		2.0	0.28	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-17-2 **Date/Time Collected:** 03/29/2007 08:59  
**Lab Sample ID:** JPL29-010 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	10		4.0	0.56	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** DUPE-1-1Q07 **Date/Time Collected:** 03/29/2007 00:00  
**Lab Sample ID:** JPL29-011 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	9.4		4.0	0.56	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** EB-3-3/29/07 **Date/Time Collected:** 03/29/2007 08:44  
**Lab Sample ID:** JPL29-012 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-18-5 **Date/Time Collected:** 03/29/2007 10:54  
**Lab Sample ID:** JPL29-014 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-18-4 **Date/Time Collected:** 03/29/2007 11:20  
**Lab Sample ID:** JPL29-015 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-18-3 **Date/Time Collected:** 03/29/2007 11:55  
**Lab Sample ID:** JPL29-016 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/09/2007	04/10/2007	R016670



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** MW-18-2 **Date/Time Collected:** 03/29/2007 12:24  
**Lab Sample ID:** JPL29-017 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/09/2007	04/10/2007	R016670

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL29  
**Sample Number:** DUPE-2-1Q07 **Date/Time Collected:** 03/29/2007 00:00  
**Lab Sample ID:** JPL29-018 **Date/Time Received:** 03/30/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/09/2007	04/10/2007	R016670

**SAMPLE DATA**

SDG # JPL30

Volatiles Analysis

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-001  
 Lab File ID: Y0413014.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 11:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-001

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413014.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 11:36

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.26	J
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-001  
 Lab File ID: Y0413014.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 11:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

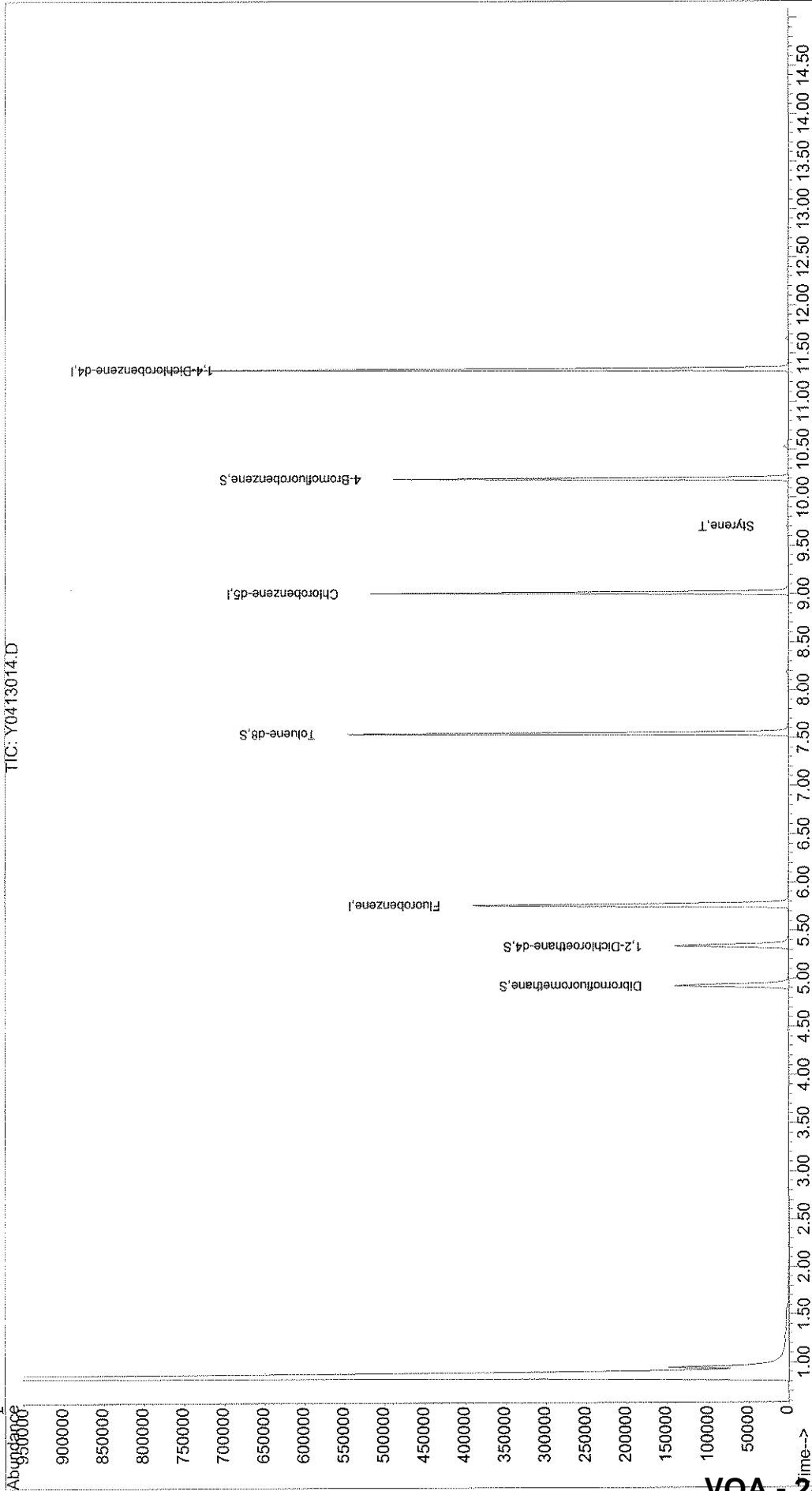
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413014.D Vial: 28  
Acq On : 13 Apr 2007 11:36 Operator: LNH  
Sample : JPL30-001 Inst : Yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:05 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413014.D  
 Acq On : 13 Apr 2007 11:36  
 Sample : JPL30-001  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:05 2007

Vial: 28  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	359058	50.00	ug/l	0.00 105.69%
50) Chlorobenzene-d5	9.01	82	132450	50.00	ug/l	0.00 82.18%
70) 1,4-Dichlorobenzene-d4	11.34	152	174645	50.00	ug/l	0.00 77.50%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101256	46.61	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	93.22%	
37) 1,2-Dichloroethane-d4	5.34	65	111404	48.34	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	96.68%	
51) Toluene-d8	7.55	98	322705	61.39	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	122.78%#	
72) 4-Bromofluorobenzene	10.20	95	127566	50.56	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.21	76	885	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	d
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413014.D  
 Acq On : 13 Apr 2007 11:36  
 Sample : JPL30-001  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:05 2007

Vial: 28  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.39	78	250		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	310		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.02	91	416		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	658		N.D.	
65) m,p-Xylene	9.29	106	310		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	1253	0.26	ug/l	92
68) Bromoform	0.00	173	0		N.D.	

*QNH 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0413014.D 8260B.M Mon Apr 16 09:05:56 2007

Quantitation Report

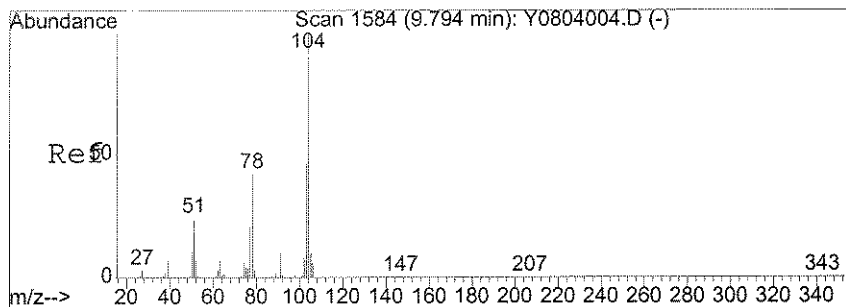
Data File : X:\MSVOA\YODA\041307\Y0413014.D  
 Acq On : 13 Apr 2007 11:36  
 Sample : JPL30-001  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:05 2007

Vial: 28  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

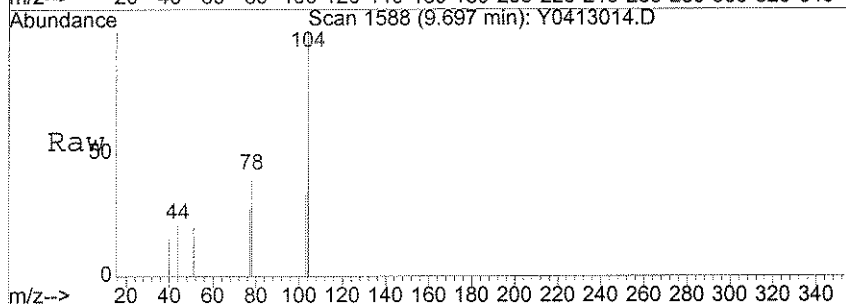
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	106		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	144		N.D.	
78) 4-Chlorotoluene	10.47	91	144		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	131		N.D.	
82) sec-butylbenzene	11.18	105	59		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	55		N.D.	
84) 4-Isopropyltoluene	11.34	119	341		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	55		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	147		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	60		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

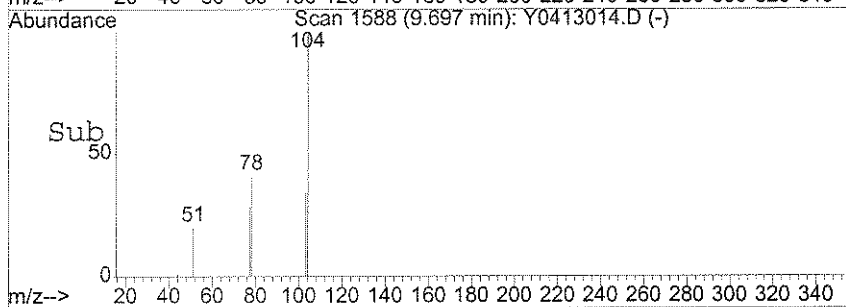
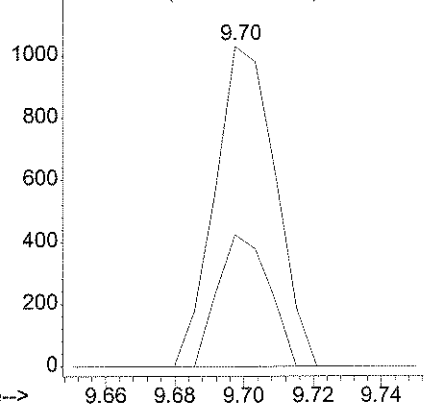


#67  
 Styrene  
 Concen: 0.26 ug/l  
 RT: 9.70 min Scan# 1588  
 Delta R.T. 0.00 min  
 Lab File: Y0413014.D  
 Acq: 13 Apr 2007 11:36

Tgt Ion: 104 Resp: 1253  
 Ion Ratio Lower Upper  
 104 100  
 78 35.0 19.7 59.7



Abundance Ion 104.00 (103.70 to 104.70): Y0413014.D  
 Ion 78.00 (77.70 to 78.70): Y0413014.D



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:01

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:01

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:01

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

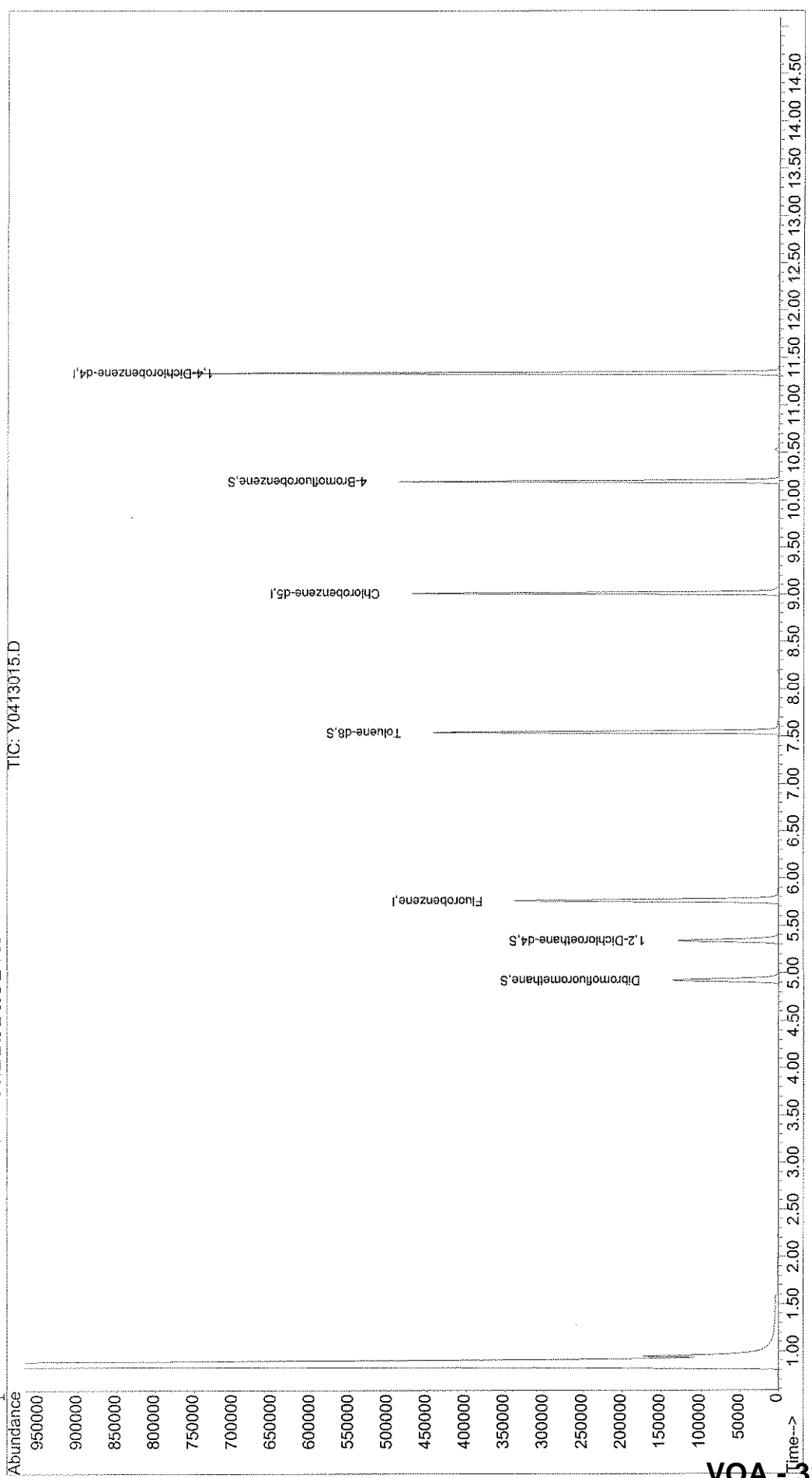
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413015.D Vial: 29  
Acq On : 13 Apr 2007 12:01 Operator: LNH  
Sample : JPL30-002 Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:06 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413015.D  
 Acq On : 13 Apr 2007 12:01  
 Sample : JPL30-002  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:06 2007

Vial: 29  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	307780	50.00	ug/l	0.00 90.59%
50) Chlorobenzene-d5	9.01	82	123268	50.00	ug/l	0.00 76.48%
70) 1,4-Dichlorobenzene-d4	11.34	152	174215	50.00	ug/l	0.00 77.31%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	98152	52.71	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	105.42%
37) 1,2-Dichloroethane-d4	5.34	65	99699	50.47	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.94%
51) Toluene-d8	7.54	98	267261	54.63	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	109.26%
72) 4-Bromofluorobenzene	10.20	95	126699	50.34	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	1010		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	2.83	53	119		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413015.D  
 Acq On : 13 Apr 2007 12:01  
 Sample : JPL30-002  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:06 2007

Vial: 29  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	209		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.02	91	408		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	219		N.D.	
65) m,p-Xylene	9.30	106	195		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	216		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413015.D 8260B.M Mon Apr 16 09:07:01 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413015.D  
 Acq On : 13 Apr 2007 12:01  
 Sample : JPL30-002  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:06 2007

Vial: 29  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	60		N.D.	
78) 4-Chlorotoluene	10.46	91	60		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	71		N.D.	
82) sec-butylbenzene	11.01	105	71		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.33	119	251		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	313		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-003  
 Lab File ID: Y0413016.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 12:26  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:26

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.43	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-003  
 Lab File ID: Y0413016.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 12:26  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

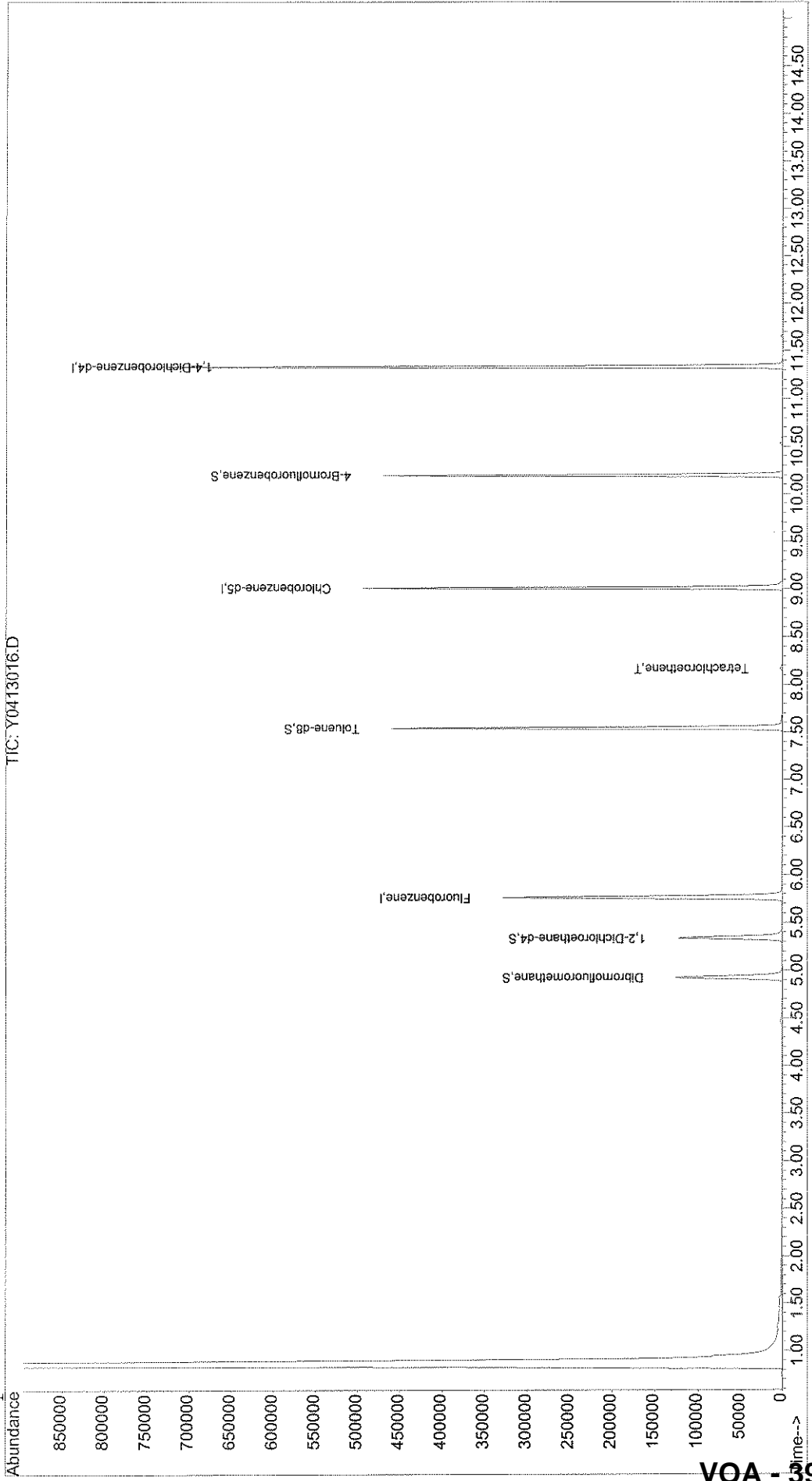
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413016.D  
Acq On : 13 Apr 2007 12:26  
Sample : JPL30-003  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:07 2007  
Vial: 30  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413016.D  
 Acq On : 13 Apr 2007 12:26  
 Sample : JPL30-003  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:07 2007

Vial: 30  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.76	96	303220	50.00	ug/l	0.00	89.25%
50) Chlorobenzene-d5	9.01	82	123663	50.00	ug/l	0.00	76.73%
70) 1,4-Dichlorobenzene-d4	11.34	152	164169	50.00	ug/l	0.00	72.85%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	93156	50.78	ug/l	0.00	
Spiked Amount	50.000	Range 85 - 115	Recovery	=	101.56%		
37) 1,2-Dichloroethane-d4	5.34	65	97380	50.03	ug/l	0.00	
Spiked Amount	50.000	Range 70 - 120	Recovery	=	100.06%		
51) Toluene-d8	7.54	98	271584	55.34	ug/l	0.00	
Spiked Amount	50.000	Range 85 - 120	Recovery	=	110.68%		
72) 4-Bromofluorobenzene	10.20	95	123717	52.16	ug/l	0.00	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	291	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.	d	
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413016.D 8260B.M Mon Apr 16 09:07:53 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413016.D  
 Acq On : 13 Apr 2007 12:26  
 Sample : JPL30-003  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:07 2007

Vial: 30  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.40	78	55		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.62	92	163		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	730	0.43	ug/l #✓	80
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	368		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	401		N.D.	
65) m,p-Xylene	9.29	106	82		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	453		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413016.D 8260B.M Mon Apr 16 09:07:53 2007



Quantitation Report

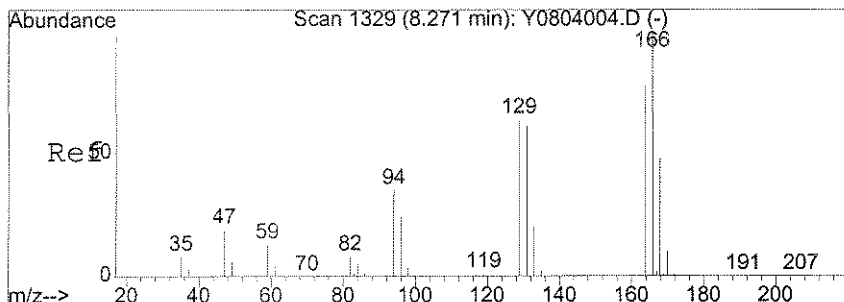
Data File : X:\MSVOA\YODA\041307\Y0413016.D  
 Acq On : 13 Apr 2007 12:26  
 Sample : JPL30-003  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:07 2007

Vial: 30  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

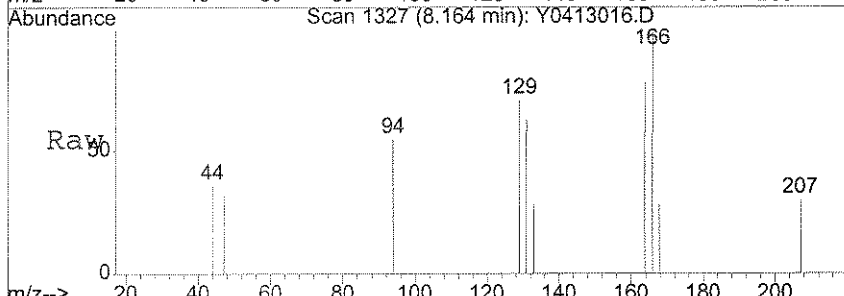
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	74		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	54		N.D.	
84) 4-Isopropyltoluene	11.34	119	269		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	54		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	257		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

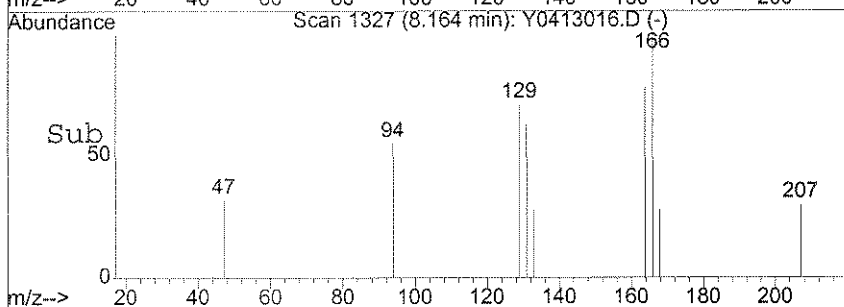
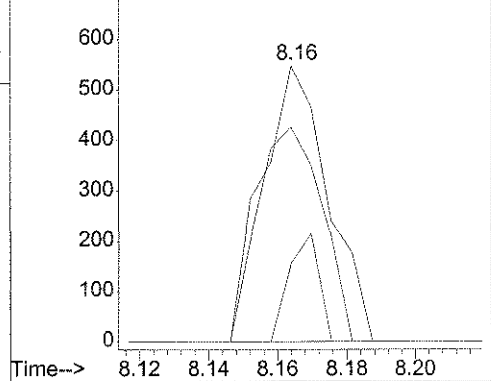


#56  
 Tetrachloroethene  
 Concen: 0.43 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. 0.00 min  
 Lab File: Y0413016.D  
 Acq: 13 Apr 2007 12:26

Tgt Ion	Resp	Lower	Upper
166	100		
164	76.0	63.3	94.9
168	17.9	39.6	59.4#



Abundance  
 Ion 165.95 (165.65 to 166.65): Y041301  
 Ion 163.95 (163.65 to 164.65): Y041301  
 Ion 167.95 (167.65 to 168.65): Y041301



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:50

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.26	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 12:50

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-004  
 Lab File ID: Y0413017.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 12:50  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

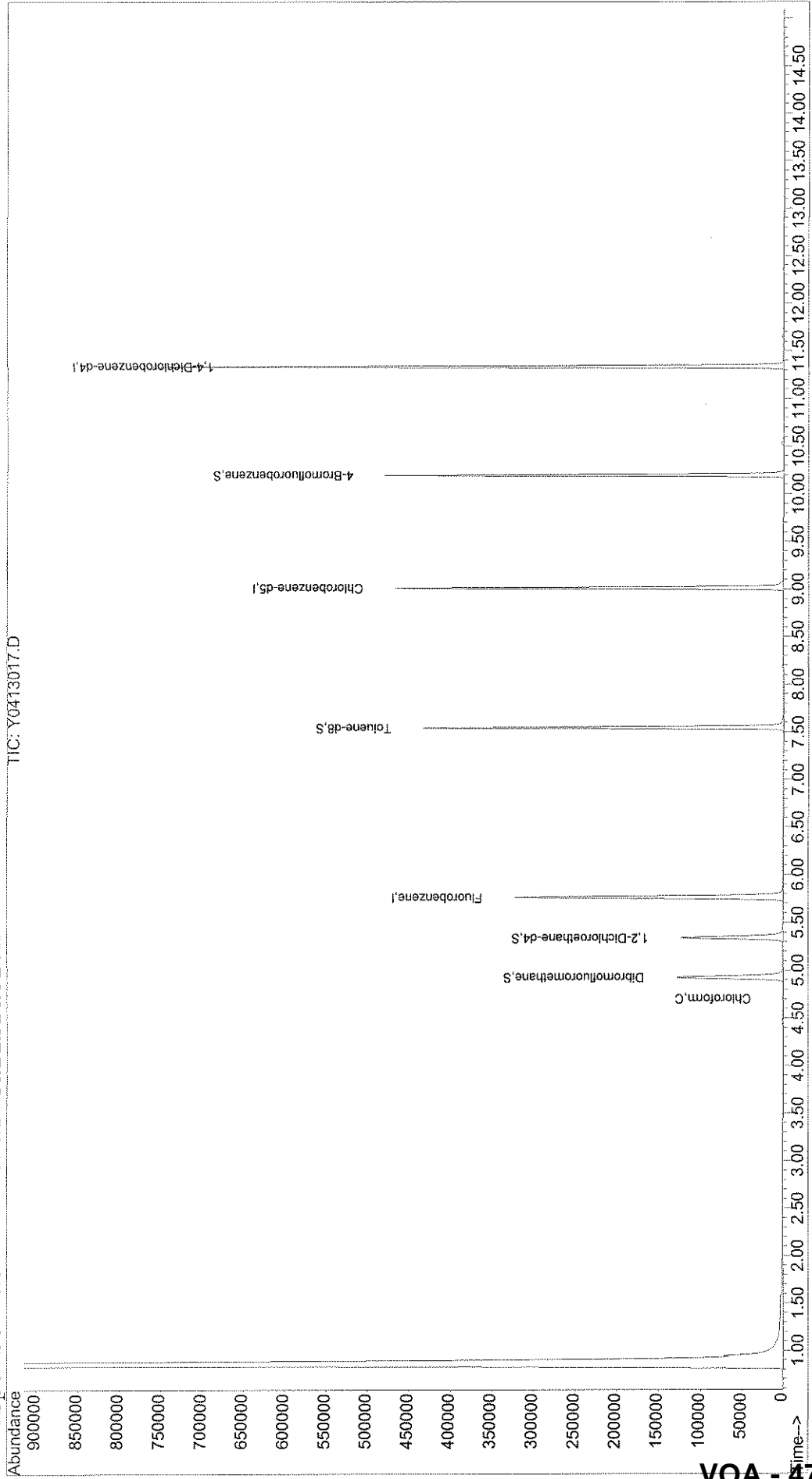
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413017.D Vial: 31  
Acq On : 13 Apr 2007 12:50 Operator: LNH  
Sample : JPL30-004 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:09 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 47

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413017.D  
 Acq On : 13 Apr 2007 12:50  
 Sample : JPL30-004  
 Misc : 5mL+IS/SS #3

Vial: 31  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:09 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.76	96	293701	50.00	ug/l	0.00	86.45%
50) Chlorobenzene-d5	9.01	82	119730	50.00	ug/l	0.00	74.29%
70) 1,4-Dichlorobenzene-d4	11.34	152	167118	50.00	ug/l	0.00	74.16%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	94015	52.91	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	105.82%	
37) 1,2-Dichloroethane-d4	5.34	65	96085	50.97	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.94%	
51) Toluene-d8	7.55	98	257381	54.17	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	108.34%	
72) 4-Bromofluorobenzene	10.20	95	122260	50.64	ug/l	0.00	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.21	76	141	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413017.D 8260B.M Mon Apr 16 09:09:32 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413017.D  
 Acq On : 13 Apr 2007 12:50  
 Sample : JPL30-004  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:09 2007

Vial: 31  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	974	0.26	ug/l ✓	94
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	317		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	138		N.D.	
65) m,p-Xylene	9.29	106	207		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*LNH 4/16/07*



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413017.D  
 Acq On : 13 Apr 2007 12:50  
 Sample : JPL30-004  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:09 2007

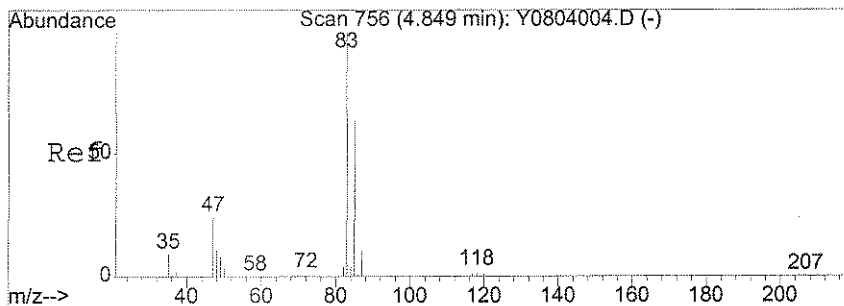
Vial: 31  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

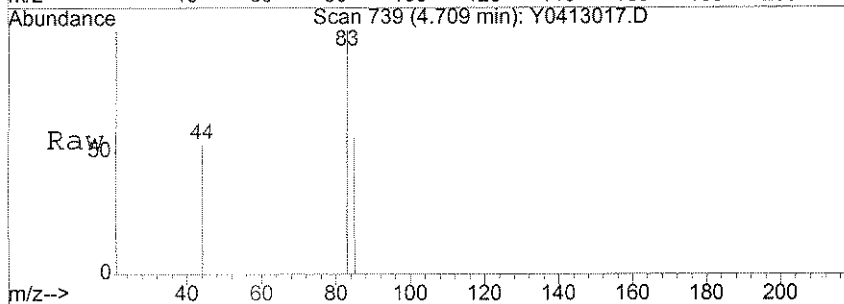
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	57		N.D.	
78) 4-Chlorotoluene	10.47	91	57		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	55		N.D.	
84) 4-Isopropyltoluene	11.34	119	210		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	55		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	235		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413017.D 8260B.M Mon Apr 16 09:09:33 2007

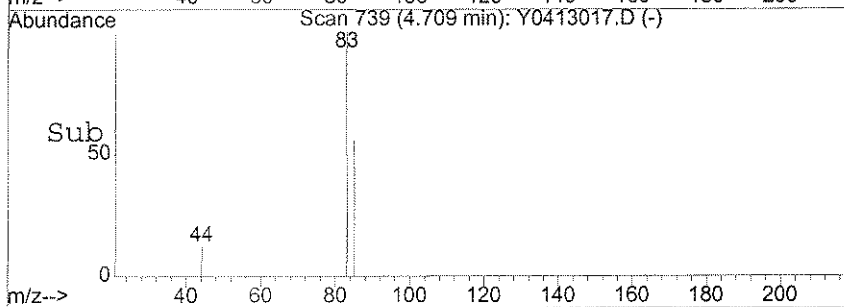
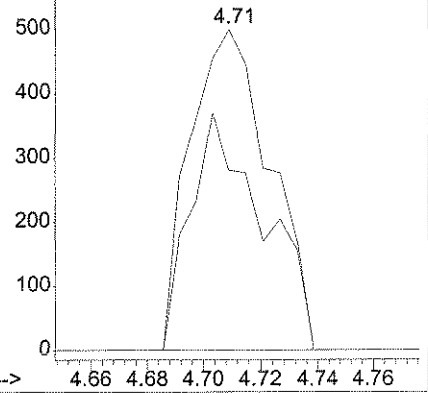


#31  
 Chloroform  
 Concen: 0.26 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0413017.D  
 Acq: 13 Apr 2007 12:50

Tgt Ion:	83	Resp:	974
Ion Ratio	Lower	Upper	
83	100		
85	67.8	43.3	83.3



Abundance Ion 83.00 (82.70 to 83.70): Y0413017.D  
 Ion 85.00 (84.70 to 85.70): Y0413017.D



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 13:14

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-005  
 Lab File ID: Y0413018.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 13:14  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-20-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 13:14

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

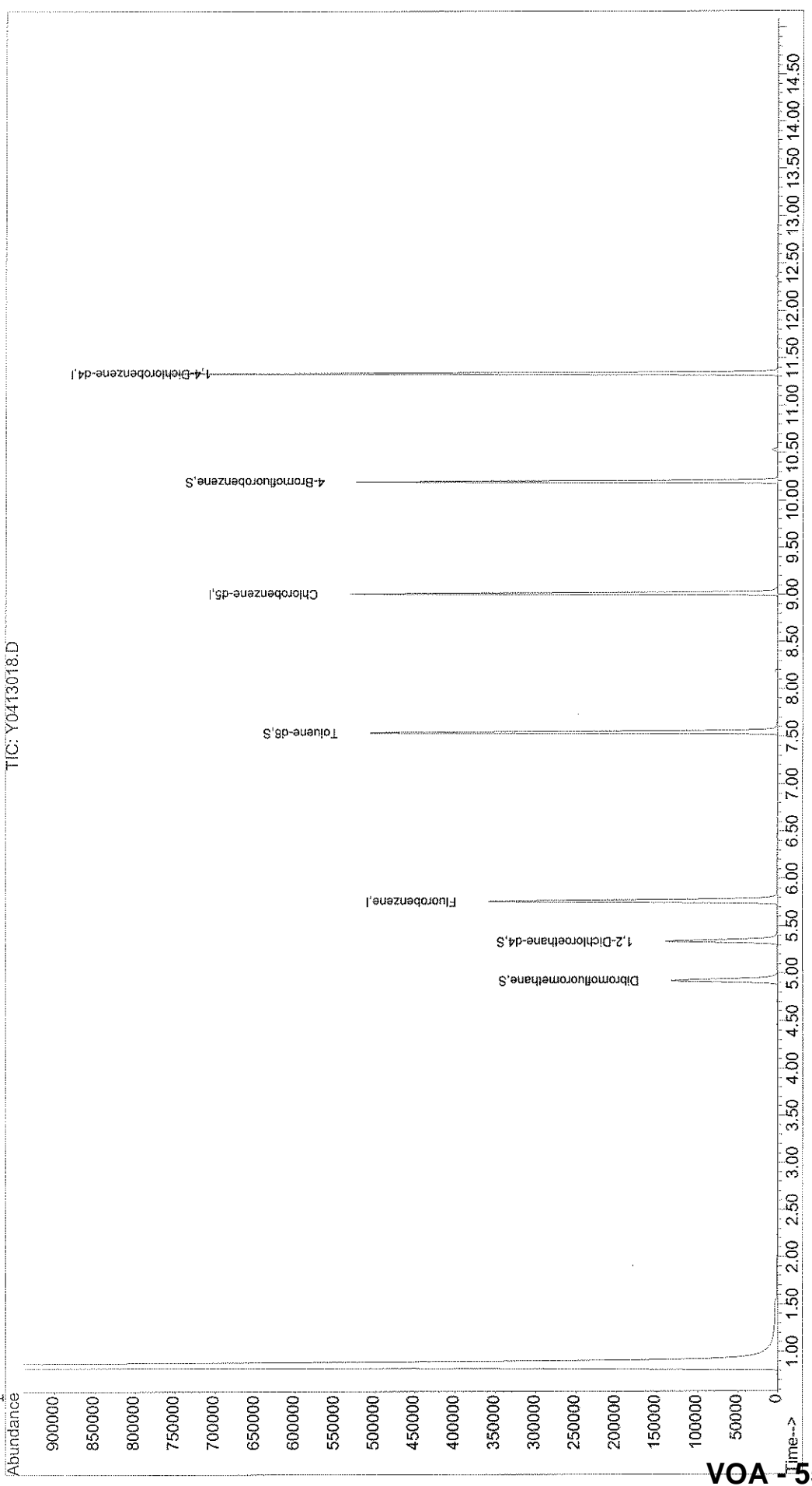
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413018.D Vial: 32  
Acq On : 13 Apr 2007 13:14 Operator: LNH  
Sample : JPL30-005 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:10 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413018.D  
 Acq On : 13 Apr 2007 13:14  
 Sample : JPL30-005  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

Vial: 32  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	333245	50.00	ug/l	0.00 98.09%
50) Chlorobenzene-d5	9.01	82	136379	50.00	ug/l	0.00 84.61%
70) 1,4-Dichlorobenzene-d4	11.34	152	174361	50.00	ug/l	0.00 77.37%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	99323	49.26	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	98.52%
37) 1,2-Dichloroethane-d4	5.34	65	106906	49.98	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.96%
51) Toluene-d8	7.55	98	305595	56.46	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.92%
72) 4-Bromofluorobenzene	10.20	95	134890	53.55	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	149		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413018.D 8260B.M Mon Apr 16 09:10:14 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413018.D  
 Acq On : 13 Apr 2007 13:14  
 Sample : JPL30-005  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

Vial: 32  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	778		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	6.79	83	53		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	433		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	132		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413018.D 8260B.M Mon Apr 16 09:10:14 2007



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413018.D  
 Acq On : 13 Apr 2007 13:14  
 Sample : JPL30-005  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

Vial: 32  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	114		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	70		N.D.	
78) 4-Chlorotoluene	10.47	91	70		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	269		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	140		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-3-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-006  
 Lab File ID: Y0413019.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 13:39  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.34	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-3-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-006  
 Lab File ID: Y0413019.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 13:39  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-3-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 13:39

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

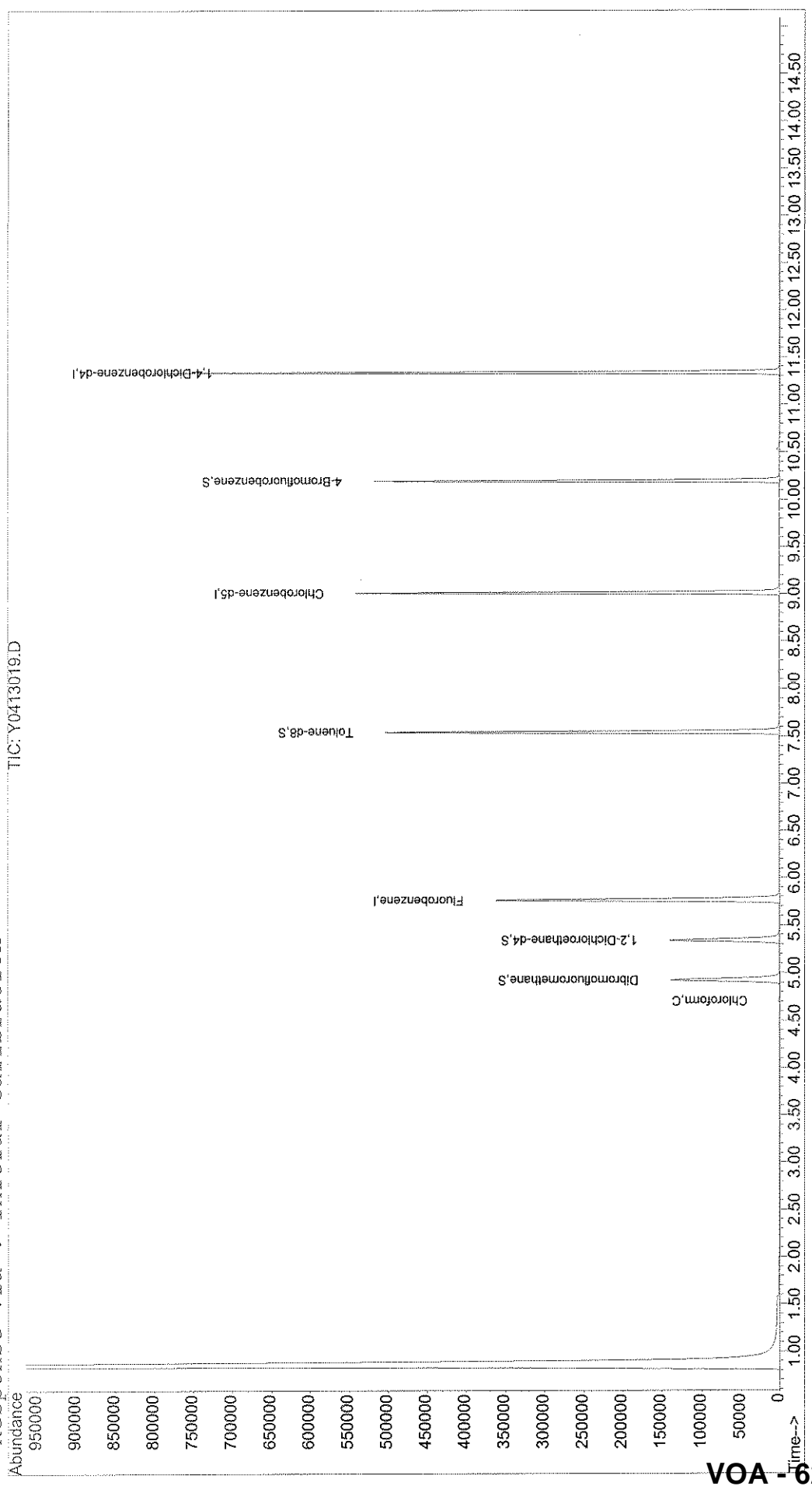
Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413019.D  
Acq On : 13 Apr 2007 13:39  
Sample : JPL30-006  
Misc : 5mL+IS/SS #2  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:10 2007

Vial: 33  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413019.D  
 Acq On : 13 Apr 2007 13:39  
 Sample : JPL30-006  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

Vial: 33  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	333454	50.00	ug/l	0.00 98.15%
50) Chlorobenzene-d5	9.01	82	137177	50.00	ug/l	0.00 85.11%
70) 1,4-Dichlorobenzene-d4	11.34	152	174904	50.00	ug/l	0.00 77.61%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	99524	49.33	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	98.66%	
37) 1,2-Dichloroethane-d4	5.34	65	108210	50.56	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	101.12%	
51) Toluene-d8	7.55	98	301048	55.30	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	110.60%	
72) 4-Bromofluorobenzene	10.20	95	133775	52.94	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	65	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413019.D 8260B.M Mon Apr 16 09:11:01 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413019.D  
 Acq On : 13 Apr 2007 13:39  
 Sample : JPL30-006  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

Vial: 33  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.71	83	1465	0.34 ug/l ✓	93
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	0.00	78	0	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	0.00	130	0	N.D.	
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	0.00	92	0	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	8.17	166	200	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	9.01	91	392	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.17	91	64	N.D.	
65) m,p-Xylene	9.30	106	134	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413019.D 8260B.M Mon Apr 16 09:11:01 2007

*Signature*

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413019.D  
 Acq On : 13 Apr 2007 13:39  
 Sample : JPL30-006  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:10 2007

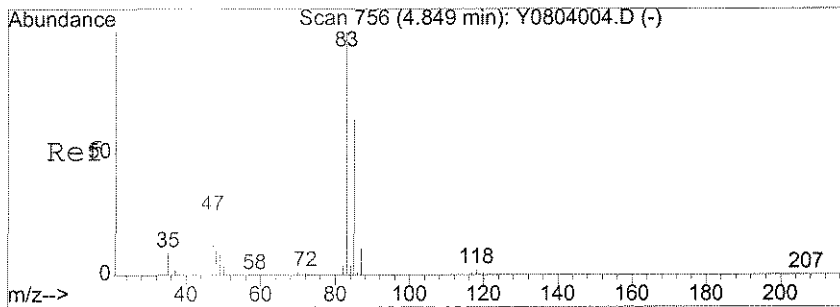
Vial: 33  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

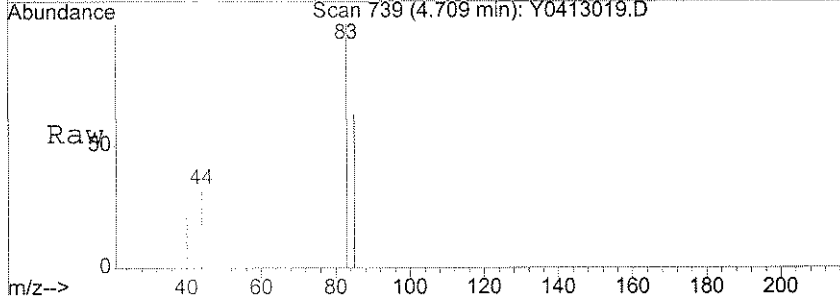
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	54		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	53		N.D.	
82) sec-butylbenzene	11.18	105	53		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	290		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	143		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



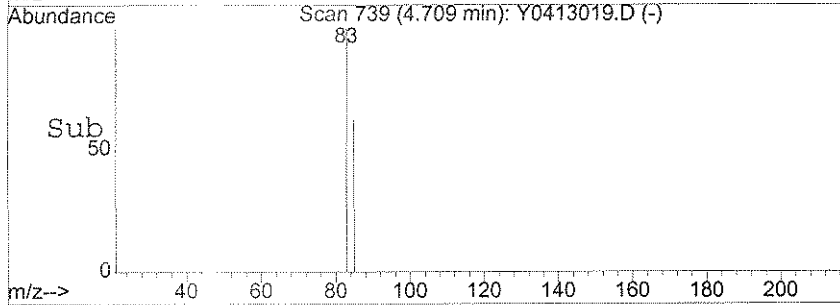
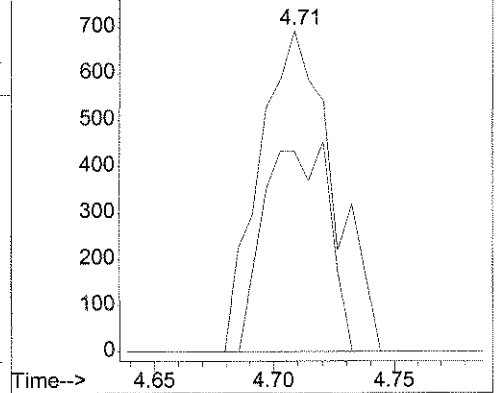


#31  
 Chloroform  
 Concen: 0.34 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0413019.D  
 Acq: 13 Apr 2007 13:39

Tgt Ion	Resp	Lower	Upper
83	1465		
83	100		
85	57.5	43.3	83.3



Abundance on 83.00 (82.70 to 83.70): Y0413019.D  
 800 Ion 85.00 (84.70 to 85.70): Y0413019.D



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-007  
 Lab File ID: Y0413009.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 09:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-007  
 Lab File ID: Y0413009.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 09:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	0.50		U
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	0.50		U
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-007  
 Lab File ID: Y0413009.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 09:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

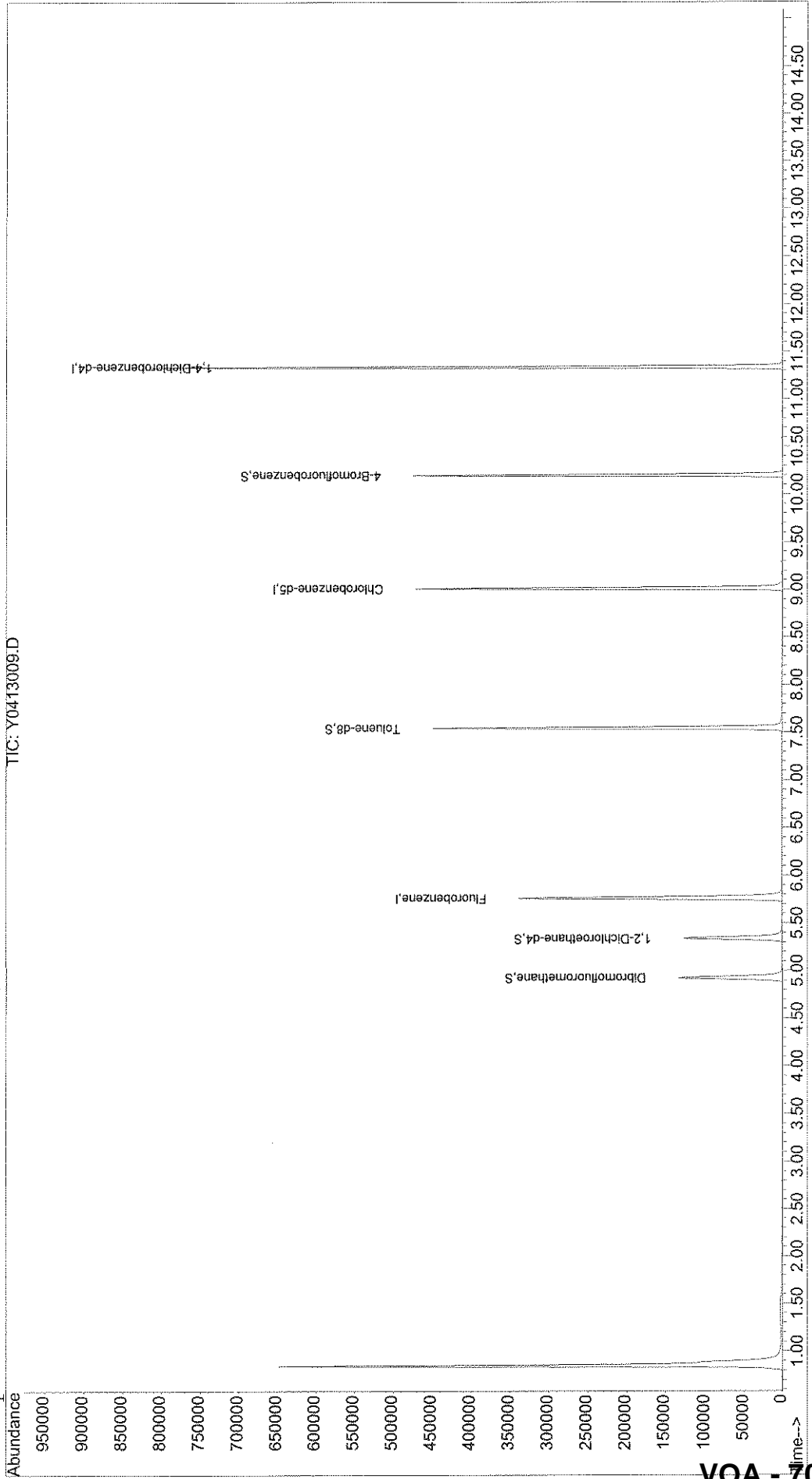
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413009.D  
Acq On : 13 Apr 2007 9:33 Vial: 24  
Sample : JPL30-007 EB Operator: LNH  
Misc : 5mL+IS/SS #1 Inst : yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 16 9:01 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 70

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413009.D  
 Acq On : 13 Apr 2007 9:33  
 Sample : JPL30-007 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:01 2007

Vial: 24  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	5.76	96	315215	50.00	ug/l	0.00 92.78%
50) Chlorobenzene-d5	9.01	82	122975	50.00	ug/l	0.00 76.30%
70) 1,4-Dichlorobenzene-d4	11.34	152	177246	50.00	ug/l	0.00 78.65%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	100158	52.52	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	105.04%
37) 1,2-Dichloroethane-d4	5.34	65	101652	50.24	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.48%
51) Toluene-d8	7.54	98	271891	55.71	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	111.42%
72) 4-Bromofluorobenzene	10.20	95	128977	50.37	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	d
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	379		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413009.D  
 Acq On : 13 Apr 2007 9:33  
 Sample : JPL30-007 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:01 2007

Vial: 24  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	4.35	43	151		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.40	78	56		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.03	112	54		N.D.	
62) 1-Chlorohexane	9.02	91	360		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	775		N.D.	
65) m,p-Xylene	9.29	106	675		N.D.	
66) o-xylene	9.69	106	112		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413009.D 8260B.M Mon Apr 16 09:01:35 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413009.D  
 Acq On : 13 Apr 2007 9:33  
 Sample : JPL30-007 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:01 2007

Vial: 24  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	54		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.53	91	122		N.D.	
78) 4-Chlorotoluene	10.64	91	133		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	134		N.D.	
80) tert-Butylbenzene	10.96	119	63		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	189		N.D.	
82) sec-butylbenzene	11.18	105	340		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	146		N.D.	
84) 4-Isopropyltoluene	11.33	119	700		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	287		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	599		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.30	180	186		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	13.55	128	358		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	145		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413009.D 8260B.M Mon Apr 16 09:01:36 2007



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413010.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 09:58

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
75-71-8	Dichlorodifluoromethane	0.50		U
74-87-3	Chloromethane	0.50		U
75-01-4	Vinyl chloride	0.50		U
74-83-9	Bromomethane	0.50		U
75-00-3	Chloroethane	0.50		U
75-69-4	Trichlorofluoromethane	0.50		U
75-35-4	1,1-Dichloroethene	0.50		U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50		U
75-09-2	Methylene chloride	0.50		U
1634-04-4	Methyl tert-butyl ether	0.50		U
156-60-5	trans-1,2-Dichloroethene	0.50		U
75-34-3	1,1-Dichloroethane	0.50		U
594-20-7	2,2-Dichloropropane	0.50		U
156-59-2	cis-1,2-Dichloroethene	0.50		U
78-93-3	2-Butanone	5.0		U
74-97-5	Bromochloromethane	0.50		U
67-66-3	Chloroform	0.50		U
71-55-6	1,1,1-Trichloroethane	0.50		U
56-23-5	Carbon tetrachloride	0.50		U
563-58-6	1,1-Dichloropropene	0.50		U
71-43-2	Benzene	0.50		U
107-06-2	1,2-Dichloroethane	0.50		U
79-01-6	Trichloroethene	0.50		U
78-87-5	1,2-Dichloropropane	0.50		U
74-95-3	Dibromomethane	0.50		U
75-27-4	Bromodichloromethane	0.50		U
10061-01-	cis-1,3-Dichloropropene	0.50		U
108-10-1	4-Methyl-2-pentanone	5.0		U
108-88-3	Toluene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-008  
 Lab File ID: Y0413010.D  
 Date Collected: 03/30/2007  
 Date/Time Analyzed: 04/13/2007 09:58  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413010.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/30/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 09:58

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

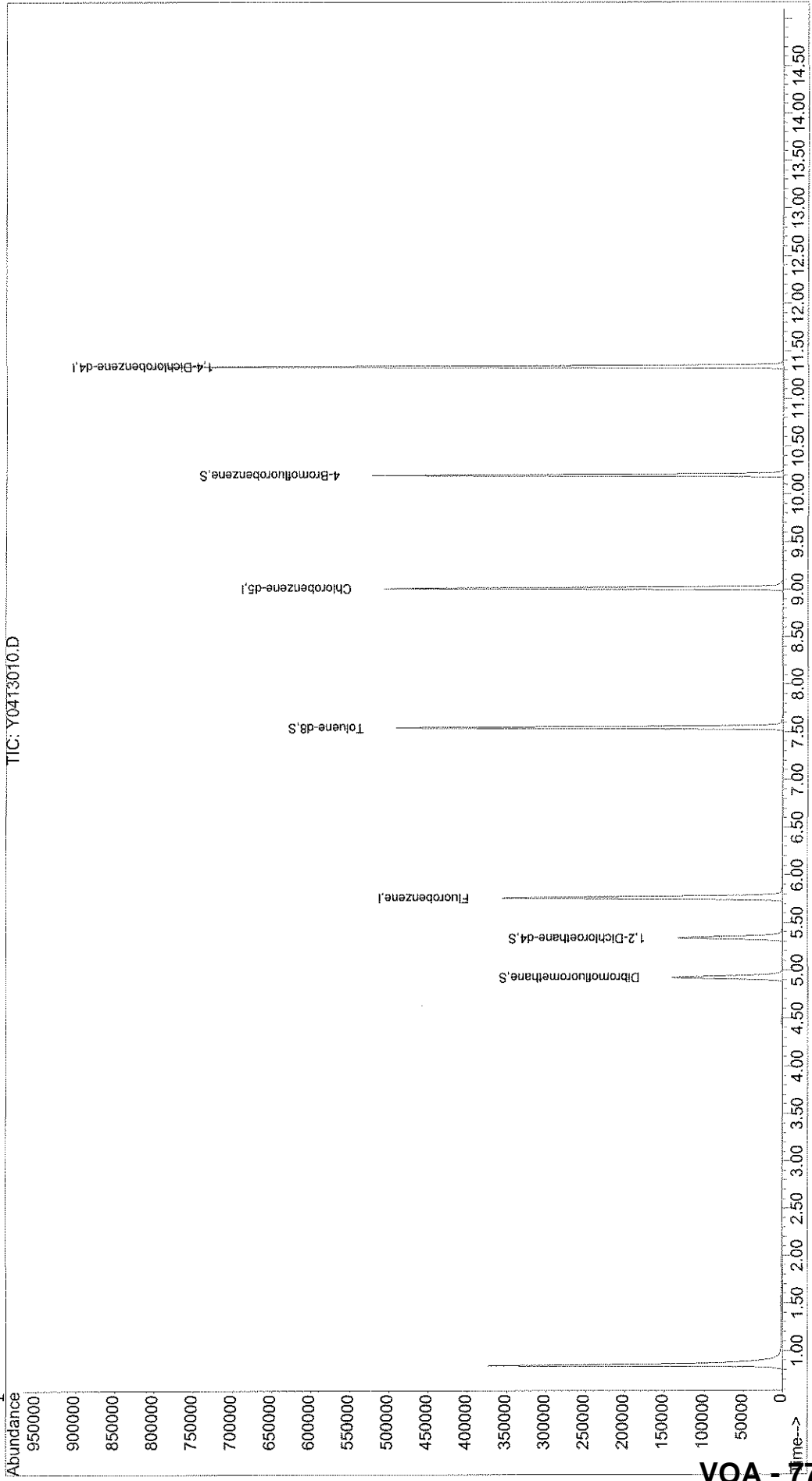
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413010.D  
Acq On : 13 Apr 2007 9:58 Vial: 25  
Sample : JPL30-008 TB Operator: LNH  
Misc : 5mL+IS/SS #1 Inst : Yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 16 9:02 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 77

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413010.D  
 Acq On : 13 Apr 2007 9:58  
 Sample : JPL30-008 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:02 2007

Vial: 25  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.76	96	326599	50.00	ug/l	0.00	96.13%
50) Chlorobenzene-d5	9.01	82	131101	50.00	ug/l	0.00	81.34%
70) 1,4-Dichlorobenzene-d4	11.34	152	182244	50.00	ug/l	0.00	80.87%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101603	51.42	ug/l	0.00	
Spiked Amount	50.000	Range 85 - 115	Recovery	=	102.84%		
37) 1,2-Dichloroethane-d4	5.34	65	105098	50.13	ug/l	0.00	
Spiked Amount	50.000	Range 70 - 120	Recovery	=	100.26%		
51) Toluene-d8	7.54	98	288636	55.48	ug/l	0.00	
Spiked Amount	50.000	Range 85 - 120	Recovery	=	110.96%		
72) 4-Bromofluorobenzene	10.20	95	136173	51.72	ug/l	0.00	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	730	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	2.51	84	261	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413010.D  
 Acq On : 13 Apr 2007 9:58  
 Sample : JPL30-008 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:02 2007

Vial: 25  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	380		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	66		N.D.	
65) m,p-Xylene	9.29	106	108		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413010.D 8260B.M Mon Apr 16 09:02:21 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413010.D  
 Acq On : 13 Apr 2007 9:58  
 Sample : JPL30-008 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:02 2007

Vial: 25  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.06	105	55		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	368		N.D.	
78) 4-Chlorotoluene	10.64	91	142		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	160		N.D.	
82) sec-butylbenzene	11.18	105	335		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	65		N.D.	
84) 4-Isopropyltoluene	11.34	119	446		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	65		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	464		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	135		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	13.54	128	179		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	69		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 14:34

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.31	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.33	J
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-009  
 Lab File ID: Y0413021.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 14:34  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	2.9	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-009  
 Lab File ID: Y0413021.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 14:34  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

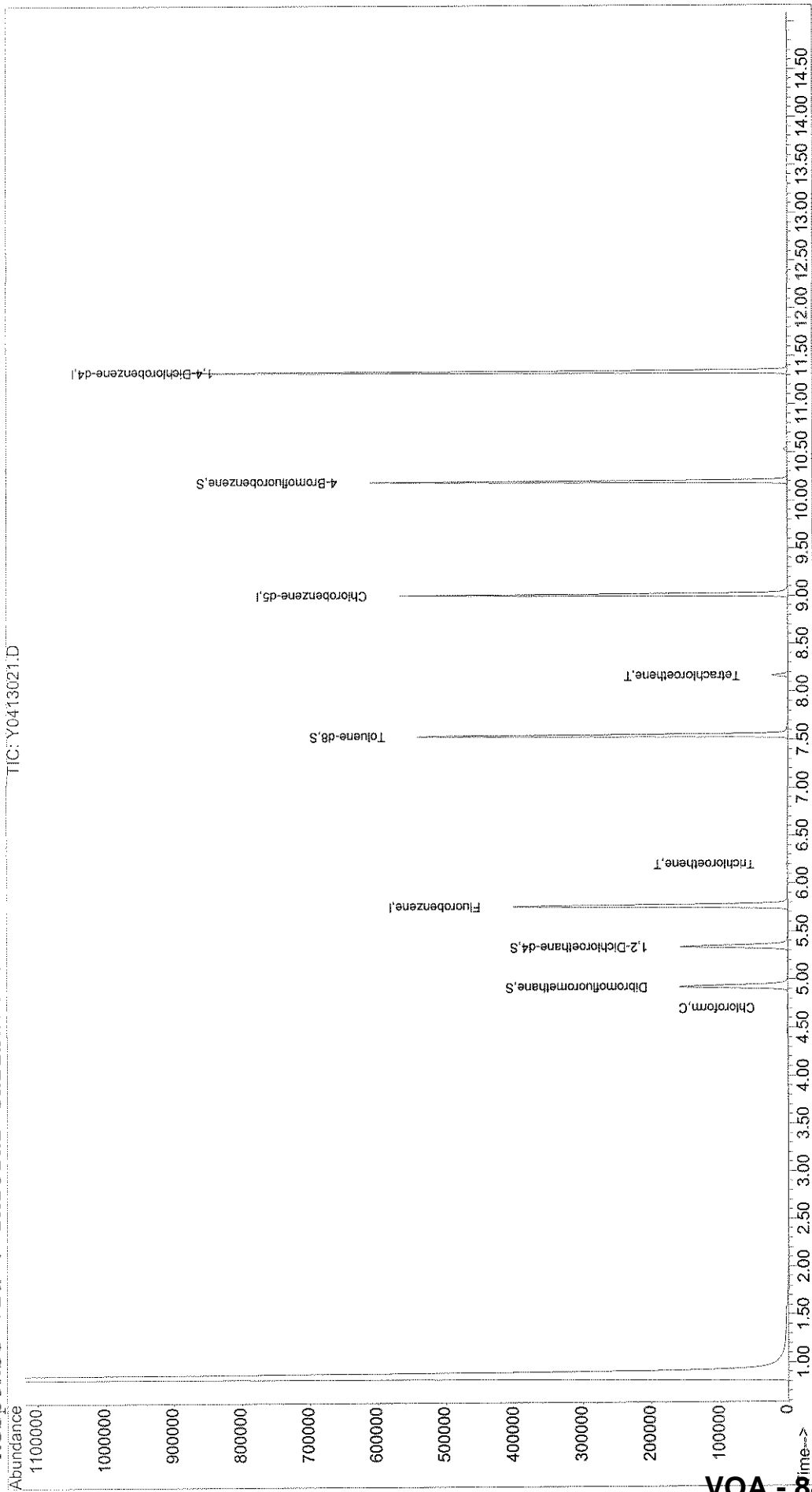
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413021.D Vial: 35  
Acq On : 13 Apr 2007 14:34 Operator: LNH  
Sample : JPL30-009 Inst : Yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:12 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413021.D  
 Acq On : 13 Apr 2007 14:34  
 Sample : JPL30-009  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:12 2007

Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	362333	50.00	ug/l	0.00 106.65%
50) Chlorobenzene-d5	9.01	82	145481	50.00	ug/l	0.00 90.26%
70) 1,4-Dichlorobenzene-d4	11.34	152	207324	50.00	ug/l	0.00 92.00%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	113156	51.62	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	103.24%	
37) 1,2-Dichloroethane-d4	5.34	65	118478	50.94	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	101.88%	
51) Toluene-d8	7.54	98	320360	55.49	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	110.98%	
72) 4-Bromofluorobenzene	10.20	95	151324	50.52	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	76		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	3.32	63	484		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413021.D 8260B.M Mon Apr 16 09:12:29 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413021.D  
 Acq On : 13 Apr 2007 14:34  
 Sample : JPL30-009  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:12 2007

Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	4.70	83	1464	0.31	ug/l ✓	92
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	5.38	78	73	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.		
41) Trichloroethene	6.20	130	836	0.33	ug/l # ✓	53
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	0.00	41	0	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	0.00	92	0	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.		
56) Tetrachloroethene	8.16	166	5773	2.87	ug/l ✓	97
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.		
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	0.00	112	0	N.D.		
62) 1-Chlorohexane	9.02	91	488	N.D.		
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.29	91	256	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.		
66) o-xylene	0.00	106	0	N.D.		
67) Styrene	0.00	104	0	N.D.		
68) Bromoform	0.00	173	0	N.D.		

*LNH 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0413021.D 8260B.M Mon Apr 16 09:12:29 2007

Quantitation Report

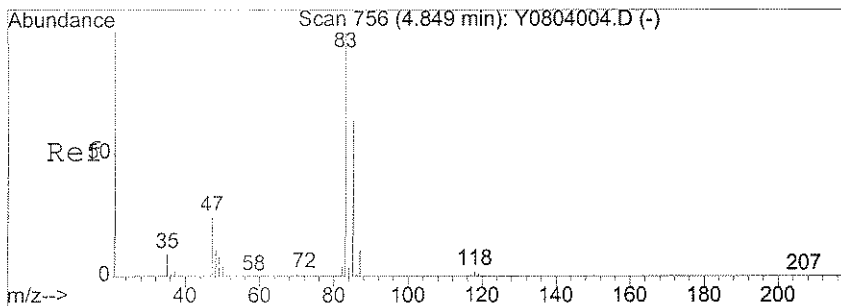
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 Acq On : 13 Apr 2007 14:34  
 Sample : JPL30-009  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:12 2007

Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

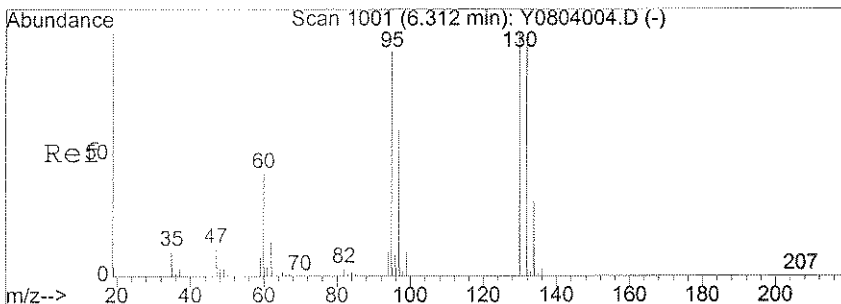
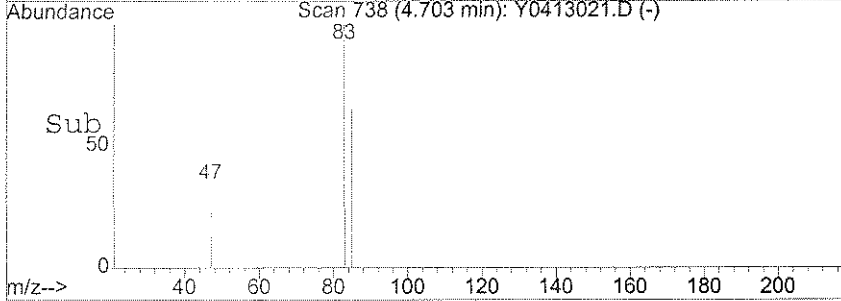
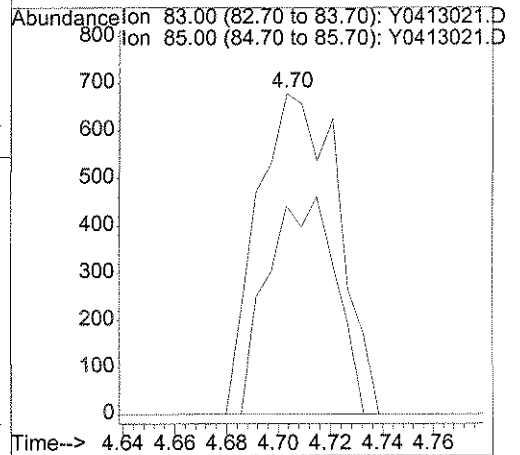
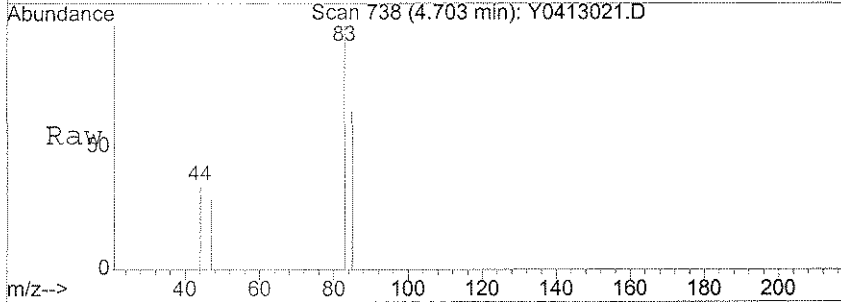
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	293		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	62		N.D.	
78) 4-Chlorotoluene	10.46	91	62		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	59		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	58		N.D.	
82) sec-butylbenzene	11.18	105	58		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	127		N.D.	
84) 4-Isopropyltoluene	11.34	119	259		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	109		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	81		N.D.	
87) n-Butylbenzene	11.74	91	186		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	61		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	53		N.D.	



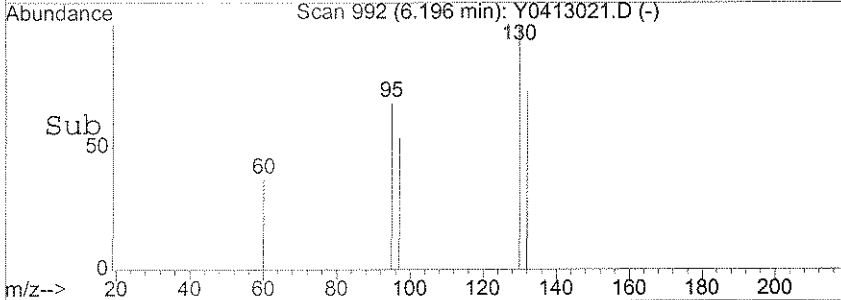
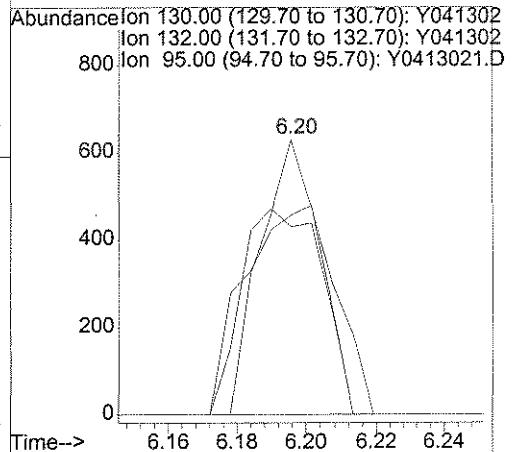
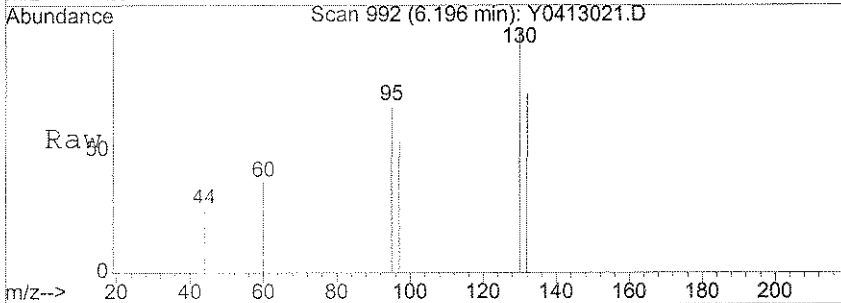
#31  
 Chloroform  
 Concen: 0.31 ug/l  
 RT: 4.70 min Scan# 738  
 Delta R.T. 0.00 min  
 Lab File: Y0413021.D  
 Acq: 13 Apr 2007 14:34

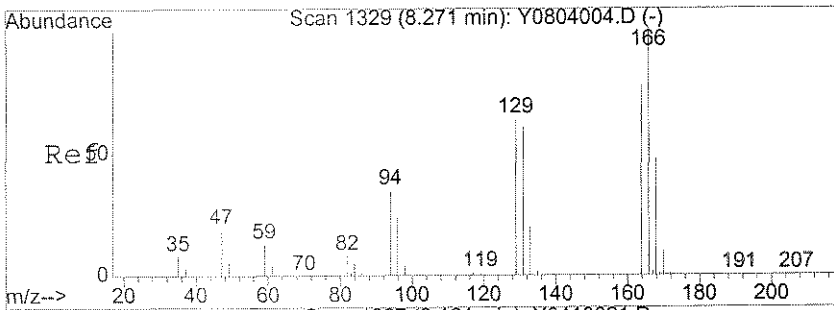
Tgt Ion: 83 Resp: 1464  
 Ion Ratio Lower Upper  
 83 100  
 85 57.0 43.3 83.3



#41  
 Trichloroethene  
 Concen: 0.33 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0413021.D  
 Acq: 13 Apr 2007 14:34

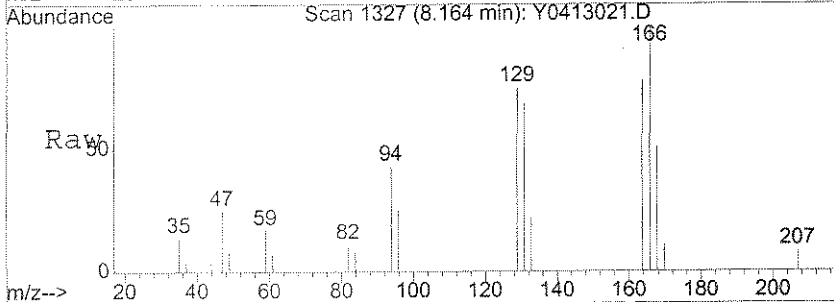
Tgt Ion: 130 Resp: 836  
 Ion Ratio Lower Upper  
 130 100  
 132 93.3 75.0 115.0  
 95 0.0 69.4 109.4#



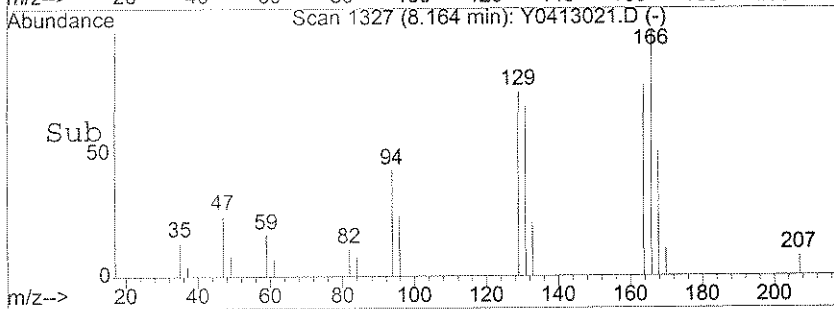
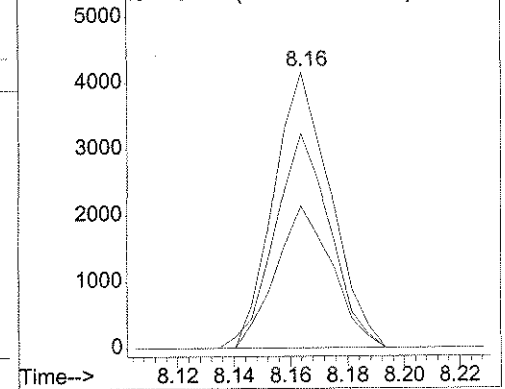


#56  
 Tetrachloroethene  
 Concen: 2.87 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. 0.00 min  
 Lab File: Y0413021.D  
 Acq: 13 Apr 2007 14:34

Tgt Ion	Resp	Lower	Upper
166	5773		
166	100		
164	75.9	63.3	94.9
168	51.1	39.6	59.4



Abundance  
 Ion 165.95 (165.65 to 166.65): Y041302  
 Ion 163.95 (163.65 to 164.65): Y041302  
 Ion 167.95 (167.65 to 168.65): Y041302





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 14:59

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.28	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-010  
 Lab File ID: Y0413022.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 14:59  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.71	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 14:59

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

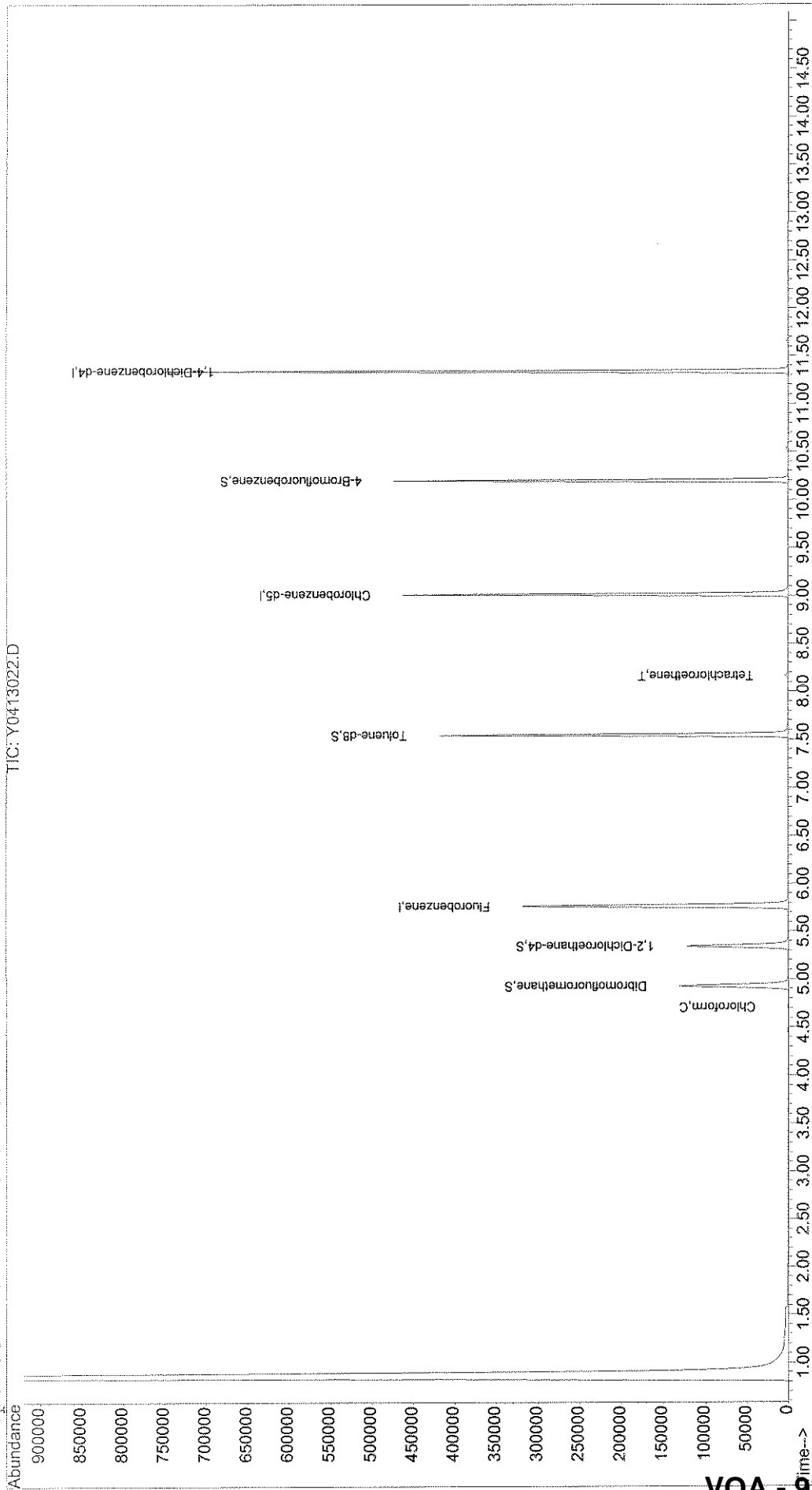
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413022.D  
Acq On : 13 Apr 2007 14:59 Vial: 36  
Sample : JPL30-010 Operator: LNH  
Misc : 5mL+IS/SS #1 Inst : Yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 16 9:13 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413022.D  
 Acq On : 13 Apr 2007 14:59  
 Sample : JPL30-010  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	292922	50.00	ug/l	0.00 86.22%
50) Chlorobenzene-d5	9.01	82	119109	50.00	ug/l	0.00 73.90%
70) 1,4-Dichlorobenzene-d4	11.34	152	168880	50.00	ug/l	0.00 74.94%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	94342	53.23	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	106.46%
37) 1,2-Dichloroethane-d4	5.34	65	95965	51.04	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.08%
51) Toluene-d8	7.55	98	253644	53.66	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	107.32%
72) 4-Bromofluorobenzene	10.20	95	122149	50.06	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.32	63	59	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413022.D 8260B.M Mon Apr 16 09:13:15 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413022.D  
 Acq On : 13 Apr 2007 14:59  
 Sample : JPL30-010  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.71	83	1073	0.28 ug/l ✓	98
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	0.00	78	0	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	6.20	130	269	N.D.	
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	0.00	92	0	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	8.17	166	1174	0.71 ug/l ✓	85
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	9.01	91	278	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.17	91	70	N.D.	
65) m,p-Xylene	9.29	106	109	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D.	

*QNH 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0413022.D 8260B.M Mon Apr 16 09:13:16 2007

Quantitation Report

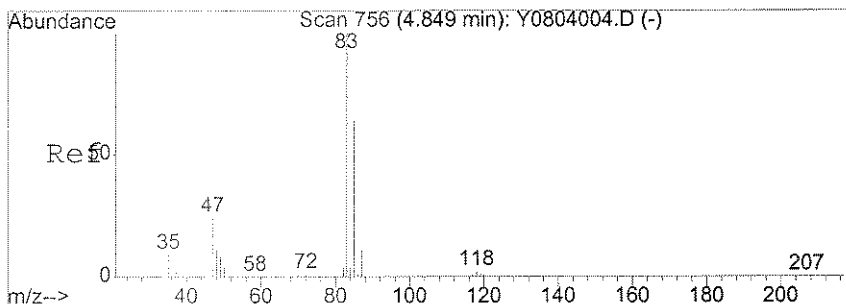
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 Acq On : 13 Apr 2007 14:59  
 Sample : JPL30-010  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

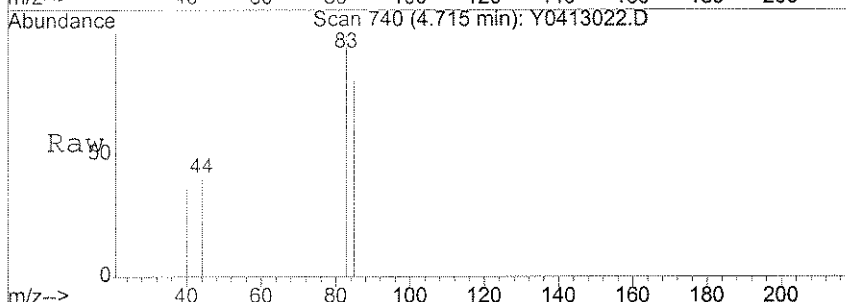
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	127		N.D.	
84) 4-Isopropyltoluene	11.33	119	188		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	111		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

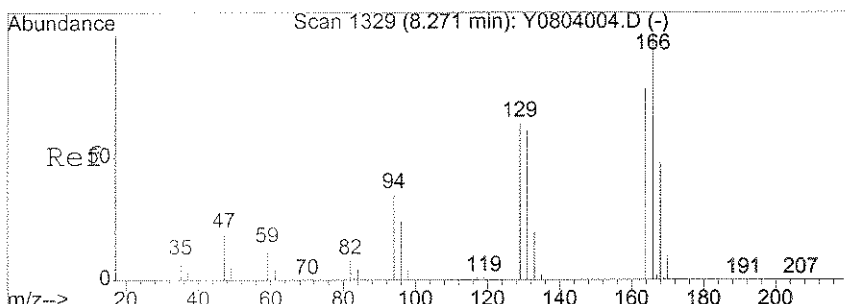
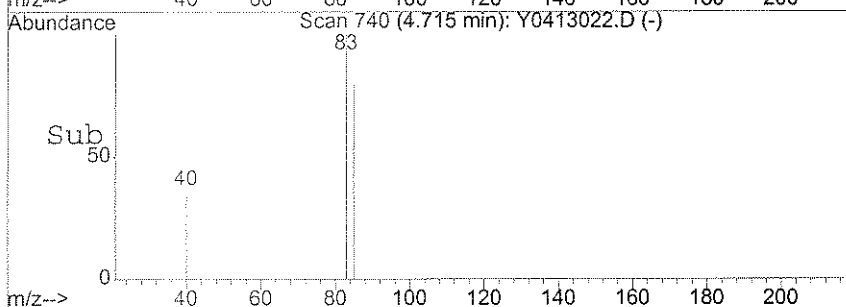
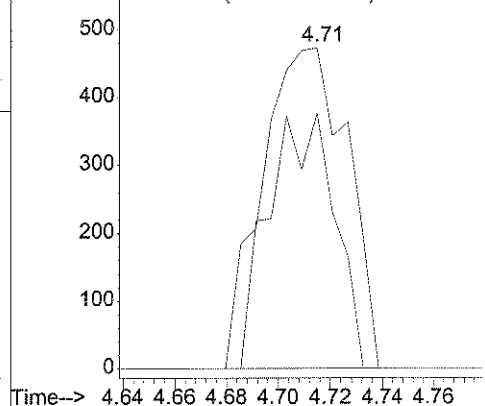


#31  
 Chloroform  
 Concen: 0.28 ug/l  
 RT: 4.71 min Scan# 740  
 Delta R.T. 0.01 min  
 Lab File: Y0413022.D  
 Acq: 13 Apr 2007 14:59

Tgt Ion	Resp	Lower	Upper
83	1073		
85	61.8	43.3	83.3

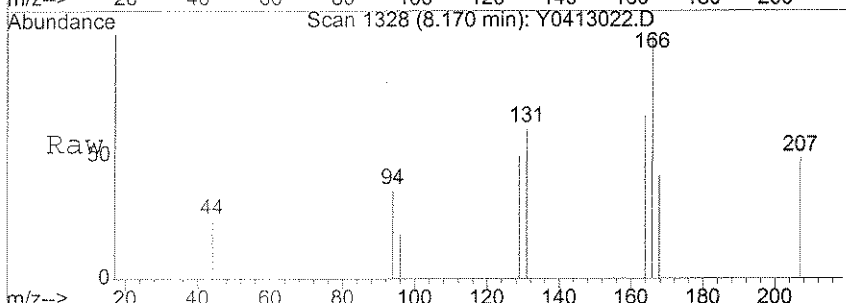


Abundance Ion 83.00 (82.70 to 83.70): Y0413022.D  
 Ion 85.00 (84.70 to 85.70): Y0413022.D

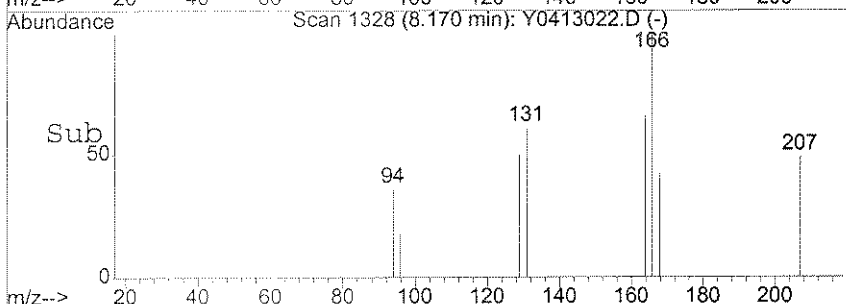
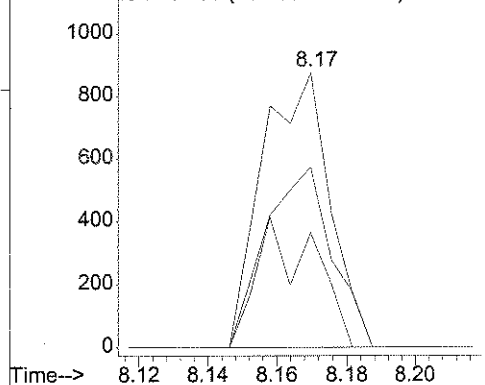


#56  
 Tetrachloroethene  
 Concen: 0.71 ug/l  
 RT: 8.17 min Scan# 1328  
 Delta R.T. 0.01 min  
 Lab File: Y0413022.D  
 Acq: 13 Apr 2007 14:59

Tgt Ion	Resp	Lower	Upper
166	1174		
164	64.7	63.3	94.9
168	40.3	39.6	59.4



Abundance Ion 165.95 (165.65 to 166.65): Y041302  
 Ion 163.95 (163.65 to 164.65): Y041302  
 Ion 167.95 (167.65 to 168.65): Y041302





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-011  
 Lab File ID: Y0413023.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 15:23  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-011  
 Lab File ID: Y0413023.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 15:23  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.78	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-011  
 Lab File ID: Y0413023.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 15:23  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

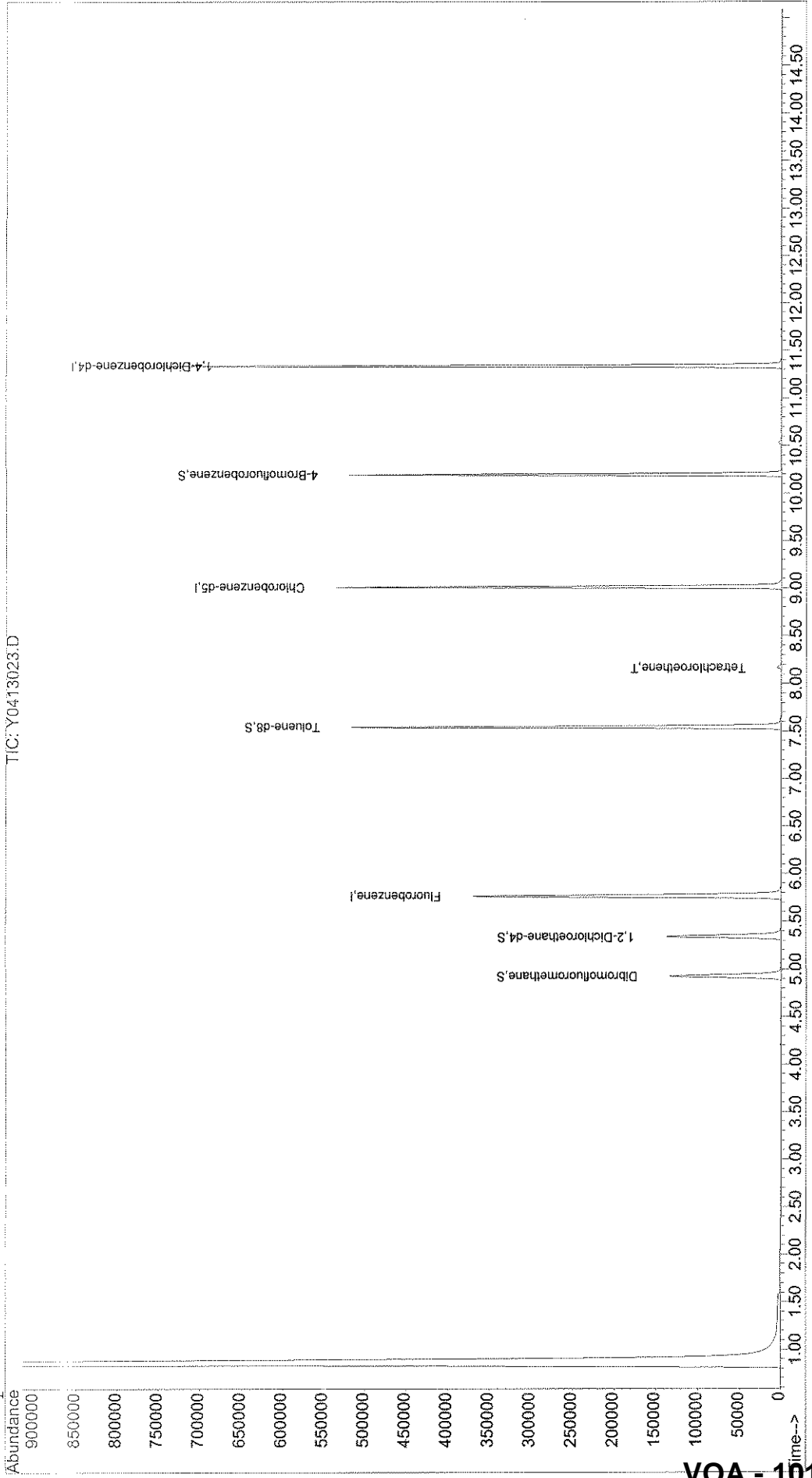
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413023.D  
Acq On : 13 Apr 2007 15:23  
Sample : JPL30-011  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:13 2007  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 101

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413023.D  
 Acq On : 13 Apr 2007 15:23  
 Sample : JPL30-011  
 Misc : 5mL+IS/SS #1

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

Response via : Initial Calibration

DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.77	96	340145	50.00	ug/l	0.00	100.12%
50) Chlorobenzene-d5	9.01	82	135137	50.00	ug/l	0.00	83.84%
70) 1,4-Dichlorobenzene-d4	11.34	152	172527	50.00	ug/l	0.00	76.56%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	97739	47.49	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	94.98%	
37) 1,2-Dichloroethane-d4	5.34	65	108158	49.54	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.08%	
51) Toluene-d8	7.55	98	303268	56.55	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	113.10%	
72) 4-Bromofluorobenzene	10.20	95	132625	53.21	ug/l	0.00	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	3.33	63	366		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration

Y0413023.D 8260B.M Mon Apr 16 09:14:03 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413023.D  
 Acq On : 13 Apr 2007 15:23  
 Sample : JPL30-011  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	4.21	96	600		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	767		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	205		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	1459	0.78	ug/l ✓	93
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	471		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	208		N.D.	
65) m,p-Xylene	9.29	106	53		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*LNH 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0413023.D 8260B.M Mon Apr 16 09:14:04 2007

Quantitation Report

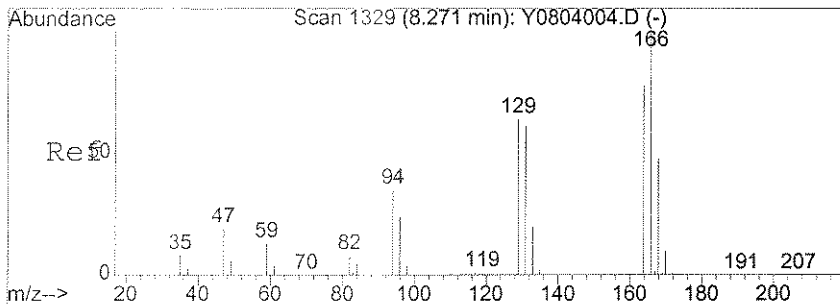
Data File : X:\MSVOA\YODA\041307\Y0413023.D  
 Acq On : 13 Apr 2007 15:23  
 Sample : JPL30-011  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:13 2007

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

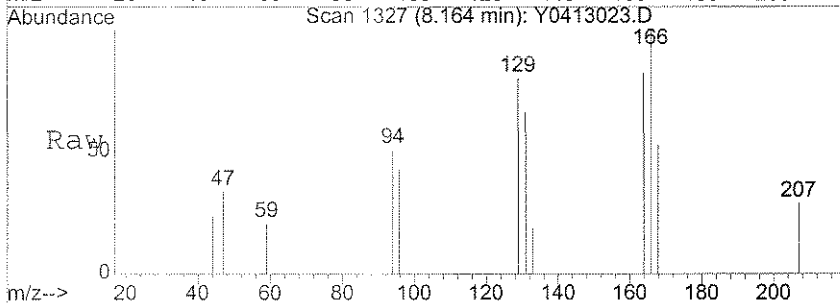
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	320		N.D.	
84) 4-Isopropyltoluene	11.33	119	196		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	238		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	283		N.D.	
87) n-Butylbenzene	11.74	91	131		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	65		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	61		N.D.	

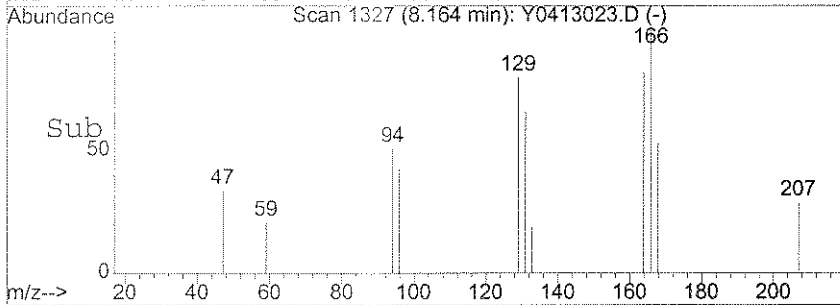
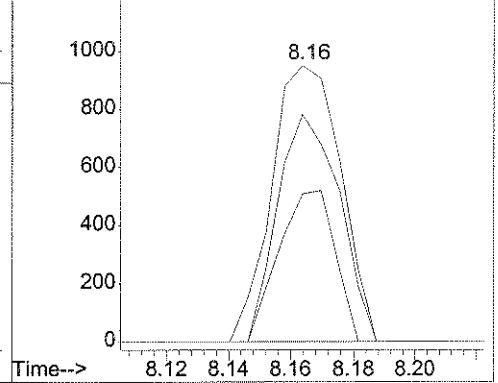


#56  
 Tetrachloroethene  
 Concen: 0.78 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. 0.00 min  
 Lab File: Y0413023.D  
 Acq: 13 Apr 2007 15:23

Tgt Ion	Resp	Lower	Upper
166	1459		
166	100		
164	73.5	63.3	94.9
168	44.3	39.6	59.4



Abundance  
 Ion 165.95 (165.65 to 166.65): Y041302  
 Ion 163.95 (163.65 to 164.65): Y041302  
 Ion 167.95 (167.65 to 168.65): Y041302





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 15:48

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.28	J
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.45	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.2	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-012  
 Lab File ID: Y0413024.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 15:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.62	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-012  
 Lab File ID: Y0413024.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 15:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

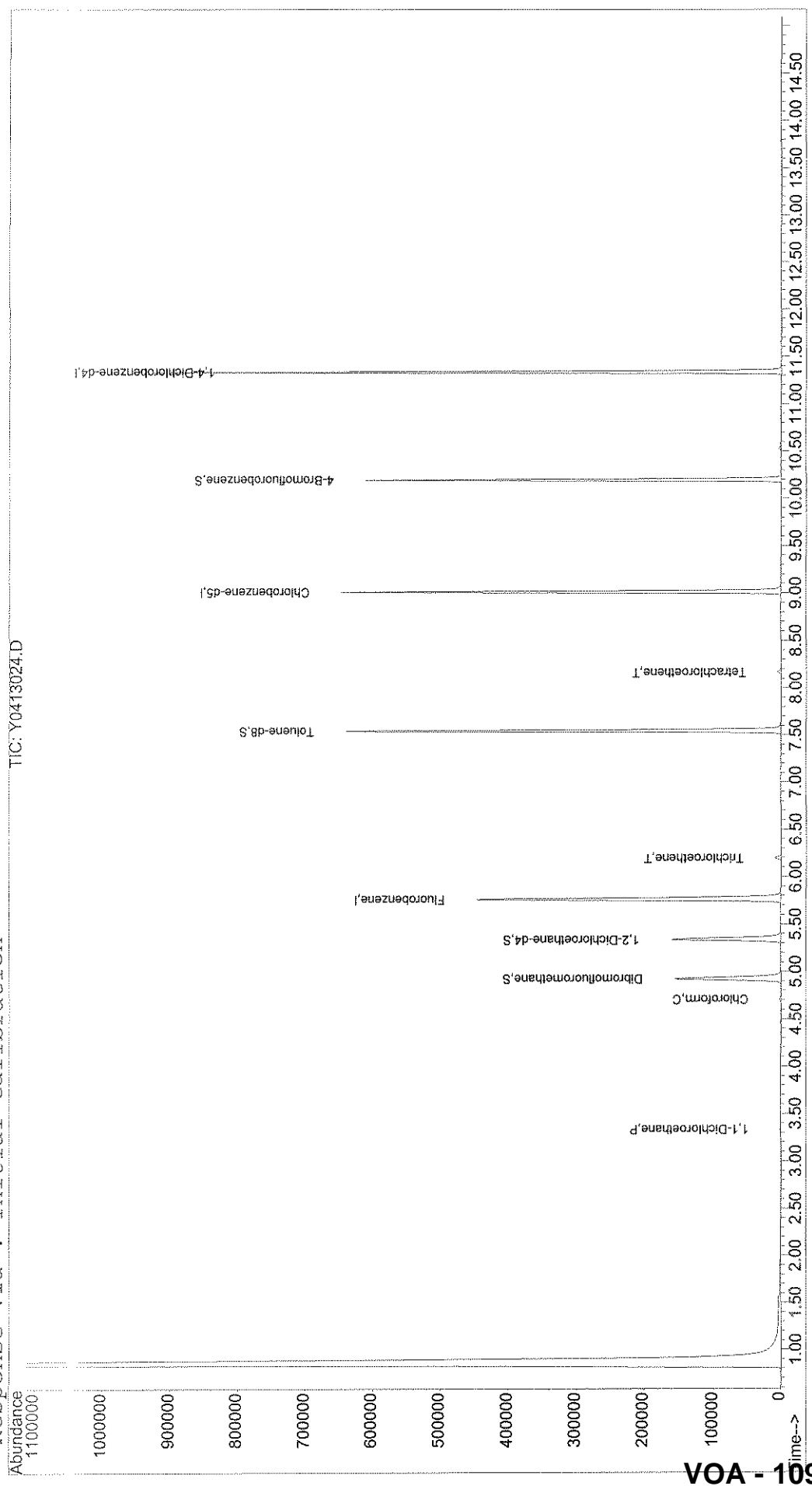
Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413024.D  
Acq On : 13 Apr 2007 15:48  
Sample : JPL30-012  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:14 2007

Vial: 38  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413024.D  
 Acq On : 13 Apr 2007 15:48  
 Sample : JPL30-012  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:14 2007

Vial: 38  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	408547	50.00	ug/l	0.00 120.25%
50) Chlorobenzene-d5	9.01	82	165232	50.00	ug/l	0.00 102.52%
70) 1,4-Dichlorobenzene-d4	11.34	152	202556	50.00	ug/l	0.00 89.88%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	112789	45.63	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	91.26%		
37) 1,2-Dichloroethane-d4	5.34	65	125837	47.99	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	95.98%		
51) Toluene-d8	7.54	98	371612	56.67	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	113.34%		
72) 4-Bromofluorobenzene	10.20	95	158385	54.12	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.33	63	1616	0.28	ug/l	72
23) Vinyl acetate	0.00	43	0	N.D.		

*JMT 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0413024.D 8260B.M Mon Apr 16 09:14:56 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413024.D  
 Acq On : 13 Apr 2007 15:48  
 Sample : JPL30-012  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:14 2007

Vial: 38  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	4.20	96	524	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.70	83	2387	0.45 ug/l ✓	95
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	0.00	78	0	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	6.20	130	3334	1.16 ug/l ✓	97
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	6.79	83	403	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	0.00	92	0	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	8.16	166	1419	0.62 ug/l ✓	93
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	8.43	129	61	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	0.00	91	0	N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.17	91	60	N.D.	
65) m,p-Xylene	9.29	106	133	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D.	

*LNH 4/16/07*

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413024.D  
 Acq On : 13 Apr 2007 15:48  
 Sample : JPL30-012  
 Misc : 5mL+IS/SS #1

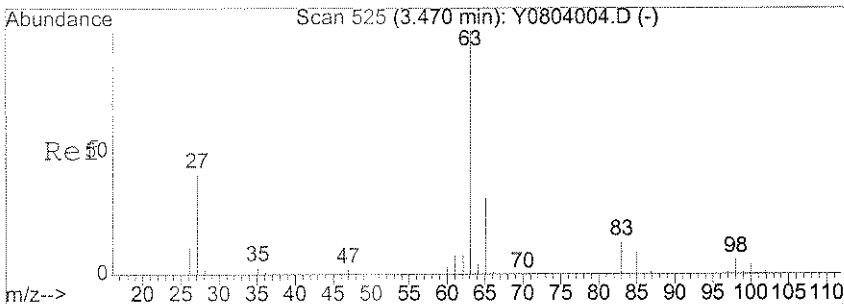
Vial: 38  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:14 2007

Quant Results File: 8260B.RES

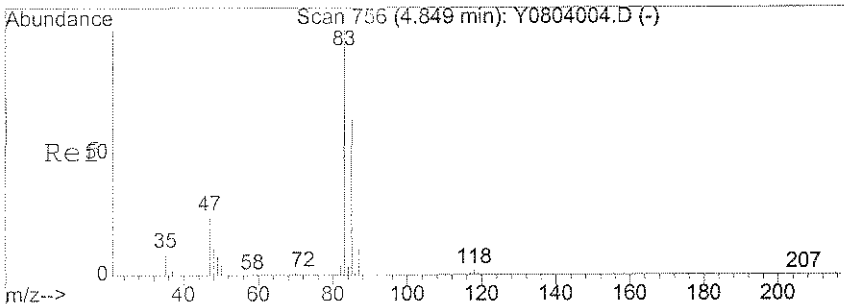
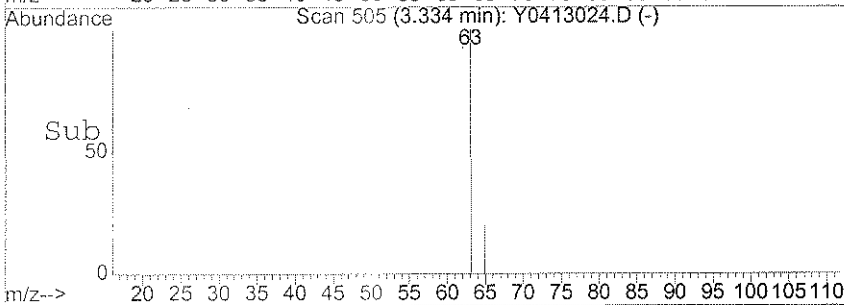
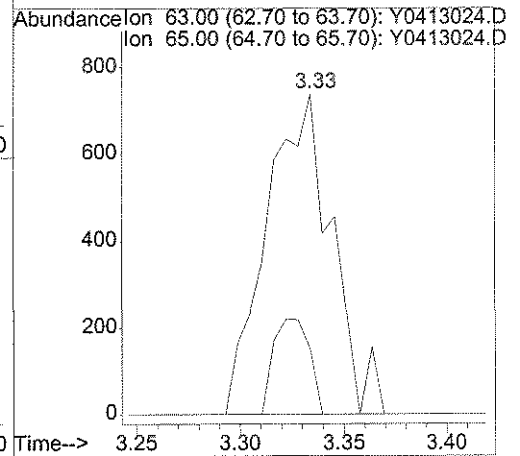
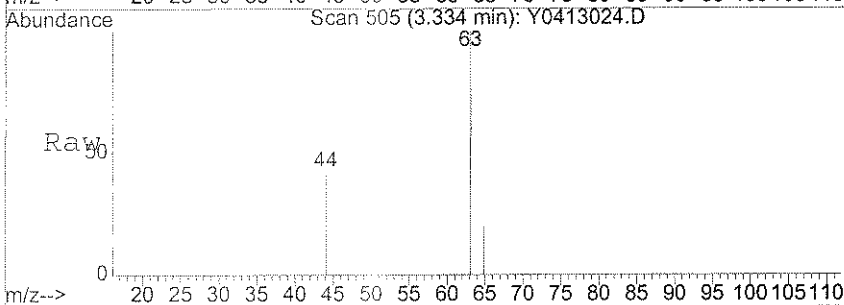
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	118		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	700		N.D.	
84) 4-Isopropyltoluene	11.33	119	204		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	285		N.D.	
86) 1,2-Dichlorobenzene	11.71	146	330		N.D.	
87) n-Butylbenzene	11.74	91	71		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	149		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	392		N.D.	



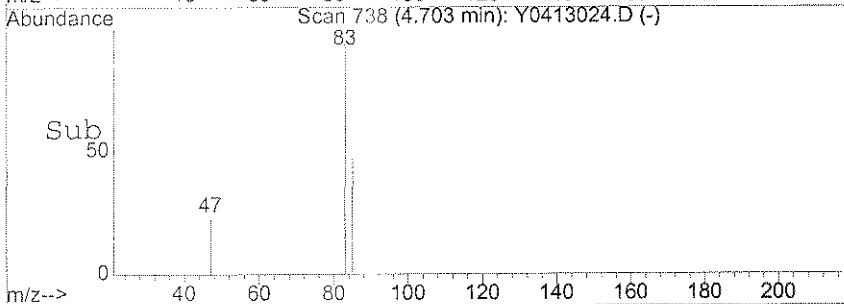
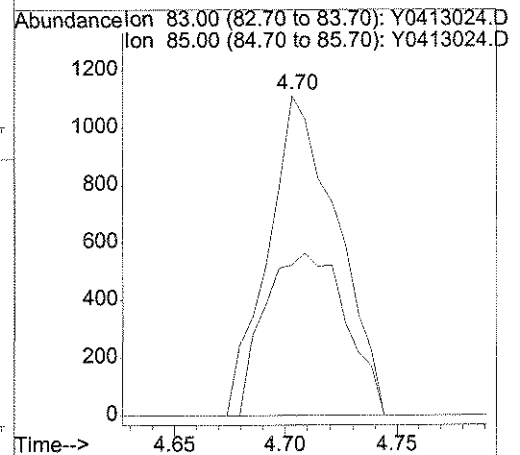
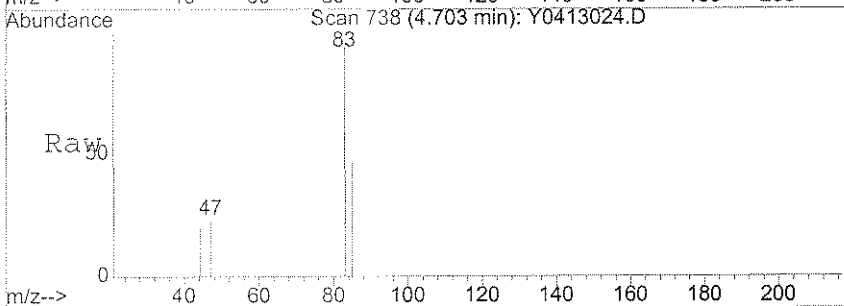
#22  
 1,1-Dichloroethane  
 Concen: 0.28 ug/l  
 RT: 3.33 min Scan# 505  
 Delta R.T. 0.01 min  
 Lab File: Y0413024.D  
 Acq: 13 Apr 2007 15:48

Tgt Ion: 63 Resp: 1616  
 Ion Ratio Lower Upper  
 63 100  
 65 16.7 12.3 52.3

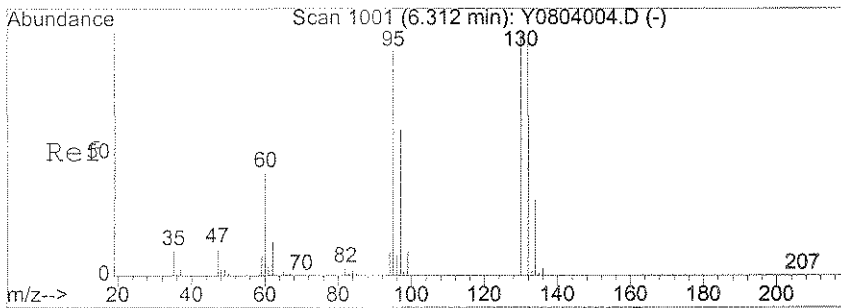


#31  
 Chloroform  
 Concen: 0.45 ug/l  
 RT: 4.70 min Scan# 738  
 Delta R.T. -0.00 min  
 Lab File: Y0413024.D  
 Acq: 13 Apr 2007 15:48

Tgt Ion: 83 Resp: 2387  
 Ion Ratio Lower Upper  
 83 100  
 85 59.2 43.3 83.3

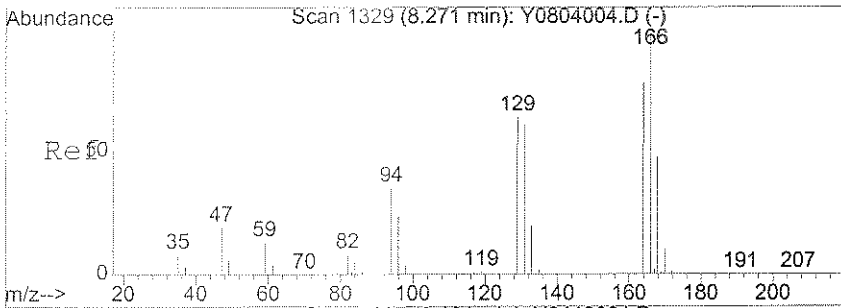
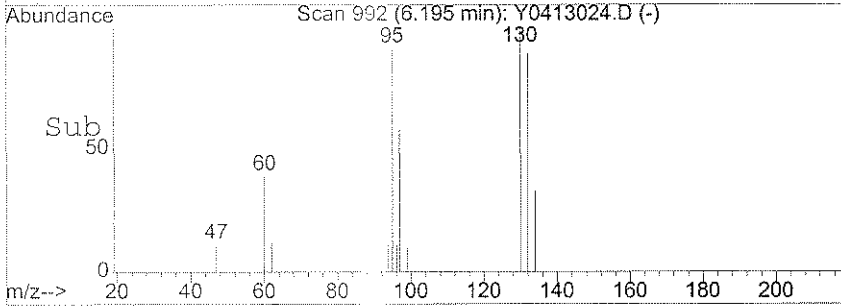
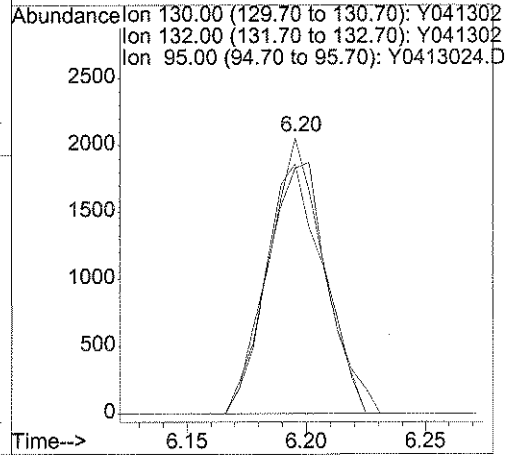
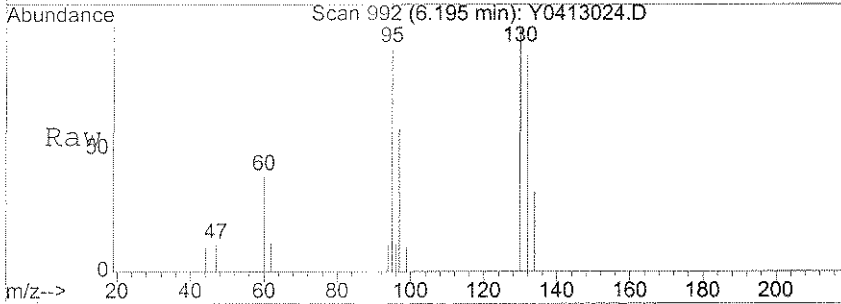






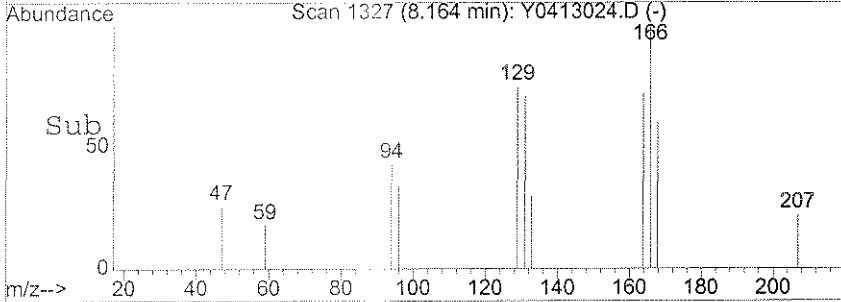
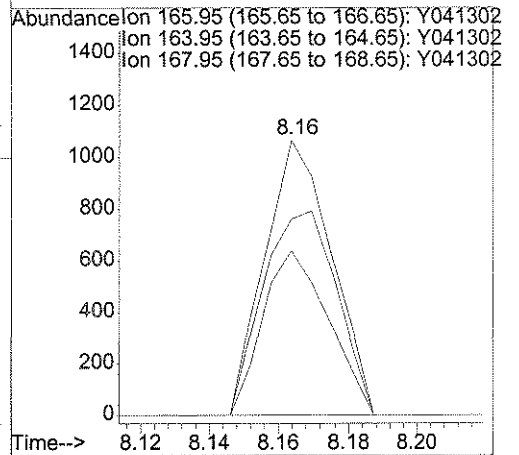
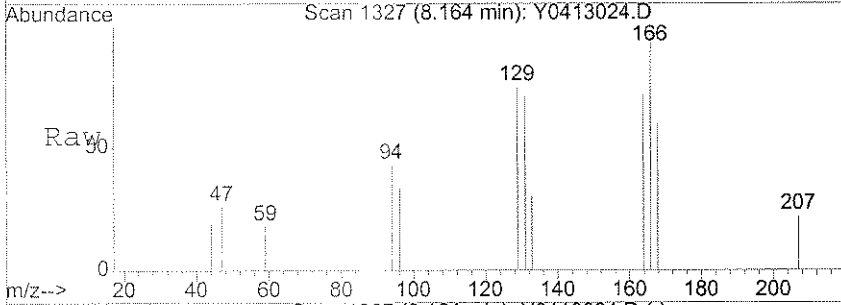
#41  
 Trichloroethene  
 Concen: 1.16 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0413024.D  
 Acq: 13 Apr 2007 15:48

Tgt Ion	Resp	Lower	Upper
130	100		
132	96.2	75.0	115.0
95	93.4	69.4	109.4



#56  
 Tetrachloroethene  
 Concen: 0.62 ug/l  
 RT: 8.16 min Scan# 1327  
 Delta R.T. -0.00 min  
 Lab File: Y0413024.D  
 Acq: 13 Apr 2007 15:48

Tgt Ion	Resp	Lower	Upper
166	100		
164	81.8	63.3	94.9
168	58.8	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 16:13

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 16:13

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-19-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 16:13

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

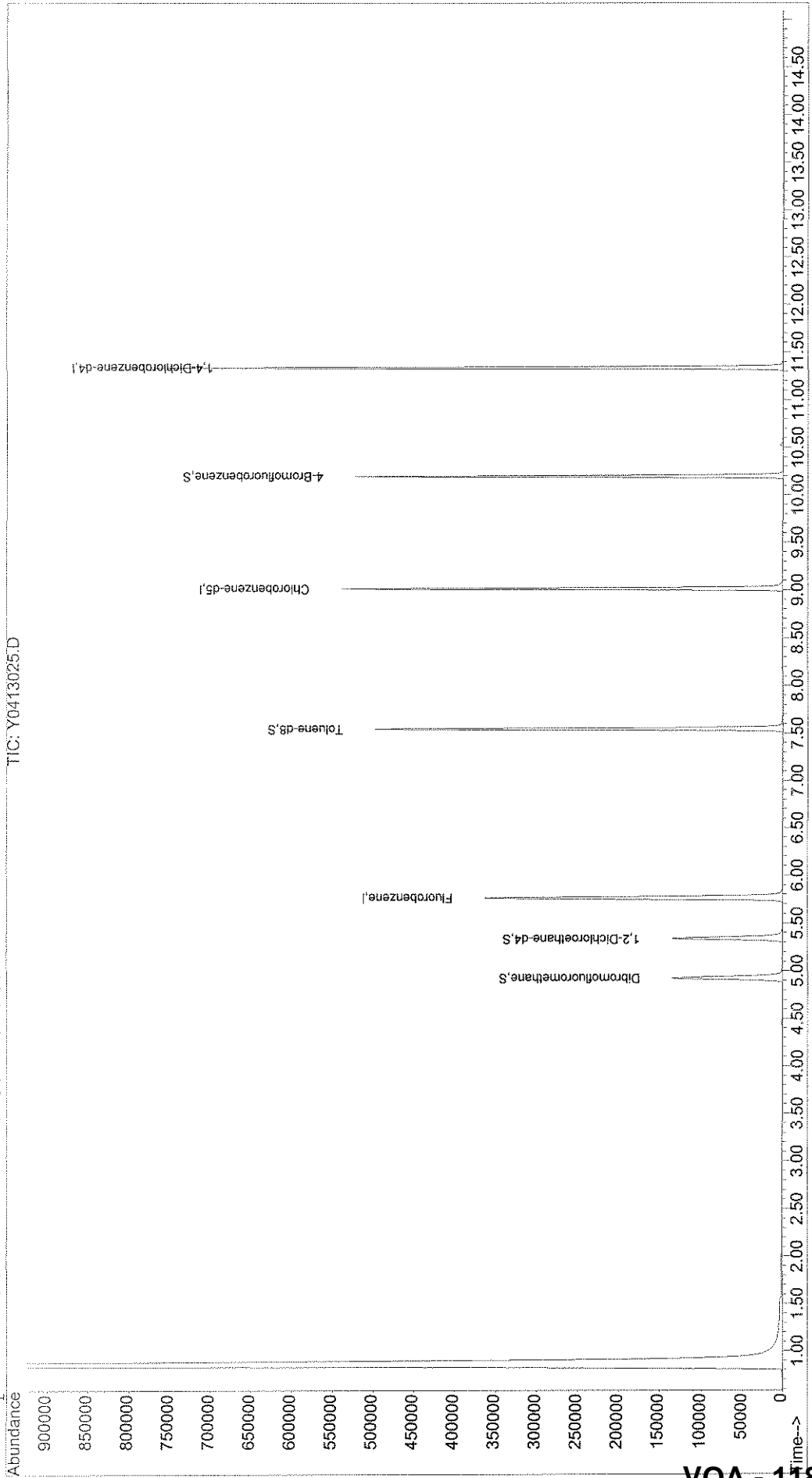
Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413025.D  
Acq On : 13 Apr 2007 16:13  
Sample : JPL30-013  
Misc : 5mL+IS/SS #2  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:15 2007

Vial: 39  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413025.D  
 Acq On : 13 Apr 2007 16:13  
 Sample : JPL30-013  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:15 2007

Vial: 39  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	5.77	96	334011	50.00	ug/l	0.00 98.32%
50) Chlorobenzene-d5	9.01	82	133677	50.00	ug/l	0.00 82.94%
70) 1,4-Dichlorobenzene-d4	11.34	152	170155	50.00	ug/l	0.00 75.51%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	98692	48.84	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	97.68%
37) 1,2-Dichloroethane-d4	5.34	65	106329	49.60	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.20%
51) Toluene-d8	7.55	98	294797	55.57	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	111.14%
72) 4-Bromofluorobenzene	10.20	95	132618	53.95	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	2.87	73	196		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413025.D 8260B.M Mon Apr 16 09:15:34 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413025.D  
 Acq On : 13 Apr 2007 16:13  
 Sample : JPL30-013  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:15 2007

Vial: 39  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	445		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	134		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413025.D 8260B.M Mon Apr 16 09:15:35 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413025.D  
 Acq On : 13 Apr 2007 16:13  
 Sample : JPL30-013  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:15 2007

Vial: 39  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.19	105	57		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	147		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413025.D 8260B.M Mon Apr 16 09:15:35 2007



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-4-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-014  
 Lab File ID: Y0413026.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 16:37  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-4-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-014  
 Lab File ID: Y0413026.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 16:37  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	0.50		U
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	0.50		U
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-4-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 16:37

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

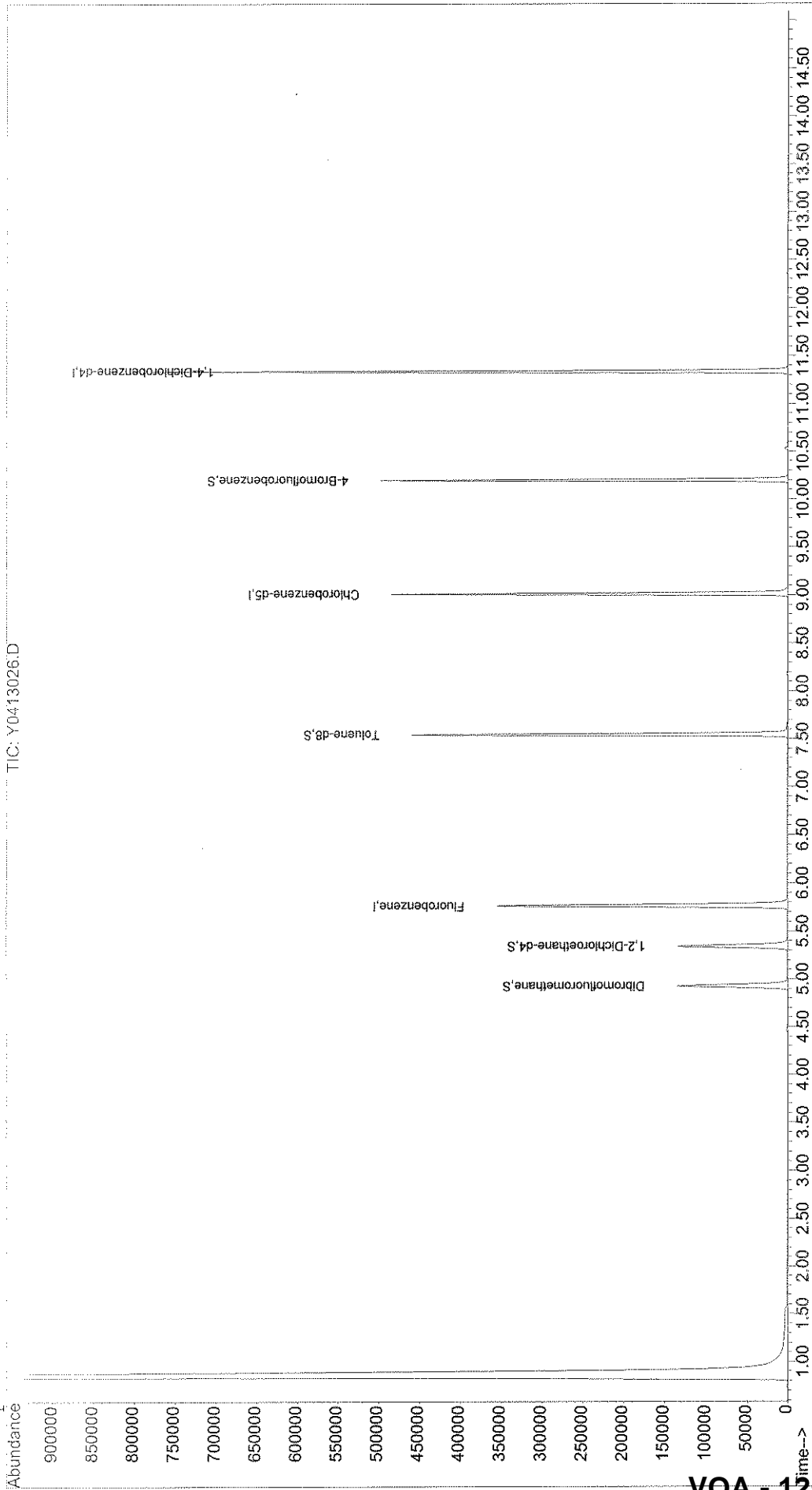
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413026.D  
Acq On : 13 Apr 2007 16:37 Vial: 40  
Sample : JPL30-014 Operator: LNH  
Misc : 5mL+IS/SS #3 Inst : Yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 16 9:16 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413026.D  
 Acq On : 13 Apr 2007 16:37  
 Sample : JPL30-014  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Vial: 40  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	5.77	96	325341	50.00	ug/l	0.00 95.76%
50) Chlorobenzene-d5	9.01	82	124349	50.00	ug/l	0.00 77.15%
70) 1,4-Dichlorobenzene-d4	11.34	152	167051	50.00	ug/l	0.00 74.13%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	96479	49.01	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	98.02%
37) 1,2-Dichloroethane-d4	5.34	65	103479	49.55	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.10%
51) Toluene-d8	7.55	98	274220	55.57	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	111.14%
72) 4-Bromofluorobenzene	10.20	95	125867	52.15	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	2.89	73	316		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413026.D 8260B.M Mon Apr 16 09:16:12 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413026.D  
 Acq On : 13 Apr 2007 16:37  
 Sample : JPL30-014  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Vial: 40  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	421		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	77		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413026.D 8260B.M Mon Apr 16 09:16:13 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413026.D  
 Acq On : 13 Apr 2007 16:37  
 Sample : JPL30-014  
 Misc : 5mL+IS/SS #3

Vial: 40  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	64		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	190		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-015  
 Lab File ID: Y0413011.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 10:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-015  
 Lab File ID: Y0413011.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 10:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-015  
 Lab File ID: Y0413011.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 10:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

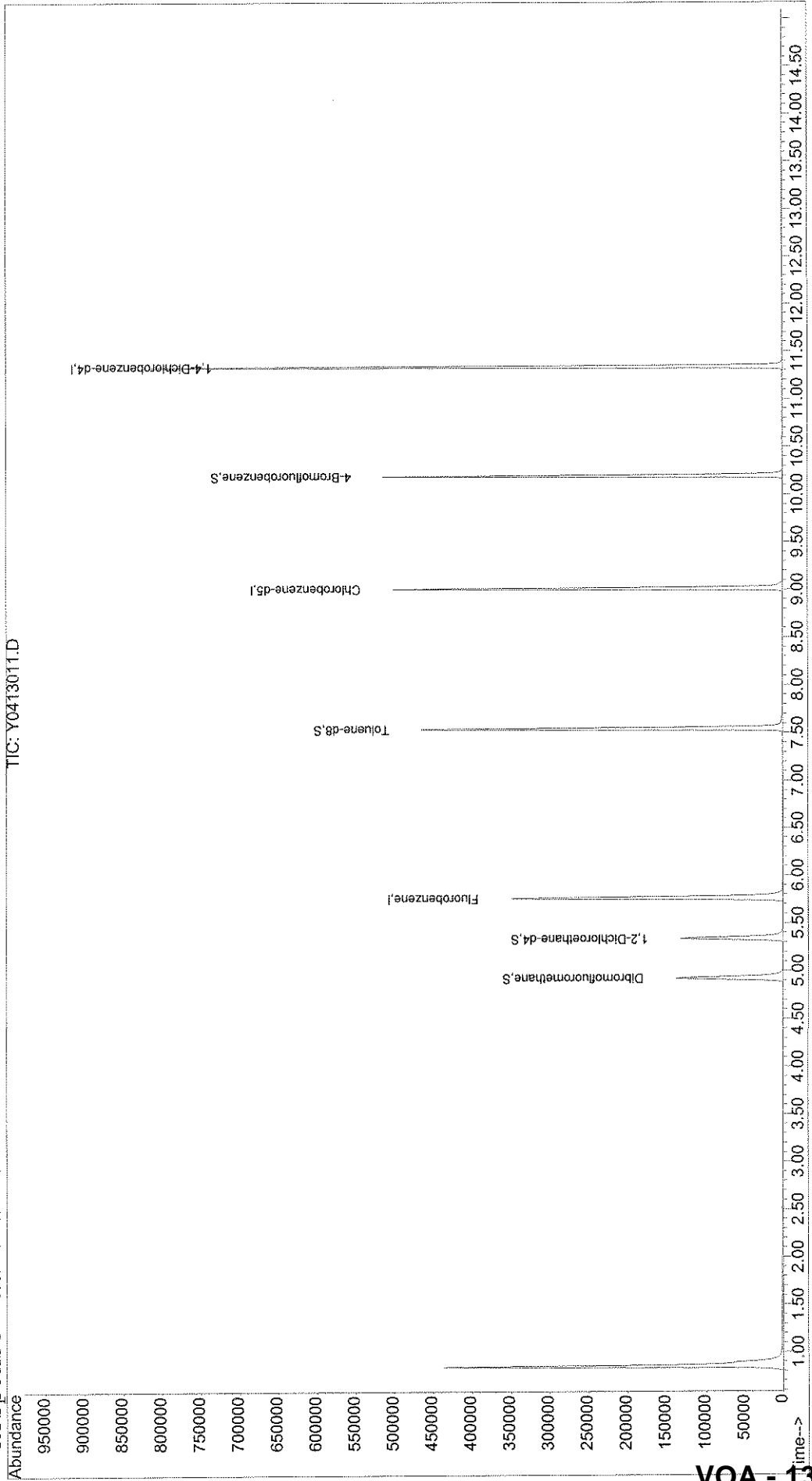
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413011.D Vial: 26  
Acq On : 13 Apr 2007 10:22 Operator: LNH  
Sample : JPL30-015 TB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:03 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413011.D  
 Acq On : 13 Apr 2007 10:22  
 Sample : JPL30-015 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:03 2007

Vial: 26  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	327892	50.00	ug/l	0.00 96.51%
50) Chlorobenzene-d5	9.01	82	129000	50.00	ug/l	0.00 80.04%
70) 1,4-Dichlorobenzene-d4	11.34	152	179922	50.00	ug/l	0.00 79.84%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	100144	50.48	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.96%
37) 1,2-Dichloroethane-d4	5.34	65	104266	49.54	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.08%
51) Toluene-d8	7.55	98	289068	56.47	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.94%
72) 4-Bromofluorobenzene	10.20	95	132810	51.09	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	88	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	2.51	84	476	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413011.D 8260B.M Mon Apr 16 09:03:24 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413011.D  
 Acq On : 13 Apr 2007 10:22  
 Sample : JPL30-015 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:03 2007

Vial: 26  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	398		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	54		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413011.D 8260B.M Mon Apr 16 09:03:24 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413011.D  
 Acq On : 13 Apr 2007 10:22  
 Sample : JPL30-015 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:03 2007

Vial: 26  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	117		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	252		N.D.	
78) 4-Chlorotoluene	10.65	91	55		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.02	105	123		N.D.	
82) sec-butylbenzene	11.18	105	59		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	53		N.D.	
84) 4-Isopropyltoluene	11.34	119	379		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	147		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	357		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.32	180	56		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-016

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413027.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 17:01

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-016  
 Lab File ID: Y0413027.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 17:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	0.50		U
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	0.50		U
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-016  
 Lab File ID: Y0413027.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 17:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

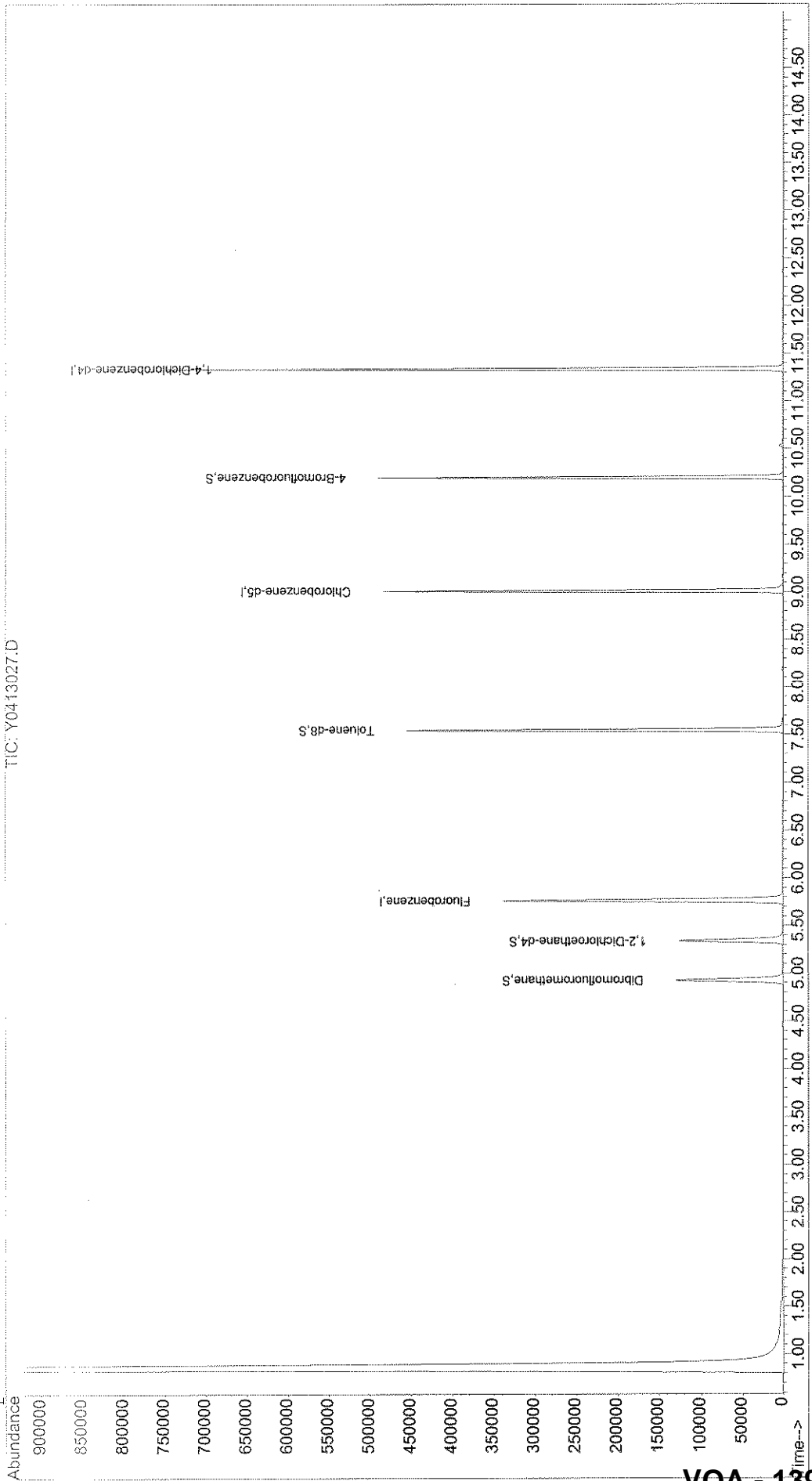
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	Q
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413027.D  
Acq On : 13 Apr 2007 17:01  
Sample : JPL30-016  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:16 2007  
Vial: 41  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413027.D  
 Acq On : 13 Apr 2007 17:01  
 Sample : JPL30-016  
 Misc : 5mL+IS/SS #1

Vial: 41  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	309977	50.00	ug/l	0.00 91.24%
50) Chlorobenzene-d5	9.01	82	124270	50.00	ug/l	0.00 77.10%
70) 1,4-Dichlorobenzene-d4	11.34	152	169026	50.00	ug/l	0.00 75.01%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	96072	51.23	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	102.46%	
37) 1,2-Dichloroethane-d4	5.34	65	101572	51.05	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	102.10%	
51) Toluene-d8	7.55	98	271904	55.13	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	110.26%	
72) 4-Bromofluorobenzene	10.20	95	125850	51.53	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413027.D 8260B.M Mon Apr 16 09:17:03 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413027.D  
 Acq On : 13 Apr 2007 17:01  
 Sample : JPL30-016  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Vial: 41  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	358		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	959		N.D.	
65) m,p-Xylene	9.17	106	185		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	302		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413027.D 8260B.M Mon Apr 16 09:17:04 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413027.D  
 Acq On : 13 Apr 2007 17:01  
 Sample : JPL30-016  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:16 2007

Vial: 41  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.21	105	73		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	154		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 17:26

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 17:26

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-017  
 Lab File ID: Y0413028.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 17:26  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

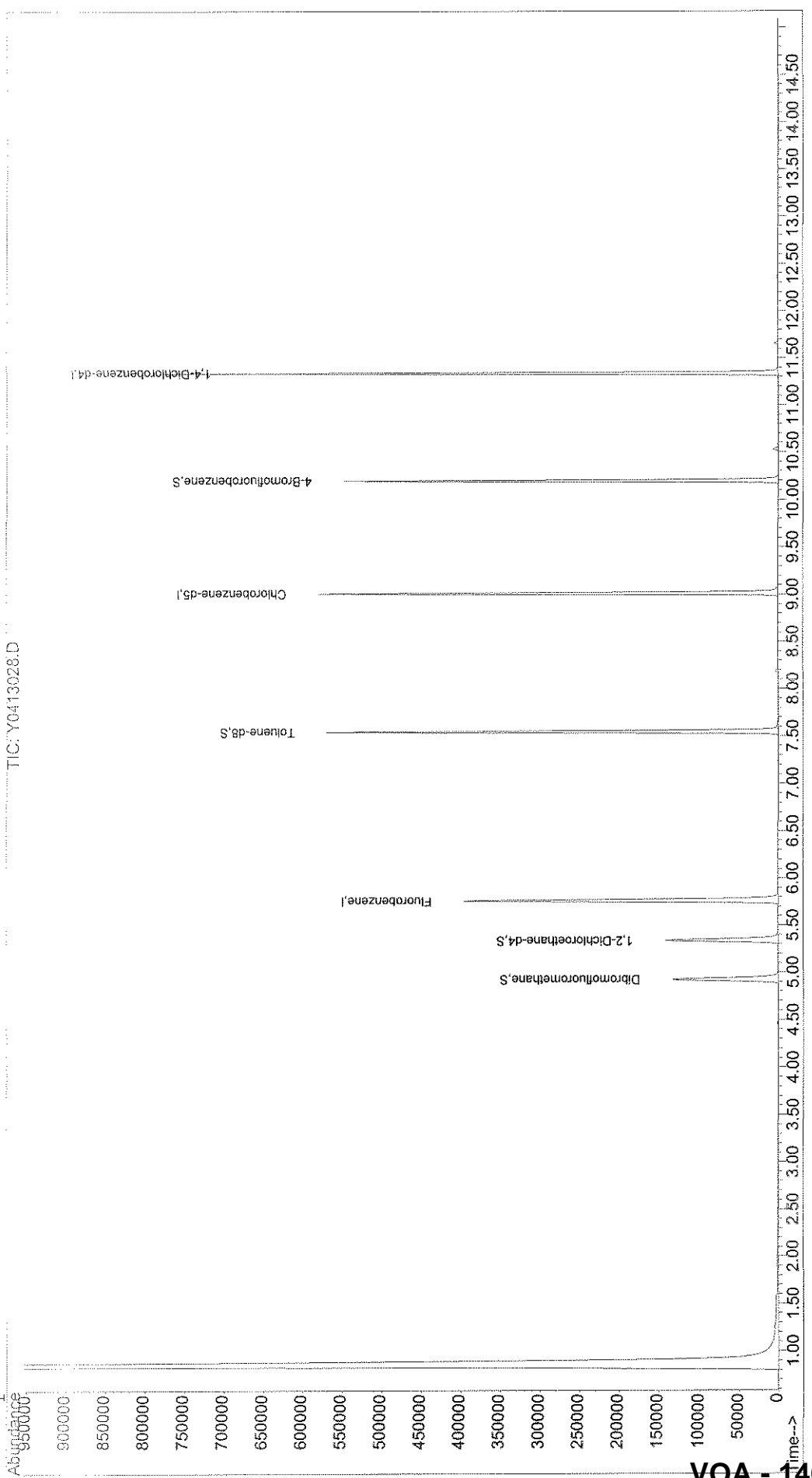
Comments:



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413028.D Vial: 42  
Acq On : 13 Apr 2007 17:26 Operator: LNH  
Sample : JPL30-017 Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rstein.p  
Quant Time: Apr 16 9:17 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413028.D  
 Acq On : 13 Apr 2007 17:26  
 Sample : JPL30-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:17 2007

Vial: 42  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	354186	50.00	ug/l	0.00 104.25%
50) Chlorobenzene-d5	9.01	82	145581	50.00	ug/l	0.00 90.32%
70) 1,4-Dichlorobenzene-d4	11.34	152	171857	50.00	ug/l	0.00 76.26%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	98569	46.00	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	92.00%
37) 1,2-Dichloroethane-d4	5.34	65	112119	49.32	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.64%
51) Toluene-d8	7.54	98	330727	57.25	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	114.50%
72) 4-Bromofluorobenzene	10.20	95	138414	55.75	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	2.81	53	169	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.32	63	291	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413028.D 8260B.M Mon Apr 16 09:17:46 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413028.D  
 Acq On : 13 Apr 2007 17:26  
 Sample : JPL30-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:17 2007

Vial: 42  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.39	78	121		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	57		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.17	166	116		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	507		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	740		N.D.	
65) m,p-Xylene	9.29	106	136		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	387		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413028.D 8260B.M Mon Apr 16 09:17:47 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413028.D  
 Acq On : 13 Apr 2007 17:26  
 Sample : JPL30-017  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:17 2007

Vial: 42  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	55		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	131		N.D.	
84) 4-Isopropyltoluene	11.34	119	129		N.D.	
85) 1,4-Dichlorobenzene	11.27	146	131		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413029.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 17:50

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.51	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	1.3	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.82	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-018  
 Lab File ID: Y0413029.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 17:50  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-3-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-018  
 Lab File ID: Y0413029.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 17:50  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

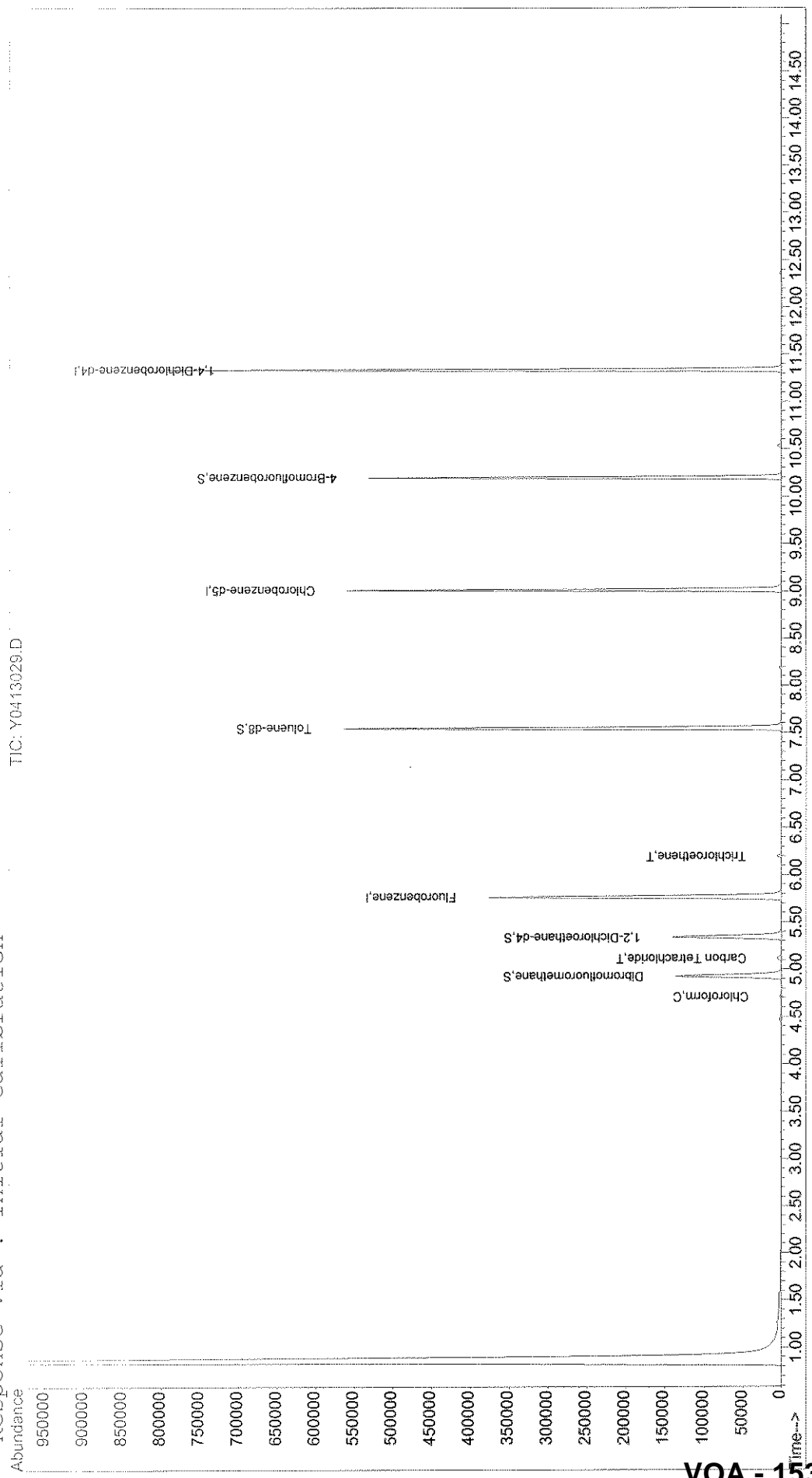
Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413029.D  
Acq On : 13 Apr 2007 17:50  
Sample : JPL30-018  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:18 2007

Vial: 43  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration





Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413029.D  
 Acq On : 13 Apr 2007 17:50  
 Sample : JPL30-018  
 Misc : 5mL+IS/SS #3

Vial: 43  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:18 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	345786	50.00	ug/l	0.00 101.78%
50) Chlorobenzene-d5	9.01	82	143555	50.00	ug/l	0.00 89.07%
70) 1,4-Dichlorobenzene-d4	11.34	152	175387	50.00	ug/l	0.00 77.83%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	97391	46.55	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	93.10%
37) 1,2-Dichloroethane-d4	5.34	65	109215	49.21	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.42%
51) Toluene-d8	7.55	98	332547	58.37	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	116.74%
72) 4-Bromofluorobenzene	10.20	95	136244	53.77	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Endomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	0.00	76	0	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0413029.D 8260B.M Mon Apr 16 09:18:36 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413029.D  
 Acq On : 13 Apr 2007 17:50  
 Sample : JPL30-018  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:18 2007

Vial: 43  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	0.00	43	0	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	4.71	83	2266	0.51	ug/l ✓	98
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	5.12	117	4028	1.29	ug/l ✓	95
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	5.39	78	55	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.		
41) Trichloroethene	6.20	130	2003	0.82	ug/l ✓	88
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	0.00	41	0	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	7.61	92	57	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.		
56) Tetrachloroethene	8.16	166	55	N.D.		
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.		
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	0.00	112	0	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.28	91	57	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.		
66) o-xylene	0.00	106	0	N.D.		
67) Styrene	0.00	104	0	N.D.		
68) Bromoform	0.00	173	0	N.D.		

*LNH 4/16/07*

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413029.D  
 Acq On : 13 Apr 2007 17:50  
 Sample : JPL30-018  
 Misc : 5mL+IS/SS #3

Vial: 43  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

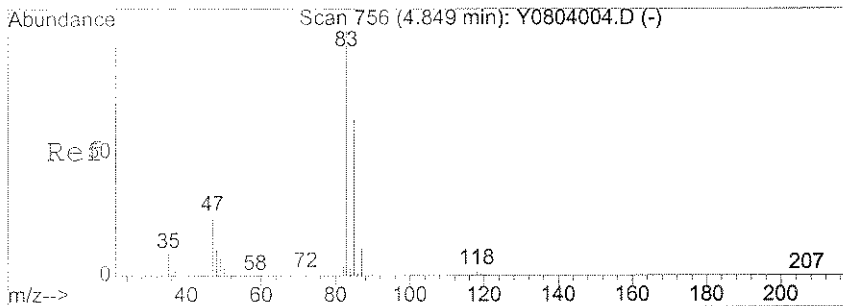
MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:18 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

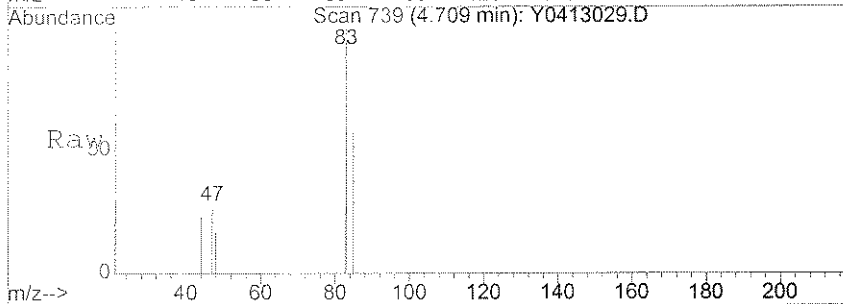
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	69		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.33	119	128		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413029.D 8260B.M Mon Apr 16 09:18:36 2007

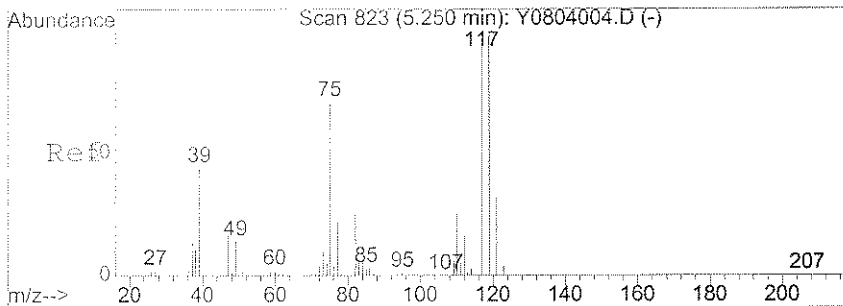
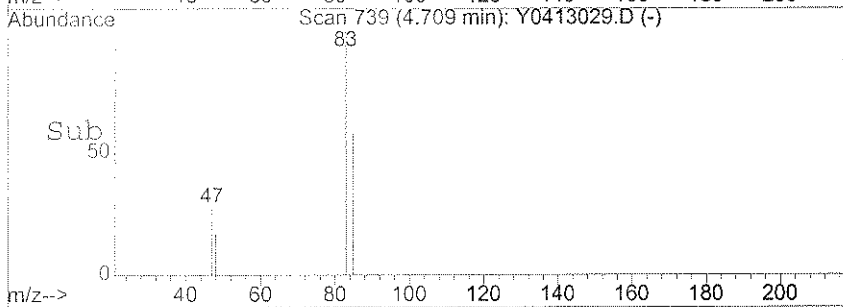
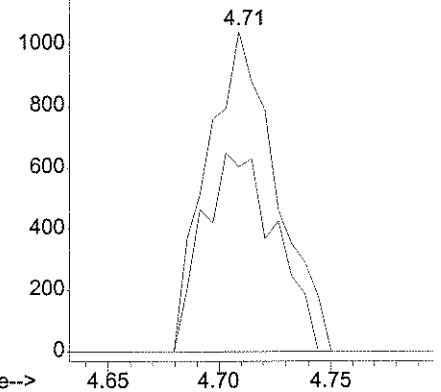


#31  
 Chloroform  
 Concen: 0.51 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0413029.D  
 Acq: 13 Apr 2007 17:50

Tgt Ion:	83	Resp:	2266
Ion Ratio	Lower	Upper	
83	100		
85	65.2	43.3	83.3

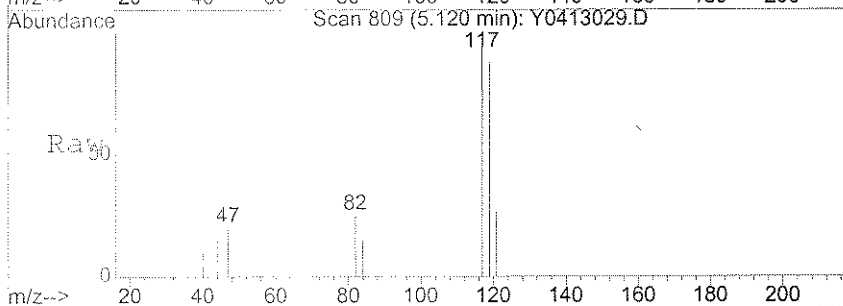


Abundance Ion 83.00 (82.70 to 83.70): Y0413029.D  
 1200 Ion 85.00 (84.70 to 85.70): Y0413029.D

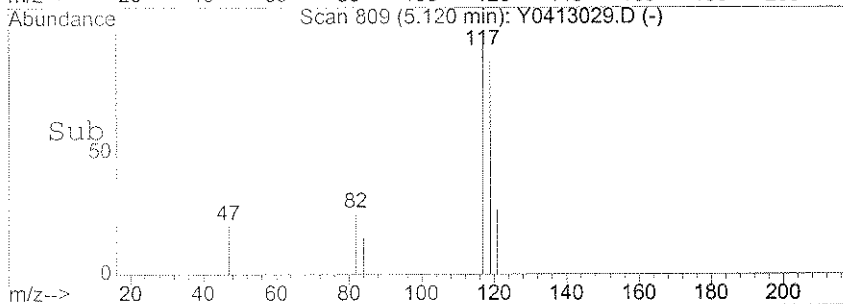
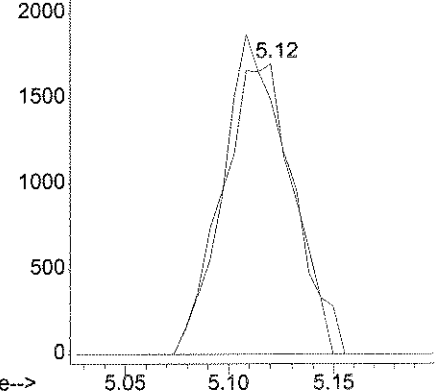


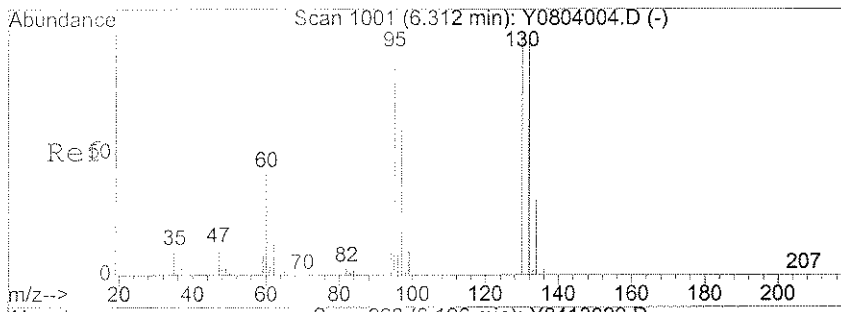
#35  
 Carbon Tetrachloride  
 Concen: 1.29 ug/l  
 RT: 5.12 min Scan# 809  
 Delta R.T. 0.01 min  
 Lab File: Y0413029.D  
 Acq: 13 Apr 2007 17:50

Tgt Ion:	117	Resp:	4028
Ion Ratio	Lower	Upper	
117	100		
119	103.0	78.2	118.2



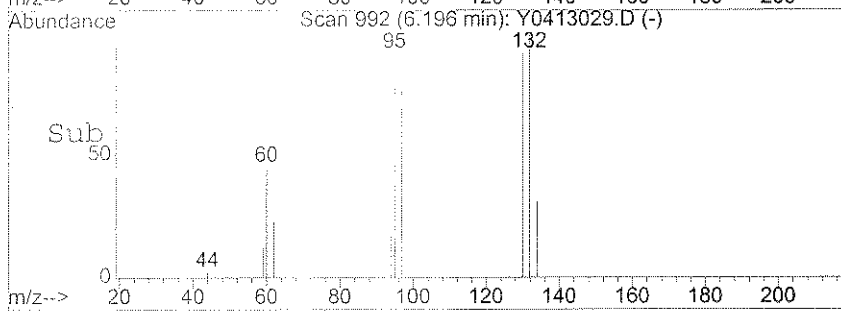
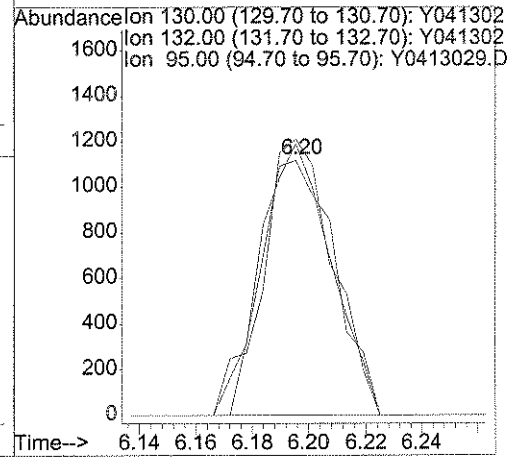
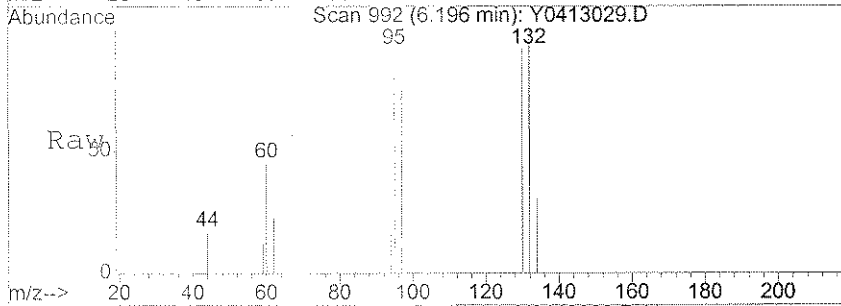
Abundance Ion 117.00 (116.70 to 117.70): Y041302  
 2000 Ion 119.00 (118.70 to 119.70): Y041302





#41  
 Trichloroethene  
 Concen: 0.82 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0413029.D  
 Acq: 13 Apr 2007 17:50

Tgt Ion	Resp	Lower	Upper
130	100		
132	104.0	75.0	115.0
95	103.8	69.4	109.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL30-019

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413012.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/02/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/13/2007 10:47

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-019  
 Lab File ID: Y0413012.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 10:47  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL30  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-019  
 Lab File ID: Y0413012.D  
 Date Collected: 04/02/2007  
 Date/Time Analyzed: 04/13/2007 10:47  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

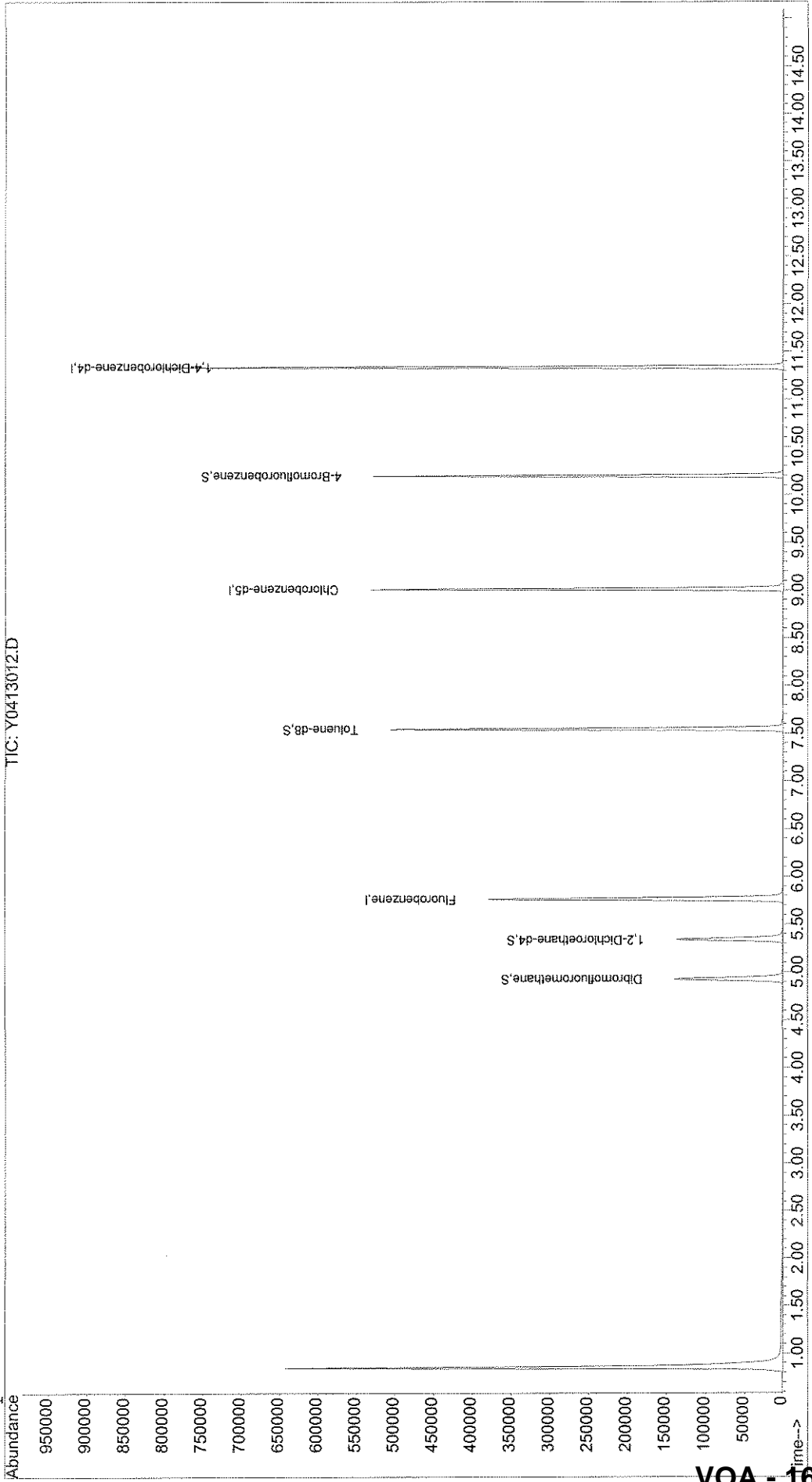
Comments:



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413012.D Vial: 27  
Acq On : 13 Apr 2007 10:47 Operator: LNH  
Sample : JPL30-019 EB Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 9:04 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413012.D  
 Acq On : 13 Apr 2007 10:47  
 Sample : JPL30-019 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:04 2007

Vial: 27  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	348620	50.00	ug/l	0.00 102.62%
50) Chlorobenzene-d5	9.01	82	133681	50.00	ug/l	0.00 82.94%
70) 1,4-Dichlorobenzene-d4	11.34	152	182353	50.00	ug/l	0.00 80.92%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101060	47.91	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	95.82%
37) 1,2-Dichloroethane-d4	5.34	65	108551	48.51	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	97.02%
51) Toluene-d8	7.54	98	299688	56.49	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.98%
72) 4-Bromofluorobenzene	10.20	95	136609	51.85	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	69	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413012.D  
 Acq On : 13 Apr 2007 10:47  
 Sample : JPL30-019 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:04 2007

Vial: 27  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	63		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.00	91	428		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	640		N.D.	
65) m,p-Xylene	9.30	106	414		N.D.	
66) o-xylene	9.68	106	53		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0413012.D 8260B.M Mon Apr 16 09:04:14 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041307\Y0413012.D  
 Acq On : 13 Apr 2007 10:47  
 Sample : JPL30-019 EB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 9:04 2007

Vial: 27  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

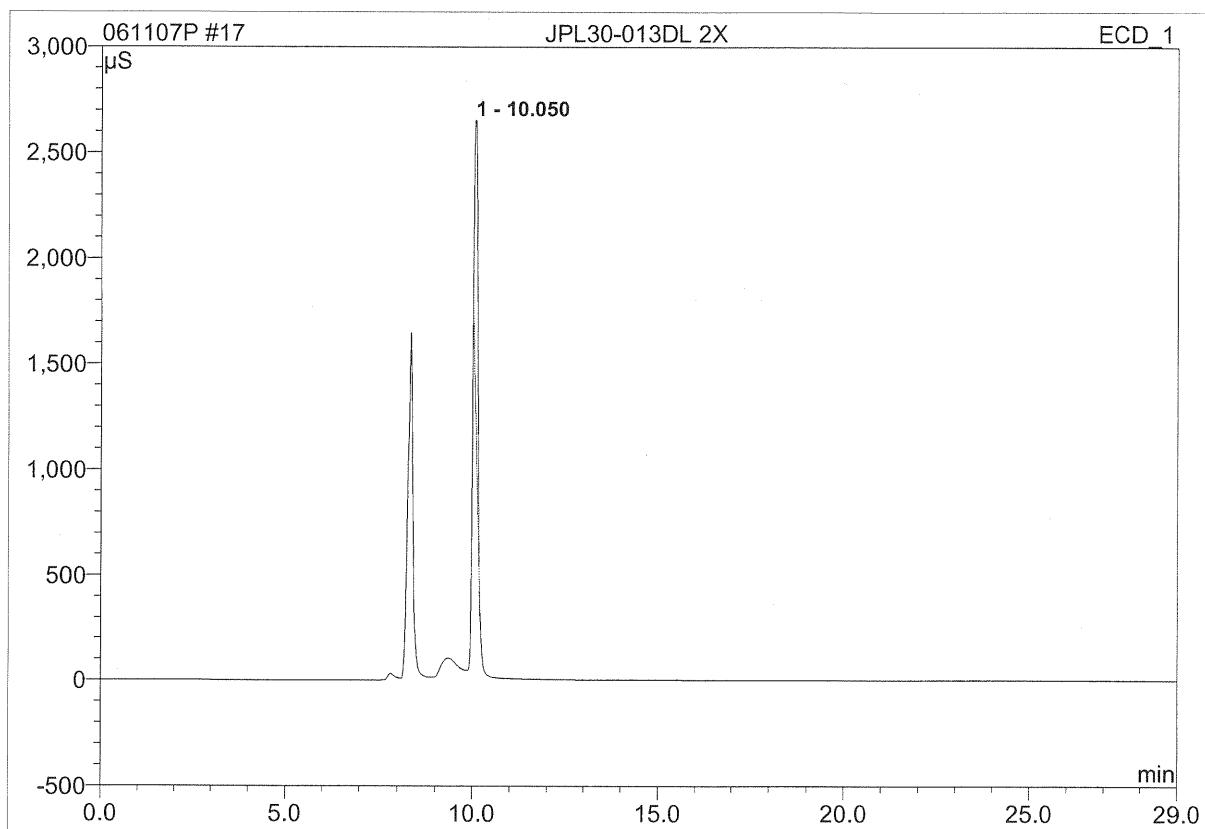
Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	248		N.D.	
78) 4-Chlorotoluene	10.64	91	111		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.02	105	125		N.D.	
82) sec-butylbenzene	11.18	105	115		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	407		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	365		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

**17 JPL30-013DL 2X****ICS-2500 DIONEX B**

Sample Name:	JPL30-013DL 2X	Injection Volume:	5.0
Vial Number:	17	Channel:	ECD_1
Sample Type:	unknown	Wavelength:	n.a.
Control Program:	perchlorate	Bandwidth:	n.a.
Quantif. Method:	perc	Dilution Factor:	1.0000
Recording Time:	6/11/2007 23:10	Sample Weight:	1.0000
Run Time (min):	29.00	Sample Amount:	1.0000

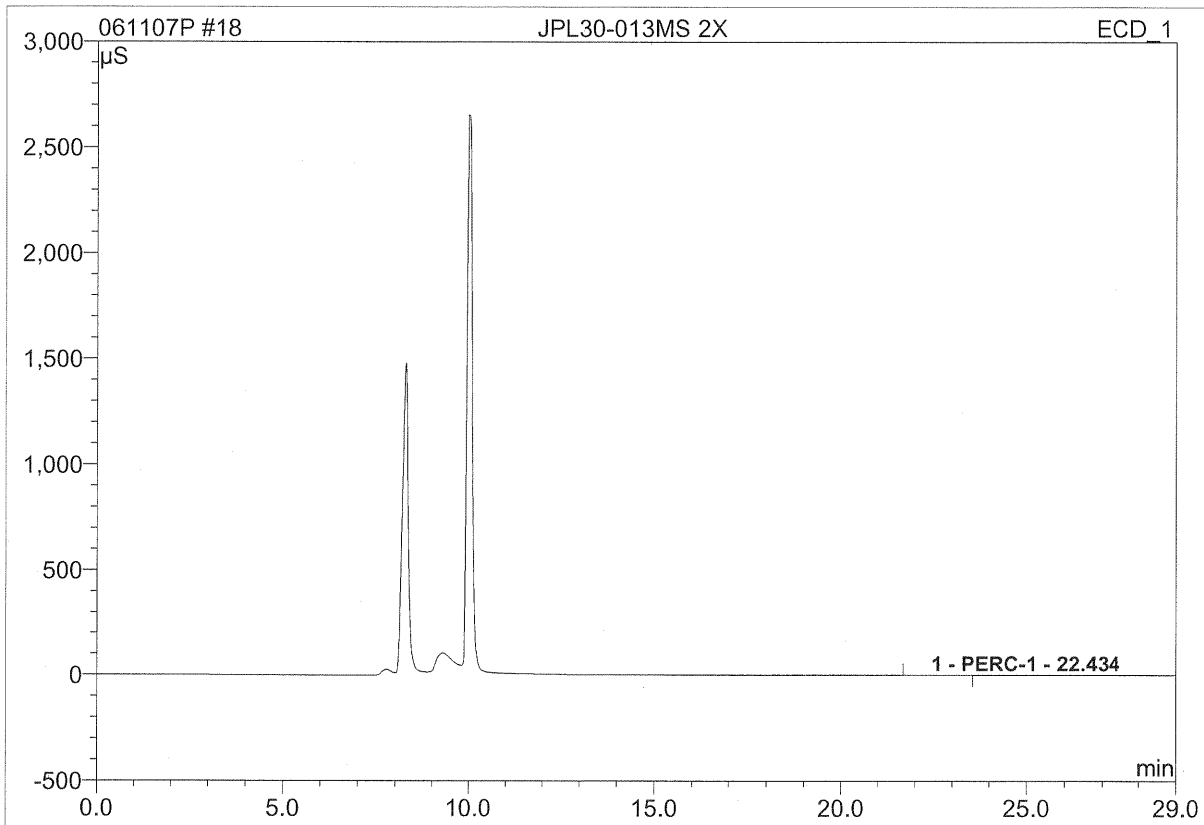


No.	Ret.Time min	Peak Name	Height μS	Area μS*min	Rel.Area %	Amount ppb	Type
1	10.05	n.a.	1204.128	149.899	100.00	n.a.	BMB
<b>Total:</b>			1204.128	149.899	100.00	0.000	

# 18 JPL30-013MS 2X

## ICS-2500 DIONEX B (250 uL IC-7-24-9)

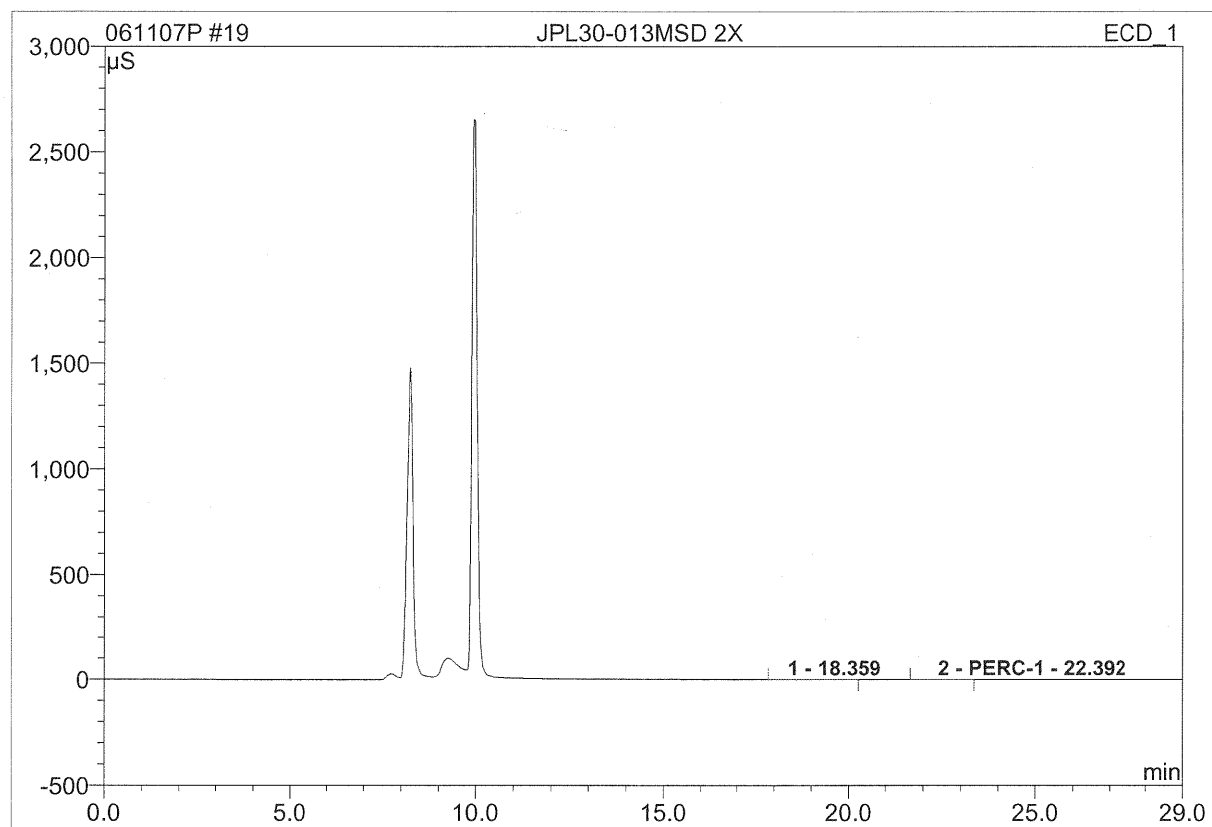
Sample Name:	JPL30-013MS 2X	Injection Volume:	5.0
Vial Number:	18	Channel:	ECD_1
Sample Type:	unknown	Wavelength:	n.a.
Control Program:	perchlorate	Bandwidth:	n.a.
Quantif. Method:	perc	Dilution Factor:	1.0000
Recording Time:	6/11/2007 23:41	Sample Weight:	1.0000
Run Time (min):	29.00	Sample Amount:	1.0000



No.	Ret.Time min	Peak Name	Height μS	Area μS*min	Rel.Area %	Amount ppb	Type
1	22.43	PERC-1	0.349	0.217	100.00	21.275	BMB
<b>Total:</b>			0.349	0.217	100.00	21.275	

**19 JPL30-013MSD 2X****ICS-2500 DIONEX B (250 uL IC-7-24-9)**

Sample Name:	JPL30-013MSD 2X	Injection Volume:	5.0
Vial Number:	19	Channel:	ECD_1
Sample Type:	unknown	Wavelength:	n.a.
Control Program:	perchlorate	Bandwidth:	n.a.
Quantif. Method:	perc	Dilution Factor:	1.0000
Recording Time:	6/12/2007 0:12	Sample Weight:	1.0000
Run Time (min):	29.00	Sample Amount:	1.0000

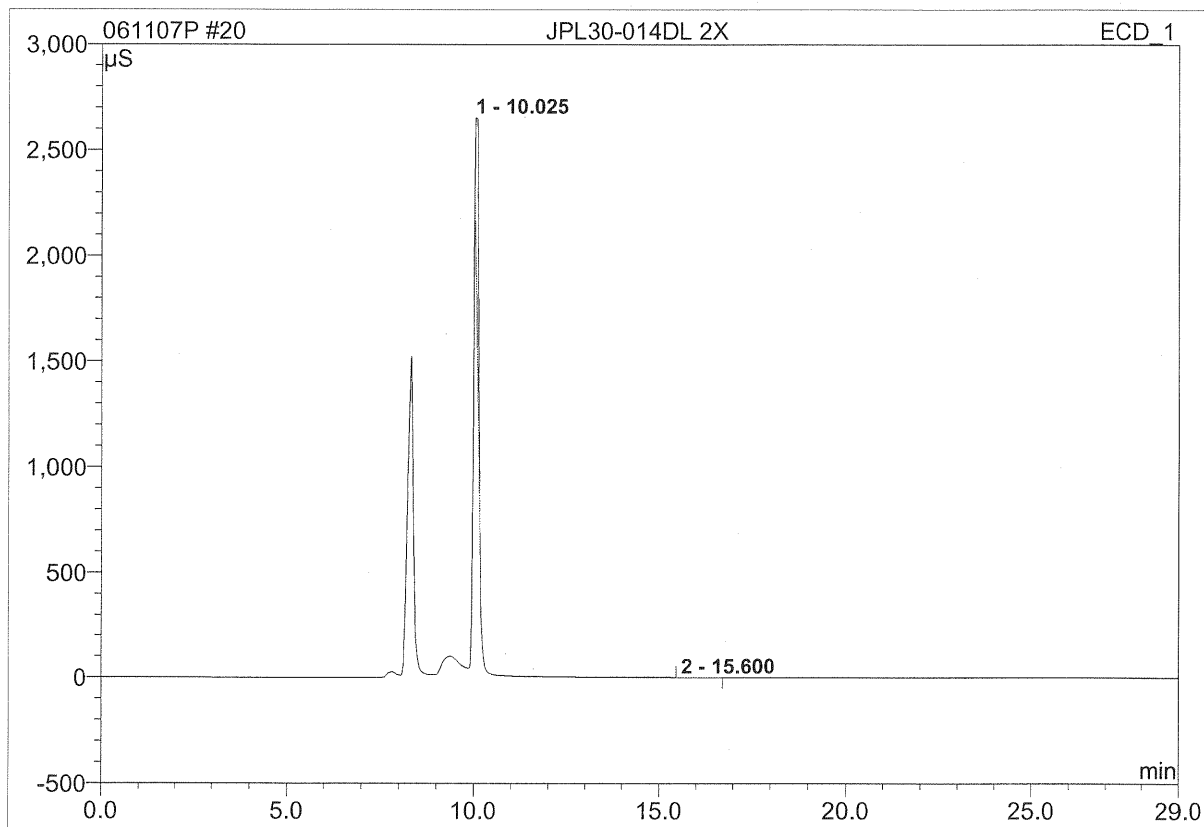


No.	Ret.Time min	Peak Name	Height $\mu\text{S}$	Area $\mu\text{S}\cdot\text{min}$	Rel.Area %	Amount ppb	Type
1	18.36	n.a.	0.086	0.105	32.16	n.a.	BMB
2	22.39	PERC-1	0.357	0.221	67.84	21.634	BMB
<b>Total:</b>			0.443	0.326	100.00	21.634	

## 20 JPL30-014DL 2X

### ICS-2500 DIONEX B

Sample Name:	JPL30-014DL 2X	Injection Volume:	5.0
Vial Number:	20	Channel:	ECD_1
Sample Type:	unknown	Wavelength:	n.a.
Control Program:	perchlorate	Bandwidth:	n.a.
Quantif. Method:	perc	Dilution Factor:	1.0000
Recording Time:	6/12/2007 0:44	Sample Weight:	1.0000
Run Time (min):	29.00	Sample Amount:	1.0000



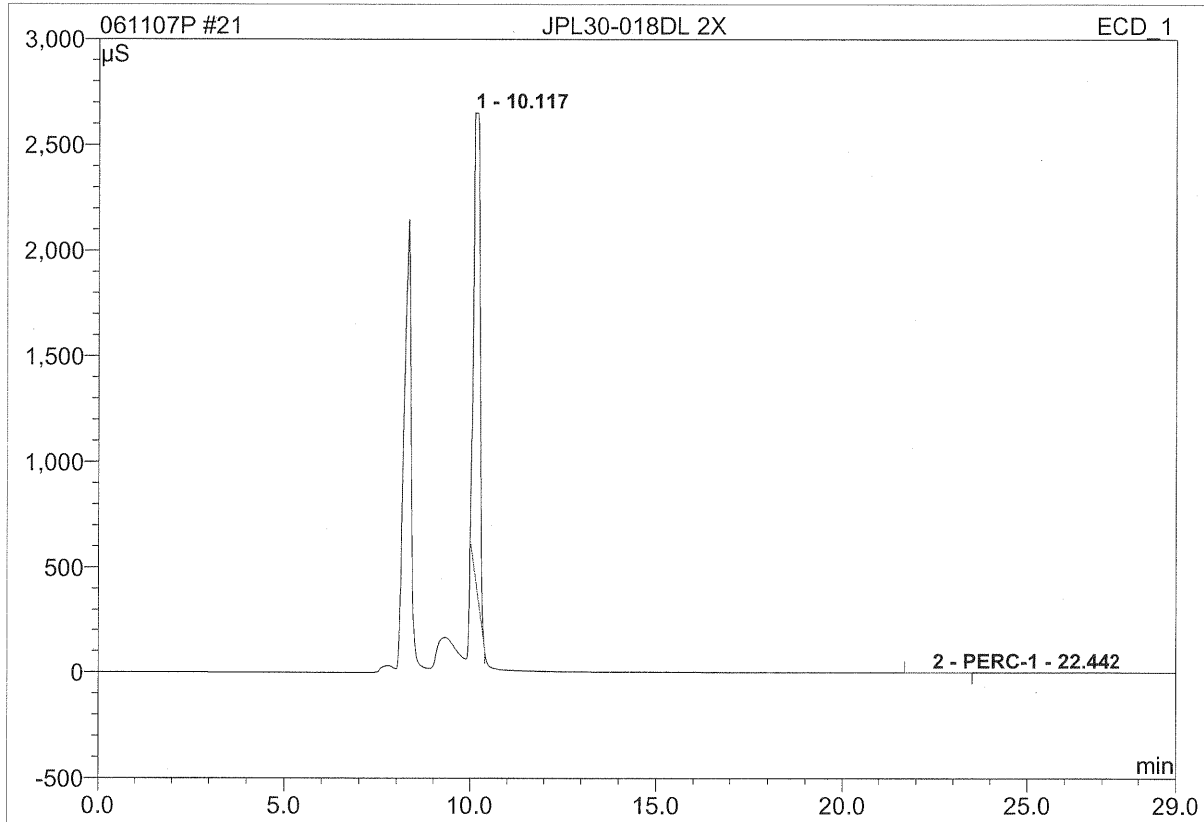
No.	Ret.Time min	Peak Name	Height µS	Area µS*min	Rel.Area %	Amount ppb	Type
1	10.03	n.a.	587.378	88.283	99.97	n.a.	BMB
2	15.60	n.a.	0.034	0.028	0.03	n.a.	BMB
<b>Total:</b>			587.412	88.311	100.00	0.000	



## 21 JPL30-018DL 2X

### ICS-2500 DIONEX B

Sample Name:	JPL30-018DL 2X	Injection Volume:	5.0
Vial Number:	21	Channel:	ECD_1
Sample Type:	unknown	Wavelength:	n.a.
Control Program:	perchlorate	Bandwidth:	n.a.
Quantif. Method:	perc	Dilution Factor:	1.0000
Recording Time:	6/12/2007 1:15	Sample Weight:	1.0000
Run Time (min):	29.00	Sample Amount:	1.0000



No.	Ret.Time min	Peak Name	Height µS	Area µS*min	Rel.Area %	Amount ppb	Type
1	10.12	n.a.	2180.857	452.228	99.94	n.a.	BMB
2	22.44	PERC-1	0.428	0.264	0.06	25.882	BMB
<b>Total:</b>			2181.285	452.492	100.00	25.882	

**TIC REPORTS**

SDG #JPL30

Volatiles Analysis

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-20-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-001  
 Lab File ID: Y0413014.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
10				
11				
12				
13				
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17				
18				
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23				
24				
25				
26				
27				
28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413014.D Vial: 28  
Acq On : 13 Apr 2007 11:36 Operator: LNH  
Sample : JPL30-001 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413014.D 8260B.M Mon Apr 16 09:06:00 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-20-4

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL30-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/31/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413015.D Vial: 29  
Acq On : 13 Apr 2007 12:01 Operator: LNH  
Sample : JPL30-002 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413015.D 8260B.M Mon Apr 16 09:07:12 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-20-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-003  
 Lab File ID: Y0413016.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413016.D Vial: 30  
Acq On : 13 Apr 2007 12:26 Operator: LNH  
Sample : JPL30-003 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413016.D 8260B.M Mon Apr 16 09:55:31 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-20-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-004  
 Lab File ID: Y0413017.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413017.D Vial: 31  
Acq On : 13 Apr 2007 12:50 Operator: LNH  
Sample : JPL30-004 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413017.D 8260B.M Mon Apr 16 09:09:37 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-20-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-005  
 Lab File ID: Y0413018.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413018.D Vial: 32  
Acq On : 13 Apr 2007 13:14 Operator: LNH  
Sample : JPL30-005 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413018.D 8260B.M Mon Apr 16 09:10:19 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-3-1Q07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL30-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 03/31/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413019.D Vial: 33  
Acq On : 13 Apr 2007 13:39 Operator: LNH  
Sample : JPL30-006 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413019.D 8260B.M Mon Apr 16 09:11:06 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-007  
 Lab File ID: Y0413009.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413009.D Vial: 24  
Acq On : 13 Apr 2007 9:33 Operator: LNH  
Sample : JPL30-007 EB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413009.D 8260B.M Mon Apr 16 09:02:35 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-4-3/30/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-008  
 Lab File ID: Y0413010.D  
 Date Collected: 03/31/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413010.D Vial: 25  
Acq On : 13 Apr 2007 9:58 Operator: LNH  
Sample : JPL30-008 TB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413010.D 8260B.M Mon Apr 16 09:02:25 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-19-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-009  
 Lab File ID: Y0413021.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413021.D Vial: 35  
Acq On : 13 Apr 2007 14:34 Operator: LNH  
Sample : JPL30-009 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413021.D 8260B.M Mon Apr 16 09:12:34 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-19-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-010  
 Lab File ID: Y0413022.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413022.D Vial: 36  
Acq On : 13 Apr 2007 14:59 Operator: LNH  
Sample : JPL30-010 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413022.D 8260B.M Mon Apr 16 09:13:22 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-19-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-011  
 Lab File ID: Y0413023.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413023.D Vial: 37  
Acq On : 13 Apr 2007 15:23 Operator: LNH  
Sample : JPL30-011 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413023.D 8260B.M Mon Apr 16 09:14:10 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-19-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-012  
 Lab File ID: Y0413024.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413024.D Vial: 38  
Acq On : 13 Apr 2007 15:48 Operator: LNH  
Sample : JPL30-012 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413024.D 8260B.M Mon Apr 16 09:15:01 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-19-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-013  
 Lab File ID: Y0413025.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413025.D Vial: 39  
Acq On : 13 Apr 2007 16:13 Operator: LNH  
Sample : JPL30-013 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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Y0413025.D 8260B.M Mon Apr 16 09:15:39 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-4-1Q07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-014  
 Lab File ID: Y0413026.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
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12				
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28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413026.D Vial: 40  
Acq On : 13 Apr 2007 16:37 Operator: LNH  
Sample : JPL30-014 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413026.D 8260B.M Mon Apr 16 09:16:16 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-015  
 Lab File ID: Y0413011.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
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28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413011.D Vial: 26  
Acq On : 13 Apr 2007 10:22 Operator: LNH  
Sample : JPL30-015 TB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413011.D 8260B.M Mon Apr 16 09:03:28 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-3-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-016  
 Lab File ID: Y0413027.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
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07				
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09				
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28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413027.D Vial: 41  
Acq On : 13 Apr 2007 17:01 Operator: LNH  
Sample : JPL30-016 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413027.D 8260B.M Mon Apr 16 09:17:09 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-3-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-017  
 Lab File ID: Y0413028.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
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28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413028.D Vial: 42  
Acq On : 13 Apr 2007 17:26 Operator: LNH  
Sample : JPL30-017 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413028.D 8260B.M Mon Apr 16 09:17:51 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-3-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL30-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413029.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
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26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413029.D Vial: 43  
Acq On : 13 Apr 2007 17:50 Operator: LNH  
Sample : JPL30-018 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413029.D 8260B.M Mon Apr 16 09:18:42 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-5-4/2/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL30  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016833  
 Lab Sample ID: JPL30-019  
 Lab File ID: Y0413012.D  
 Date Collected: 04/03/2007  
 Date Analyzed: 04/13/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
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11					
12					
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29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413012.D Vial: 27  
Acq On : 13 Apr 2007 10:47 Operator: LNH  
Sample : JPL30-019 EB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413012.D 8260B.M Mon Apr 16 09:04:22 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041307MVOWY1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL30

Run Sequence: R016833

Matrix: (SOIL/WATER) Water

Lab Sample ID: B041307MVOWY1

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0413007.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/13/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
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09					
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11					
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29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041307\Y0413007.D Vial: 22  
Acq On : 13 Apr 2007 8:44 Operator: LNH  
Sample : B041307MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-1) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0413007.D 8260B.M Mon Apr 16 09:02:45 2007

Library Search Compound Report

Data File : X:\MSVOA\YODA\041607\Y0416010.D Vial: 24  
Acq On : 16 Apr 2007 11:07 Operator: LNH  
Sample : B041607MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-1) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0416010.D 8260B.M Thu Apr 19 11:18:50 2007

## **Metals Data**

**JPL30**

## COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks LaboratoriesContract: JPL Groundwater MonitorinLab Code: LAUCKSSDG No.: JPL30

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-20-5	JPL30-001
MW-20-4	JPL30-002
MW-20-4D	JPL30-002D
MW-20-4MS	JPL30-002MS
MW-20-3	JPL30-003
MW-20-2	JPL30-004
MW-20-1	JPL30-005
DUPE-3-1Q07	JPL30-006
EB-4-3/30/07	JPL30-007
MW-3-4	JPL30-016
MW-3-3	JPL30-017
MW-3-3MS	JPL30-017MS
MW-3-3MSD	JPL30-017MSD
MW-3-2	JPL30-018
EB-5-4/2/07	JPL30-019

Were ICP interelement corrections applied? Yes/No YESWere ICP background corrections applied? Yes/No NOIf yes-was raw data generated before application of background corrections? Yes/No NOComments:  
\_\_\_\_\_  
\_\_\_\_\_

I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill AmbacherName: Bill AmbacherDate: 4/24/07Title: Inorganics/Metals Manager

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-20-5

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-001

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-20-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-002

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016781

Color Before: Colorless Clarity Before: Clear Texture: \_\_\_\_\_

Color After: Colorless Clarity After: Clear Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-20-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-003

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.90			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-20-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-004

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-20-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-005

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.05			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-3-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-006

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-4-3/30/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-007

Level (low/med): LOW

Date Received: 03/31/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.42			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-3-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-016

Level (low/med): LOW

Date Received: 04/03/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.22			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-3-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-017

Level (low/med): LOW

Date Received: 04/03/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-3-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-018

Level (low/med): LOW

Date Received: 04/03/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.39			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-5-4/2/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL30

Matrix (soil/water): Water

Lab Sample ID: JPL30-019

Level (low/med): LOW

Date Received: 04/03/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.43			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL30**

**COVER PAGE - INORGANIC ANALYSES DATA PACKAGE**

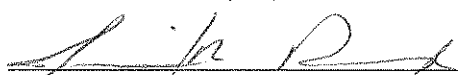
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL30

Client Identification	Lab Sample Work Order Number
MW-20-5	JPL30-001DL
MW-20-4	JPL30-002DL
MW-20-3	JPL30-003DL
MW-20-2	JPL30-004DL
MW-20-1	JPL30-005DL
DUP0E-3-1Q07	JPL30-006DL
EB-4-3/30/07	JPL30-007
MW-19-5	JPL30-009DL
MW-19-4	JPL30-010DL
MW-19-3	JPL30-011DL
MW-19-2	JPL30-012DL
MW-19-1	JPL30-013DL
DUPE-4-1Q07	JPL30-014DL
MW-3-4	JPL30-016DL
MW-3-3	JPL30-017DL
MW-3-3MS	JPL30-017MS
MW-3-3MSD	JPL30-017MSD
MW-3-2	JPL30-018DL
EB-5-4/2/07	JPL30-019

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Penner

Date: 4-20-07

Title: Inorganics Lead

## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-20-5 **Date/Time Collected:** 03/30/2007 08:05  
**Lab Sample ID:** JPL30-001 **Date/Time Received:** 03/31/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-20-4 **Date/Time Collected:** 03/30/2007 08:41  
**Lab Sample ID:** JPL30-002 **Date/Time Received:** 03/31/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-20-3 **Date/Time Collected:** 03/30/2007 09:19  
**Lab Sample ID:** JPL30-003 **Date/Time Received:** 03/31/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	2.5	U	2.5	0.35	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: MW-20-2 Date/Time Collected: 03/30/2007 09:51  
Lab Sample ID: JPL30-004 Date/Time Received: 03/31/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-20-1 **Date/Time Collected:** 03/30/2007 10:40  
**Lab Sample ID:** JPL30-005 **Date/Time Received:** 03/31/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: DUPE-3-1Q07 Date/Time Collected: 03/30/2007 00:00  
Lab Sample ID: JPL30-006 Date/Time Received: 03/31/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** EB-4-3/30/07 **Date/Time Collected:** 03/30/2007 10:29  
**Lab Sample ID:** JPL30-007 **Date/Time Received:** 03/31/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: MW-19-5 Date/Time Collected: 04/02/2007 07:48  
Lab Sample ID: JPL30-009 Date/Time Received: 04/03/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-19-4 **Date/Time Collected:** 04/02/2007 08:10  
**Lab Sample ID:** JPL30-010 **Date/Time Received:** 04/03/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	2.5	U	2.5	0.35	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-19-3 **Date/Time Collected:** 04/02/2007 08:30  
**Lab Sample ID:** JPL30-011 **Date/Time Received:** 04/03/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-19-2 **Date/Time Collected:** 04/02/2007 08:50  
**Lab Sample ID:** JPL30-012 **Date/Time Received:** 04/03/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: MW-19-1 Date/Time Collected: 04/02/2007 09:22  
Lab Sample ID: JPL30-013 Date/Time Received: 04/03/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	47		2.0	0.28	04/13/2007	04/14/2007	R016835



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** DUPE-4-1Q07 **Date/Time Collected:** 04/02/2007 00:00  
**Lab Sample ID:** JPL30-014 **Date/Time Received:** 04/03/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL30  
**Sample Number:** MW-3-4 **Date/Time Collected:** 04/02/2007 10:33  
**Lab Sample ID:** JPL30-016 **Date/Time Received:** 04/03/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: MW-3-3 Date/Time Collected: 04/02/2007 11:04  
Lab Sample ID: JPL30-017 Date/Time Received: 04/03/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/13/2007	04/14/2007	R016835

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: MW-3-2 Date/Time Collected: 04/02/2007 11:47  
Lab Sample ID: JPL30-018 Date/Time Received: 04/03/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	45		2.0	0.28	04/16/2007	04/17/2007	R016893

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL30  
Sample Number: EB-5-4/2/07 Date/Time Collected: 04/02/2007 09:07  
Lab Sample ID: JPL30-019 Date/Time Received: 04/03/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/16/2007	04/17/2007	R016893

**SAMPLE DATA**

SDG JPL31

VOLATILES ANALYSIS

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-001  
 Lab File ID: Y0415012.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 15:49  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.35	J

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-001  
 Lab File ID: Y0415012.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 15:49  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	1.7	
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-001  
 Lab File ID: Y0415012.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 15:49  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

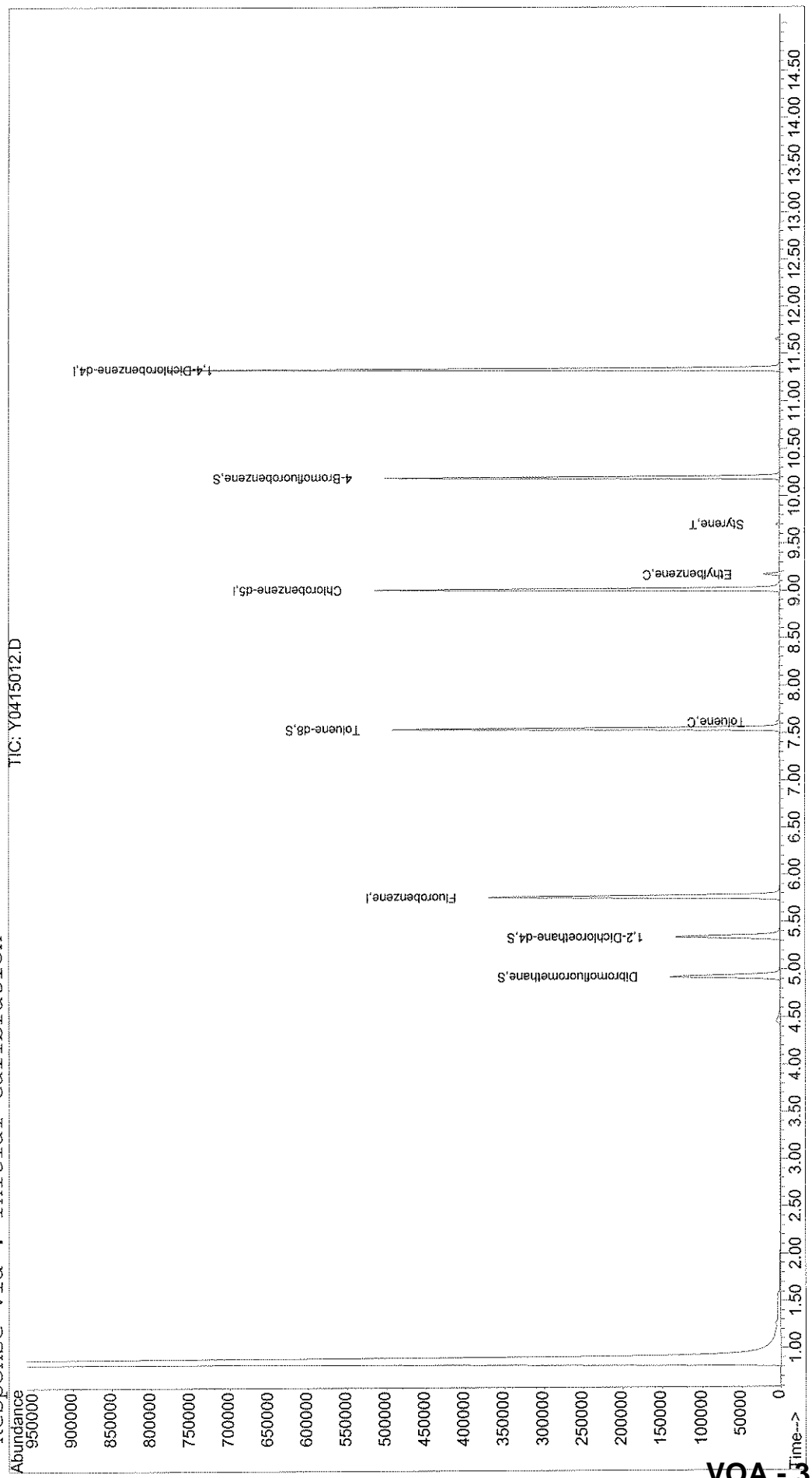
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415012.D  
Acq On : 15 Apr 2007 15:49  
Sample : JPL31-001  
Misc : 5mL+IS/SS #6  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:38 2007  
Vial: 25  
Operator: LH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415012.D  
 Acq On : 15 Apr 2007 15:49  
 Sample : JPL31-001  
 Misc : 5mL+IS/SS #6  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:38 2007

Vial: 25  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	341397	50.00	ug/l	0.00 100.49%
50) Chlorobenzene-d5	9.01	82	132251	50.00	ug/l	0.00 82.05%
70) 1,4-Dichlorobenzene-d4	11.34	152	179126	50.00	ug/l	0.00 79.49%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101130	48.96	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	97.92%	
37) 1,2-Dichloroethane-d4	5.34	65	104990	47.91	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	95.82%	
51) Toluene-d8	7.54	98	291397	55.52	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	111.04%	
72) 4-Bromofluorobenzene	10.20	95	133657	51.65	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	1.16	62	257	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	2.15	43	190	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	608	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	d
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415012.D 8260B.M Mon Apr 16 13:38:24 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415012.D  
 Acq On : 15 Apr 2007 15:49  
 Sample : JPL31-001  
 Misc : 5mL+IS/SS #6  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:38 2007

Vial: 25  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.40	78	363		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	7.61	92	1440	0.35	ug/l	98
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.04	112	288		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	13088	1.70	ug/l	99
65) m,p-Xylene	9.29	106	364		N.D.	
66) o-xylene	9.68	106	62		N.D.	
67) Styrene	9.70	104	2418	0.50	ug/l	95
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415012.D 8260B.M Mon Apr 16 13:38:24 2007

Quantitation Report

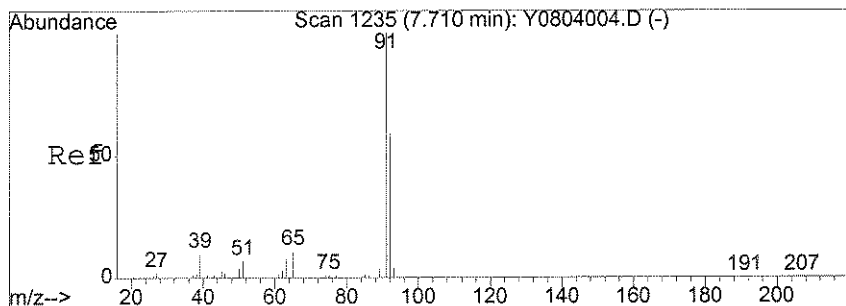
Data File : X:\MSVOA\YODA\041507\Y0415012.D  
 Acq On : 15 Apr 2007 15:49  
 Sample : JPL31-001  
 Misc : 5mL+IS/SS #6  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:38 2007

Vial: 25  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

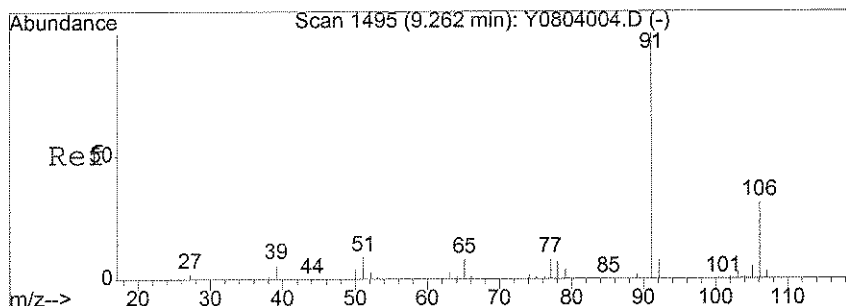
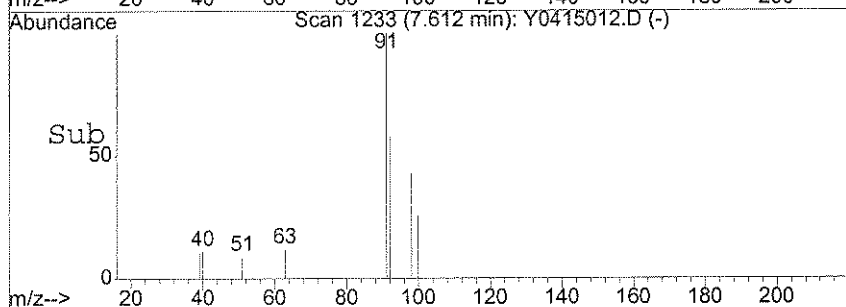
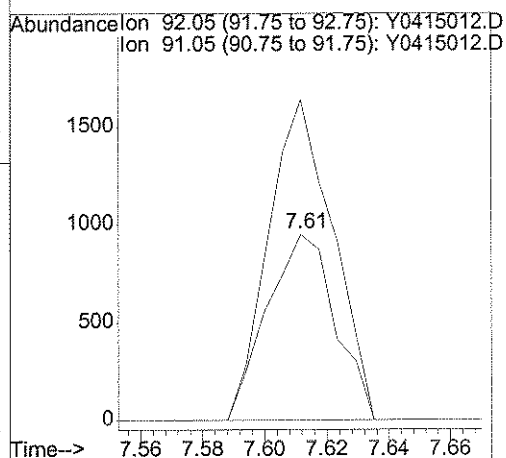
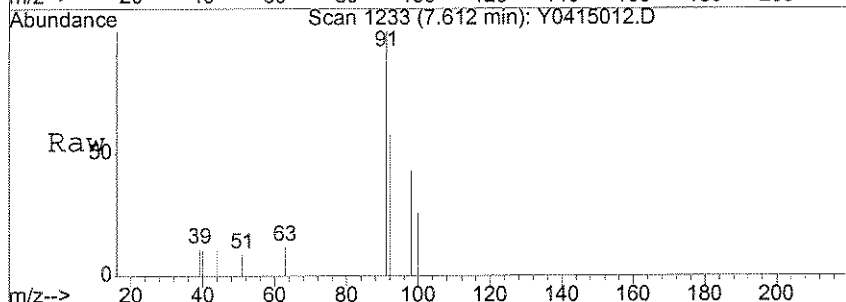
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.05	105	214		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.53	91	63		N.D.	
78) 4-Chlorotoluene	10.64	91	78		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	190		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	225		N.D.	
82) sec-butylbenzene	11.18	105	308		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	153		N.D.	
84) 4-Isopropyltoluene	11.34	119	627		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	314		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	65		N.D.	
87) n-Butylbenzene	11.74	91	618		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	151		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	13.55	128	190		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	86		N.D.	



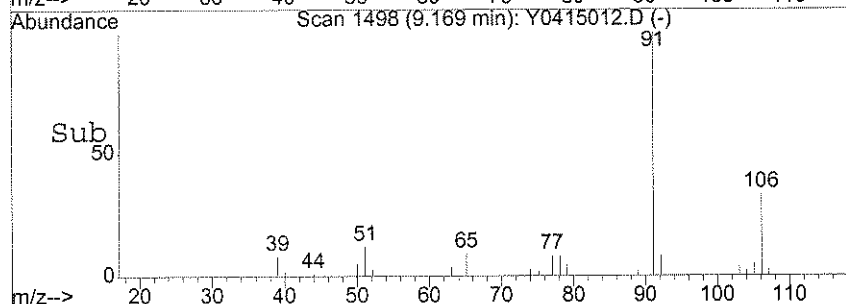
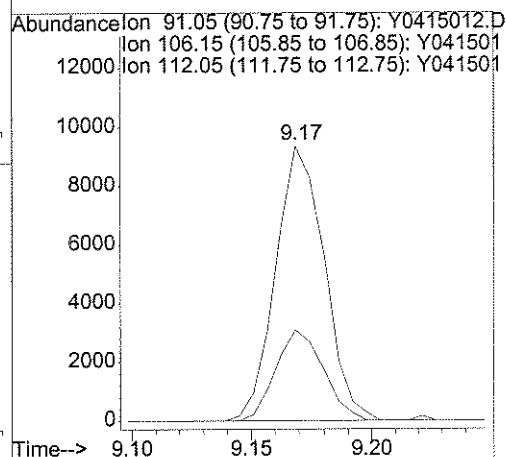
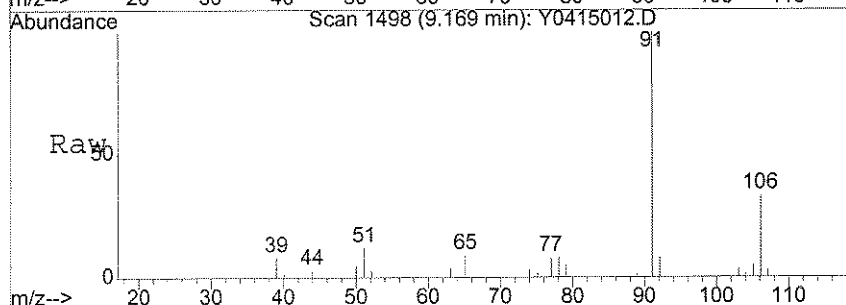
#52  
 Toluene  
 Concen: 0.35 ug/l  
 RT: 7.61 min Scan# 1233  
 Delta R.T. 0.00 min  
 Lab File: Y0415012.D  
 Acq: 15 Apr 2007 15:49

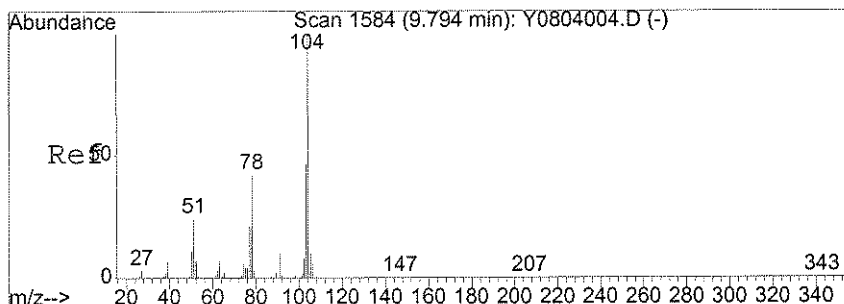
Tgt Ion: 92 Resp: 1440  
 Ion Ratio Lower Upper  
 92 100  
 91 163.8 133.7 200.5



#64  
 Ethylbenzene  
 Concen: 1.70 ug/l  
 RT: 9.17 min Scan# 1498  
 Delta R.T. 0.00 min  
 Lab File: Y0415012.D  
 Acq: 15 Apr 2007 15:49

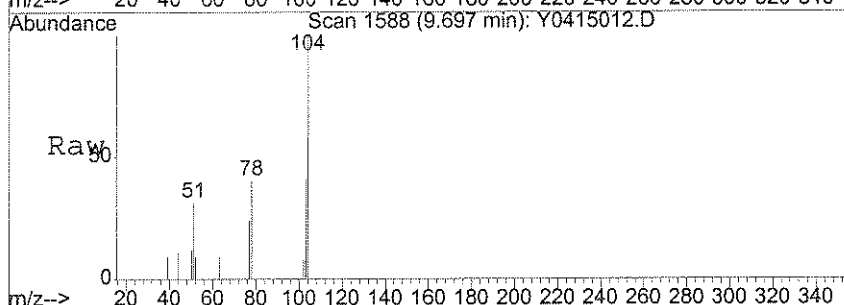
Tgt Ion: 91 Resp: 13088  
 Ion Ratio Lower Upper  
 91 100  
 106 33.0 26.1 39.1  
 112 0.0 0.0 0.0



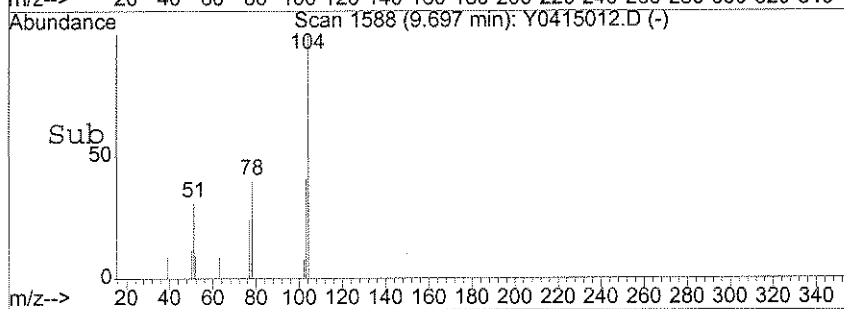
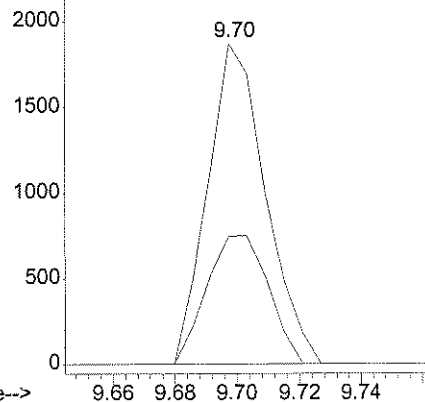


#67  
 Styrene  
 Concen: 0.50 ug/l  
 RT: 9.70 min Scan# 1588  
 Delta R.T. 0.00 min  
 Lab File: Y0415012.D  
 Acq: 15 Apr 2007 15:49

Tgt Ion	Ratio	Lower	Upper
104	100		
78	42.8	19.7	59.7



Abundance  
 Ion 104.00 (103.70 to 104.70): Y0415012.D  
 Ion 78.00 (77.70 to 78.70): Y0415012.D



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-002  
 Lab File ID: Y0415013.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 16:14  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.30	J
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.70	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 16:14

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.46	J
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-002  
 Lab File ID: Y0415013.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 16:14  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

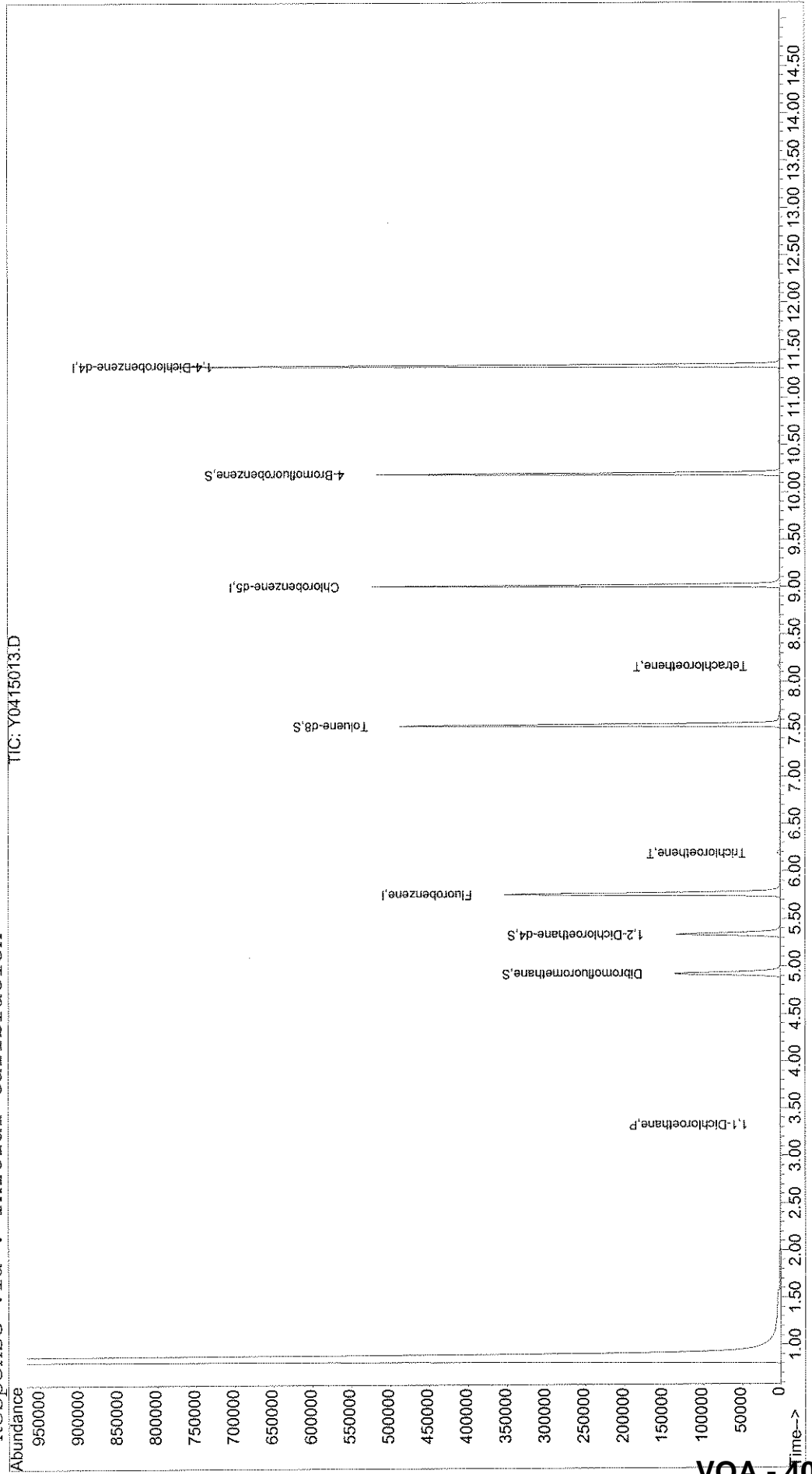
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415013.D  
Acq On : 15 Apr 2007 16:14  
Sample : JPL31-002  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:39 2007  
Vial: 26  
Operator: LH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415013.D  
 Acq On : 15 Apr 2007 16:14  
 Sample : JPL31-002  
 Misc : 5mL+IS/SS #3

Vial: 26  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	330820	50.00	ug/l	0.00 97.38%
50) Chlorobenzene-d5	9.01	82	132406	50.00	ug/l	0.00 82.15%
70) 1,4-Dichlorobenzene-d4	11.34	152	180515	50.00	ug/l	0.00 80.10%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	100616	50.27	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.54%
37) 1,2-Dichloroethane-d4	5.34	65	105058	49.48	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.96%
51) Toluene-d8	7.54	98	290283	55.24	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	110.48%
72) 4-Bromofluorobenzene	10.20	95	136328	52.27	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.21	76	86	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	3.32	63	1392	0.30	ug/l	87
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0415013.D 8260B.M Mon Apr 16 13:39:17 2007

*Handwritten:* 87  
 LW 4/16/07

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415013.D  
 Acq On : 15 Apr 2007 16:14  
 Sample : JPL31-002  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Vial: 26  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.70	83	548		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	53		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	1635	0.70	ug/l ✓	97
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	7.61	92	58		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	834	0.46	ug/l # ✓	86
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	407		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.16	91	350		N.D.	
65) m,p-Xylene	9.29	106	129		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*LH 4/16/07*

(#) = qualifier out of range (m) = manual integration  
 Y0415013.D 8260B.M Mon Apr 16 13:39:18 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415013.D  
 Acq On : 15 Apr 2007 16:14  
 Sample : JPL31-002  
 Misc : 5mL+IS/SS #3

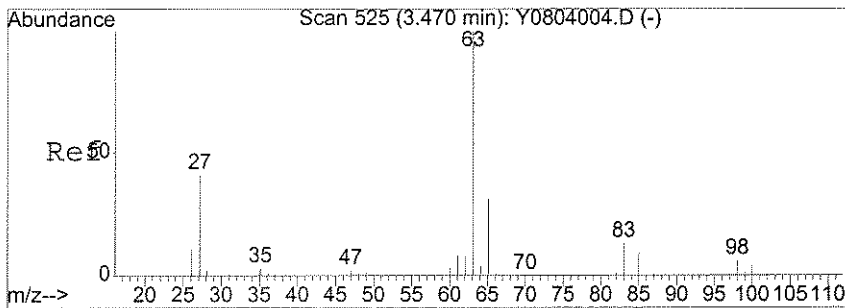
Vial: 26  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Quant Results File: 8260B.RES

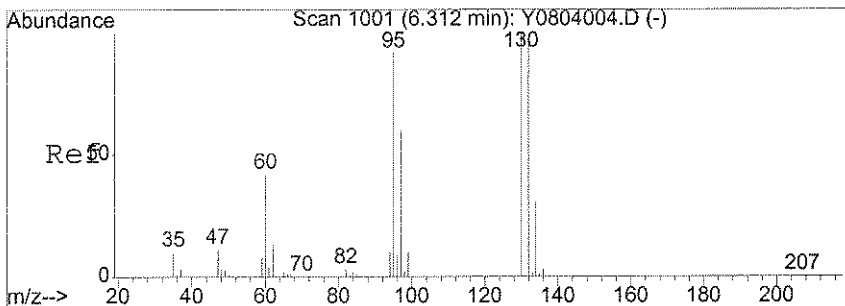
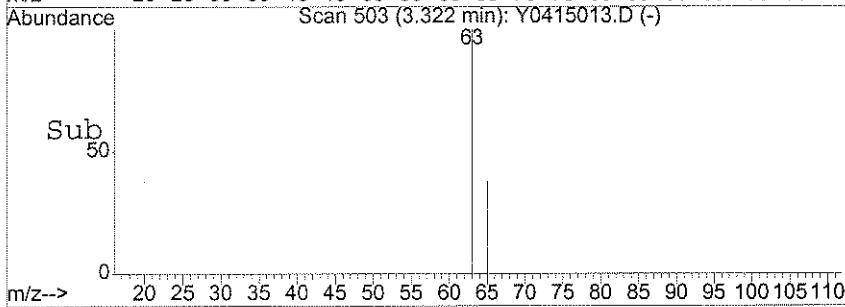
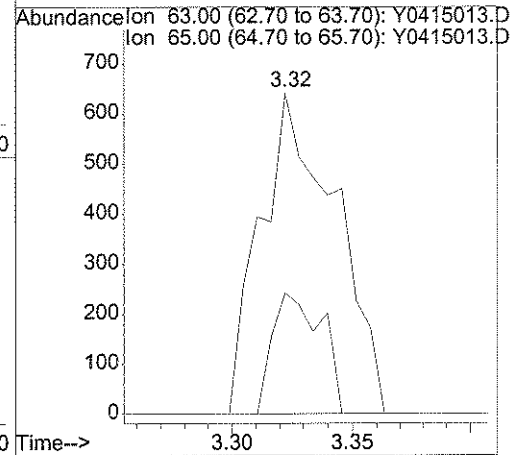
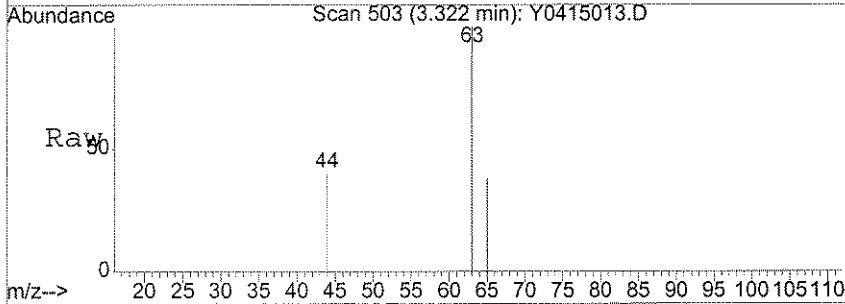
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	10.65	120	55		N.D.	
77) 2-Chlorotoluene	10.54	91	55		N.D.	
78) 4-Chlorotoluene	10.64	91	115		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	63		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.02	105	57		N.D.	
82) sec-butylbenzene	11.18	105	323		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	167		N.D.	
84) 4-Isopropyltoluene	11.34	119	448		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	434		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	277		N.D.	
87) n-Butylbenzene	11.74	91	484		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	210		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	13.55	128	58		N.D.	
92) 1,2,3-Trichlorobenzene	13.78	180	358		N.D.	



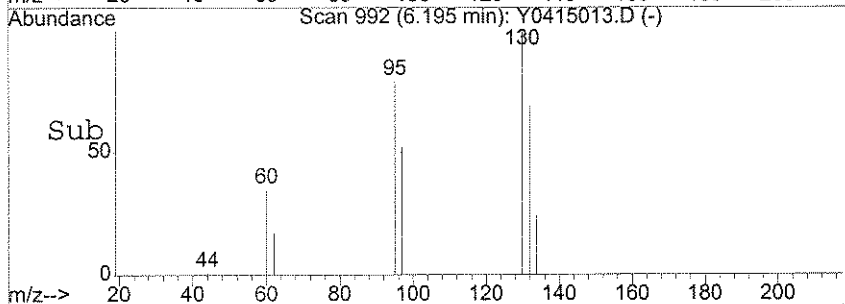
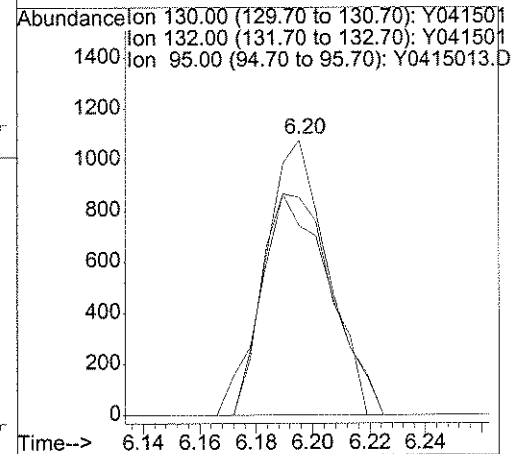
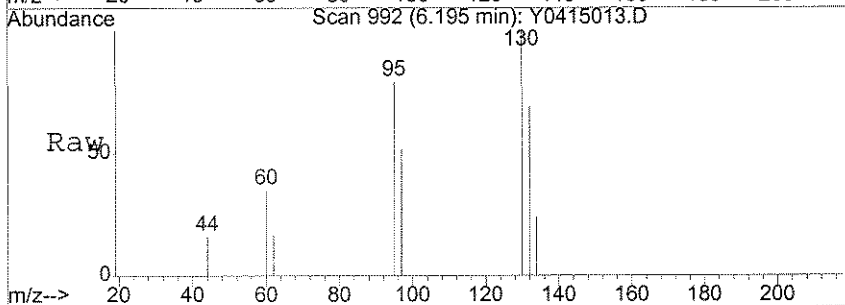
#22  
 1,1-Dichloroethane  
 Concen: 0.30 ug/l  
 RT: 3.32 min Scan# 503  
 Delta R.T. -0.00 min  
 Lab File: Y0415013.D  
 Acq: 15 Apr 2007 16:14

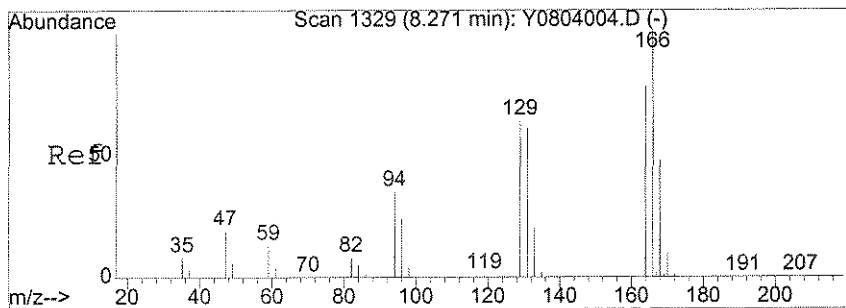
Tgt Ion	Resp	Lower	Upper
63	1392		
63	100		
65	24.9	12.3	52.3



#41  
 Trichloroethene  
 Concen: 0.70 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0415013.D  
 Acq: 15 Apr 2007 16:14

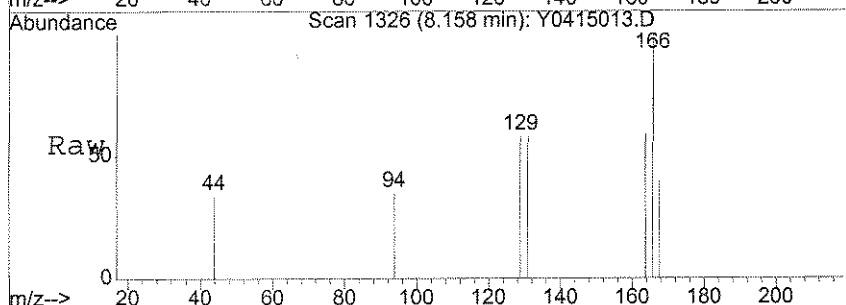
Tgt Ion	Resp	Lower	Upper
130	1635		
130	100		
132	90.8	75.0	115.0
95	88.7	69.4	109.4



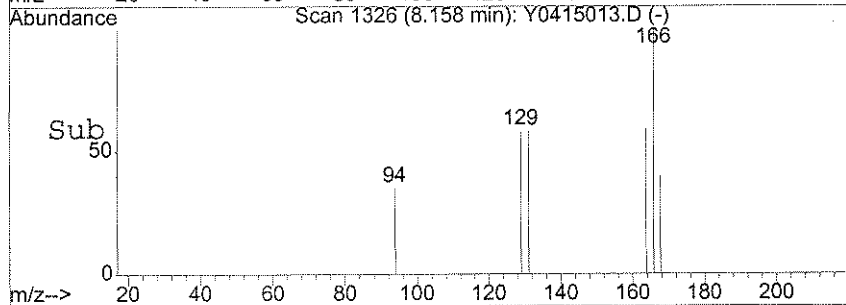
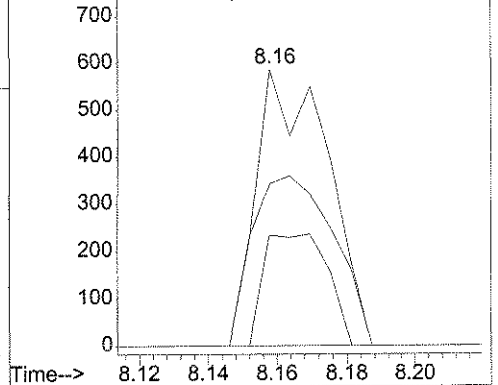


#56  
 Tetrachloroethene  
 Concen: 0.46 ug/l  
 RT: 8.16 min Scan# 1326  
 Delta R.T. -0.01 min  
 Lab File: Y0415013.D  
 Acq: 15 Apr 2007 16:14

Tgt Ion	Resp	Lower	Upper
166	100		
164	70.1	63.3	94.9
168	36.1	39.6	59.4#



Abundance  
 Ion 165.95 (165.65 to 166.65): Y041501  
 Ion 163.95 (163.65 to 164.65): Y041501  
 Ion 167.95 (167.65 to 168.65): Y041501





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-003  
 Lab File ID: Y0415014.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 16:38  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-003  
 Lab File ID: Y0415014.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 16:38  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-4-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-003  
 Lab File ID: Y0415014.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 16:38  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

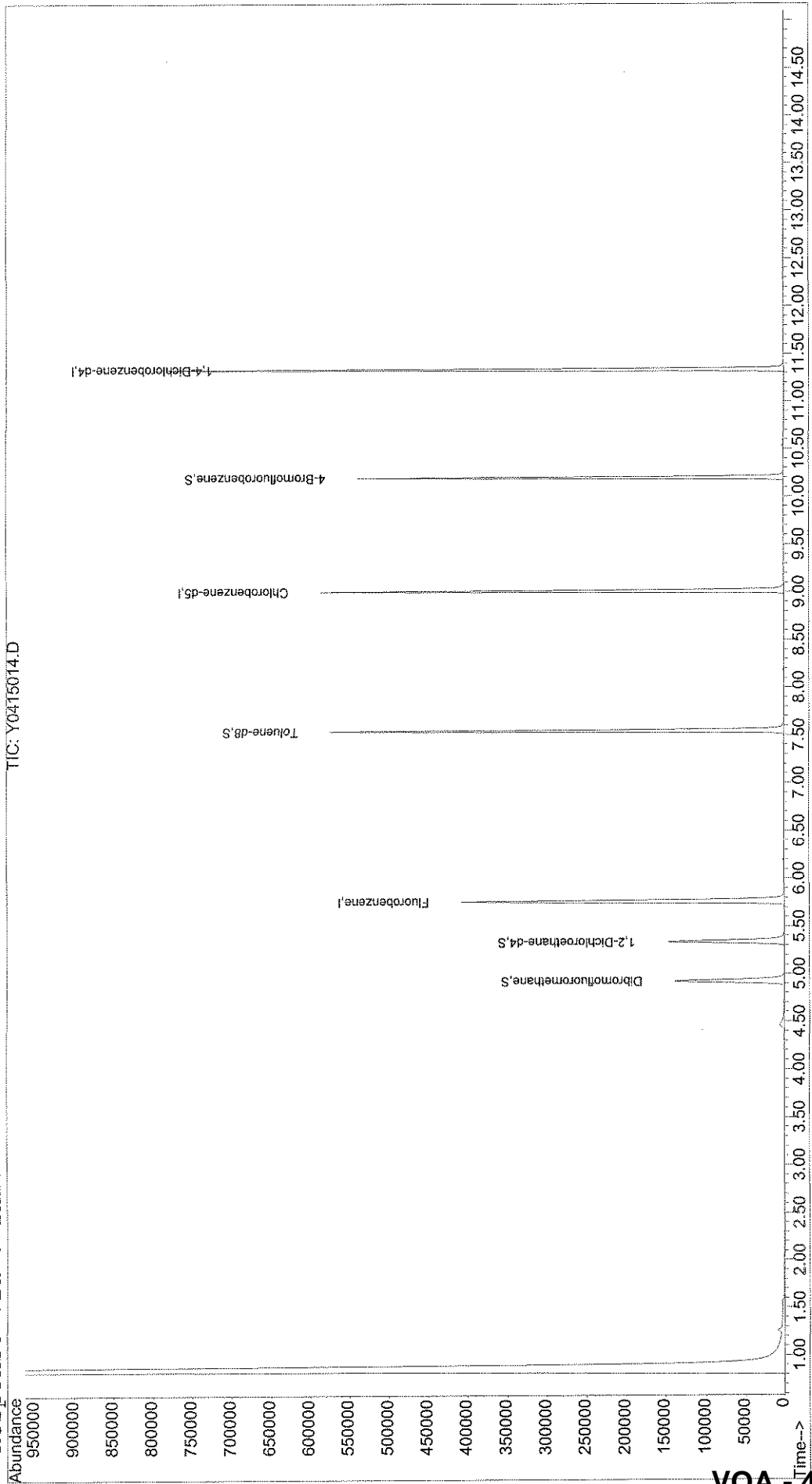
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415014.D Vial: 27  
Acq On : 15 Apr 2007 16:38 Operator: LH  
Sample : JPL31-003 Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:39 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415014.D  
 Acq On : 15 Apr 2007 16:38  
 Sample : JPL31-003  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Vial: 27  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	368612	50.00	ug/l	0.00 108.50%
50) Chlorobenzene-d5	9.01	82	148596	50.00	ug/l	0.00 92.19%
70) 1,4-Dichlorobenzene-d4	11.34	152	180808	50.00	ug/l	0.00 80.23%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101624	45.57	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	91.14%
37) 1,2-Dichloroethane-d4	5.34	65	113102	47.80	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.60%
51) Toluene-d8	7.54	98	341504	57.91	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	115.82%
72) 4-Bromofluorobenzene	10.20	95	140099	53.63	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	283	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415014.D 8260B.M Mon Apr 16 13:40:09 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415014.D  
 Acq On : 15 Apr 2007 16:38  
 Sample : JPL31-003  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Vial: 27  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	834		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	5.13	75	54		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	518		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	340		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.17	166	60		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	508		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	167		N.D.	
65) m,p-Xylene	9.29	106	179		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415014.D 8260B.M Mon Apr 16 13:40:09 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415014.D  
 Acq On : 15 Apr 2007 16:38  
 Sample : JPL31-003  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:39 2007

Vial: 27  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	73		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	289		N.D.	
78) 4-Chlorotoluene	10.64	91	53		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	56		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	218		N.D.	
82) sec-butylbenzene	11.18	105	218		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	119		N.D.	
84) 4-Isopropyltoluene	11.34	119	491		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	119		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	416		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-004  
 Lab File ID: Y0415015.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 17:02  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-004  
 Lab File ID: Y0415015.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 17:02  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 17:02

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

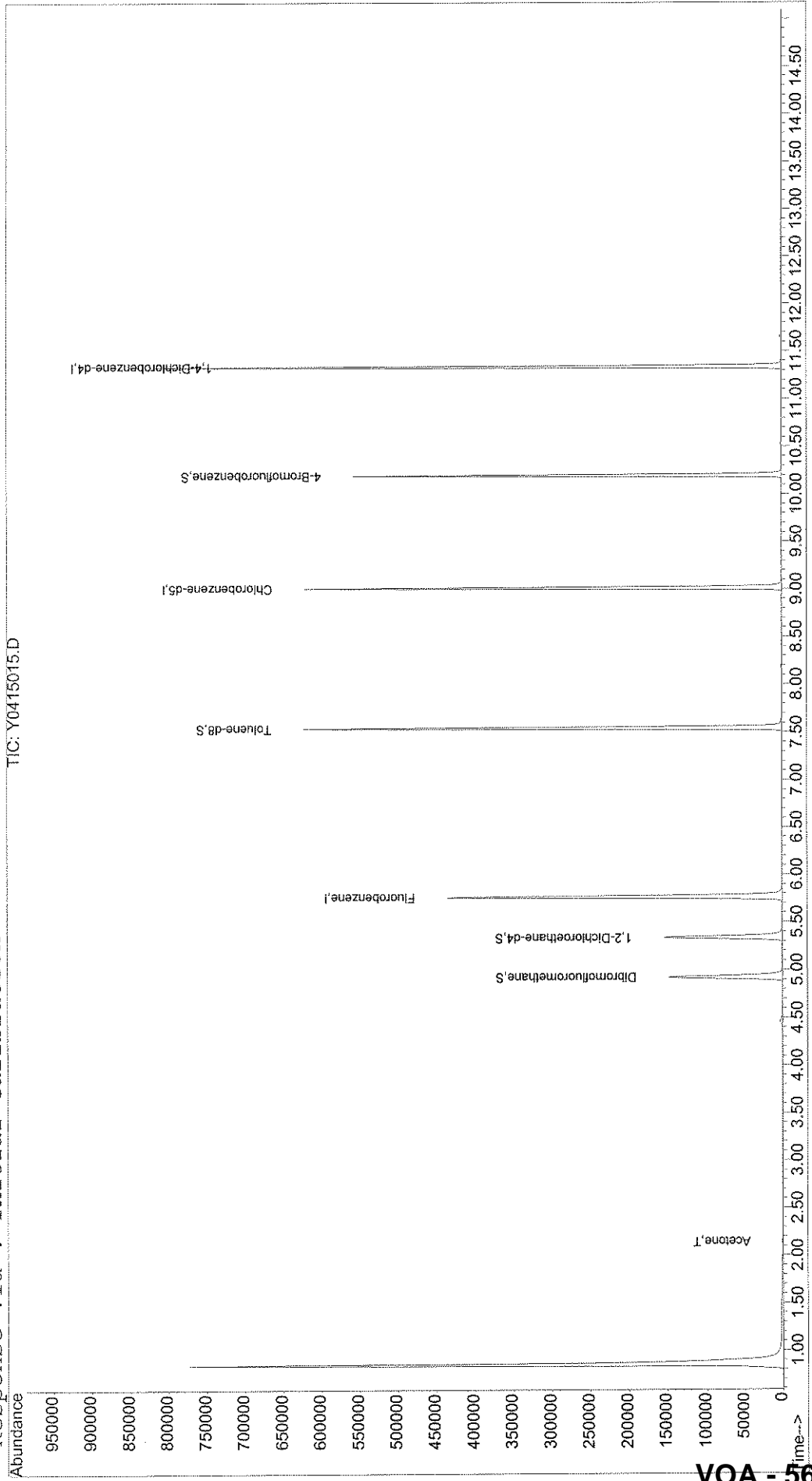
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415015.D Vial: 28  
Acq On : 15 Apr 2007 17:02 Operator: LH  
Sample : JPL31-004 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:41 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415015.D  
 Acq On : 15 Apr 2007 17:02  
 Sample : JPL31-004  
 Misc : 5mL+IS/SS #1

Vial: 28  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	5.76	96	394934	50.00	ug/l	0.00	116.25%
50) Chlorobenzene-d5	9.01	82	156535	50.00	ug/l	0.00	97.12%
70) 1,4-Dichlorobenzene-d4	11.34	152	186692	50.00	ug/l	0.00	82.84%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	107355	44.93	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	89.86%	
37) 1,2-Dichloroethane-d4	5.34	65	121132	47.79	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.58%	
51) Toluene-d8	7.54	98	366653	59.02	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	118.04%	
72) 4-Bromofluorobenzene	10.20	95	145281	53.86	ug/l	0.00	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.14	43	1972	2.19	ug/l #	48
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	136	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	2.42	40	54	N.D.		
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

LH 4/16/07

(#) = qualifier out of range (m) = manual integration  
 Y0415015.D 8260B.M Mon Apr 16 13:41:08 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415015.D  
 Acq On : 15 Apr 2007 17:02  
 Sample : JPL31-004  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Vial: 28  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	227		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	501		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	236		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415015.D 8260B.M Mon Apr 16 13:41:09 2007

Quantitation Report

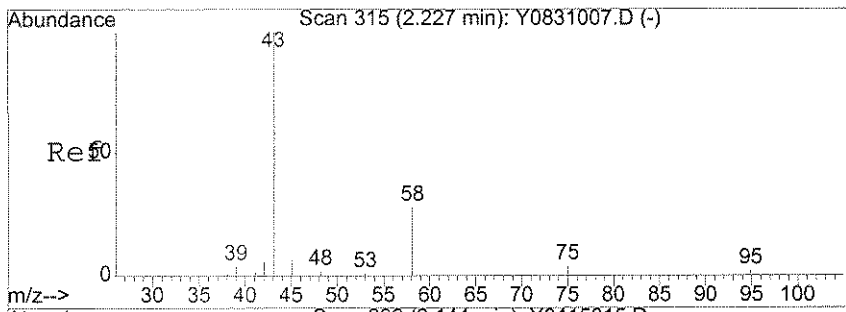
Data File : X:\MSVOA\YODA\041507\Y0415015.D  
 Acq On : 15 Apr 2007 17:02  
 Sample : JPL31-004  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Vial: 28  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

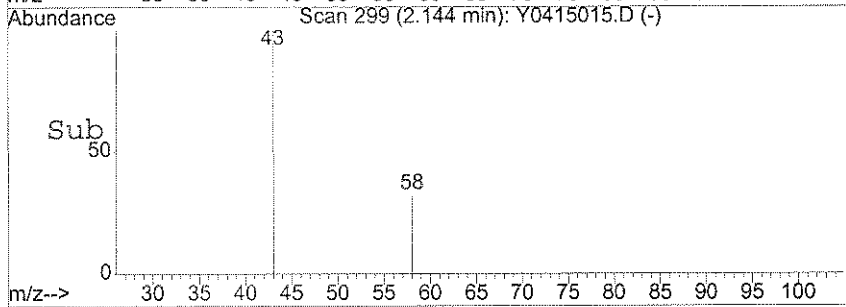
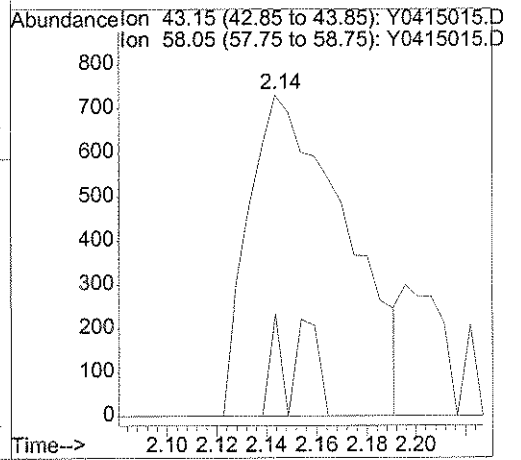
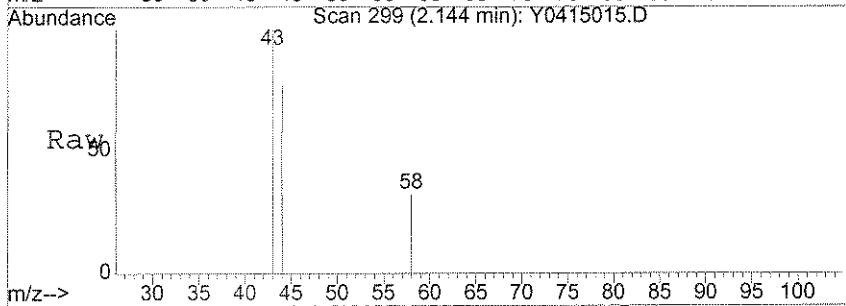
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.46	91	240		N.D.	
78) 4-Chlorotoluene	10.46	91	240		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	146		N.D.	
82) sec-butylbenzene	11.17	105	54		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	61		N.D.	
84) 4-Isopropyltoluene	11.34	119	270		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	61		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	345		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	13.54	128	53		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



#11  
 Acetone  
 Concen: 2.19 ug/l  
 RT: 2.14 min Scan# 299  
 Delta R.T. -0.01 min  
 Lab File: Y0415015.D  
 Acq: 15 Apr 2007 17:02

Tgt Ion	Resp	Lower	Upper
43	1972		
58	3.7	26.8	40.2#



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-005  
 Lab File ID: Y0415010.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 15:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-005  
 Lab File ID: Y0415010.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 15:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415010.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 15:00

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

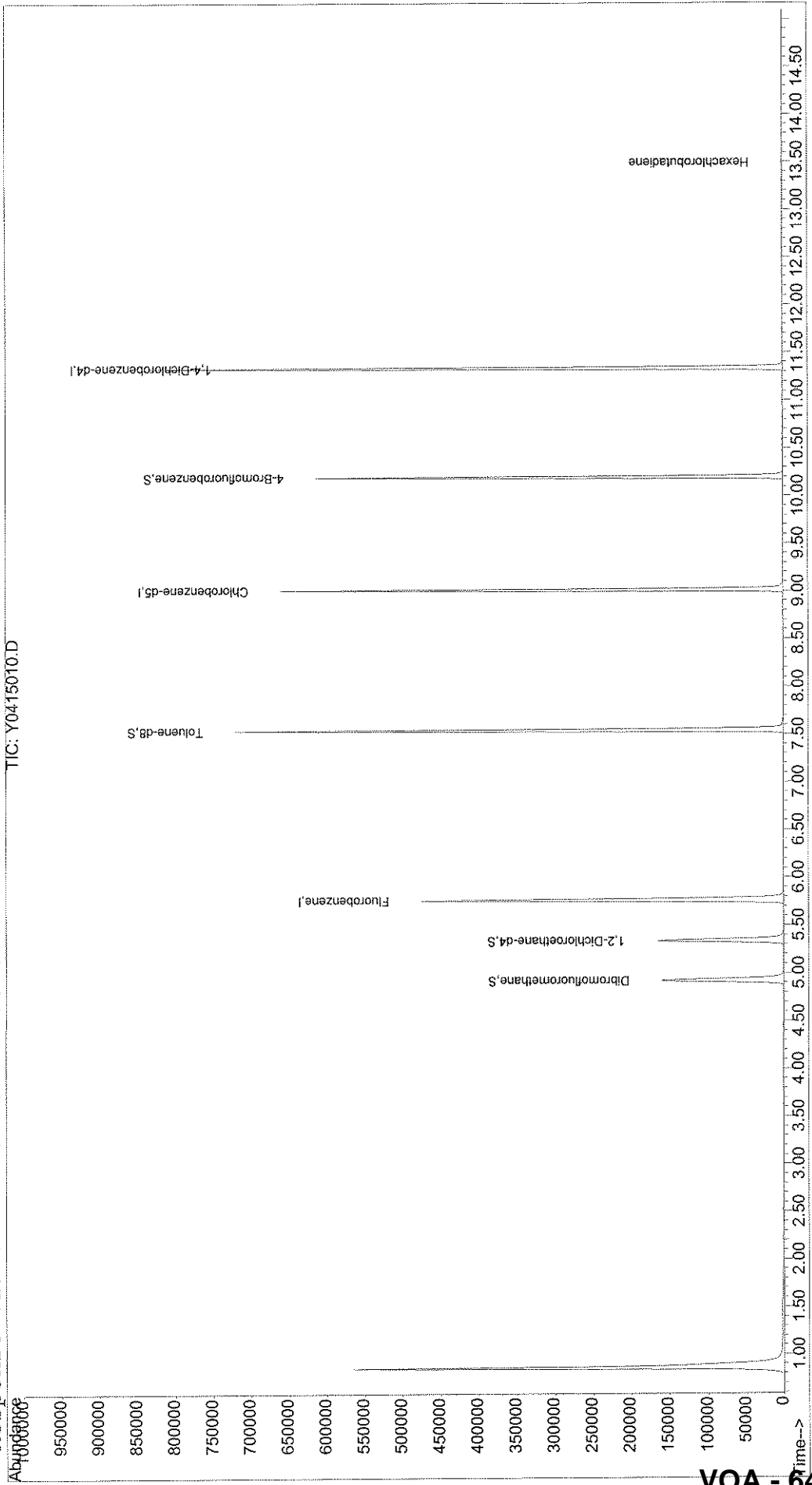
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	1.5	B
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415010.D  
Acq On : 15 Apr 2007 15:00 Vial: 23  
Sample : JPL31-005 TB Operator: LH  
Misc : 5mL+IS/SS #1 Inst : Yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 16 13:35 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415010.D  
 Acq On : 15 Apr 2007 15:00  
 Sample : JPL31-005 TB  
 Misc : 5mL+IS/SS #1

Vial: 23  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:35 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	442094	50.00	ug/l	0.00 130.13%
50) Chlorobenzene-d5	9.01	82	170253	50.00	ug/l	0.00 105.63%
70) 1,4-Dichlorobenzene-d4	11.34	152	193911	50.00	ug/l	0.00 86.05%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	123068	46.01	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	92.02%
37) 1,2-Dichloroethane-d4	5.34	65	131311	46.27	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	92.54%
51) Toluene-d8	7.54	98	419840	62.14	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	124.28%#
72) 4-Bromofluorobenzene	10.20	95	163368	58.31	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.15	43	457	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	4207	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	2.52	84	477	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0415010.D 8260B.M Mon Apr 16 13:36:00 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415010.D  
 Acq On : 15 Apr 2007 15:00  
 Sample : JPL31-005 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:35 2007

Vial: 23  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	83		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	6.38	83	213		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	121		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	198		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.03	112	89		N.D.	
62) 1-Chlorohexane	9.06	91	61		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	357		N.D.	
65) m,p-Xylene	9.29	106	223		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	64		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415010.D 8260B.M Mon Apr 16 13:36:01 2007

Quantitation Report

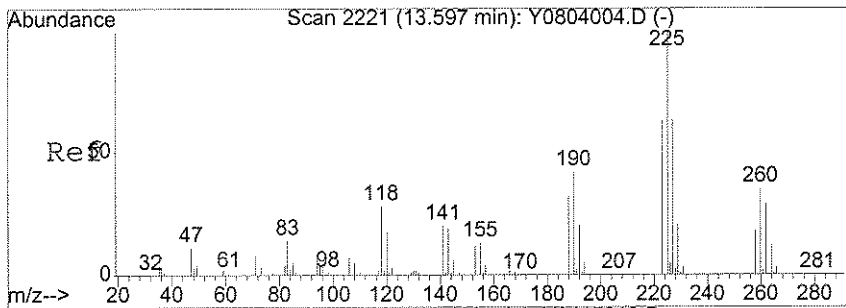
Data File : X:\MSVOA\YODA\041507\Y0415010.D  
 Acq On : 15 Apr 2007 15:00  
 Sample : JPL31-005 TB  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:35 2007

Vial: 23  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

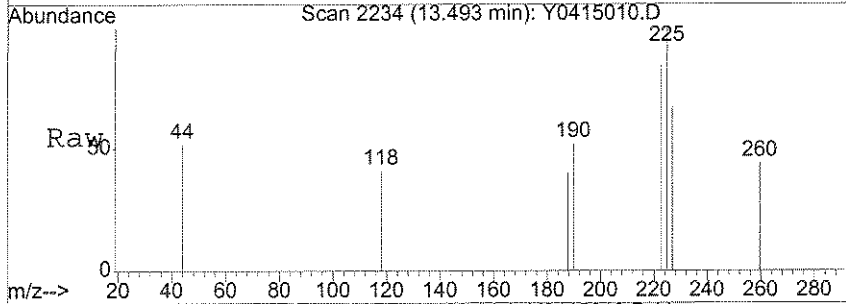
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.06	105	413		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	10.64	120	71		N.D.	
77) 2-Chlorotoluene	10.53	91	250		N.D.	
78) 4-Chlorotoluene	10.53	91	250		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	318		N.D.	
80) tert-Butylbenzene	10.96	119	325		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	402		N.D.	
82) sec-butylbenzene	11.18	105	691		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	304		N.D.	
84) 4-Isopropyltoluene	11.34	119	1064		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	404		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	72		N.D.	
87) n-Butylbenzene	11.74	91	1114		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	464		N.D.	
90) Hexachlorobutadiene	13.49	225	404	1.48	ug/l #	78
91) Naphthalene	13.55	128	470		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	440		N.D.	

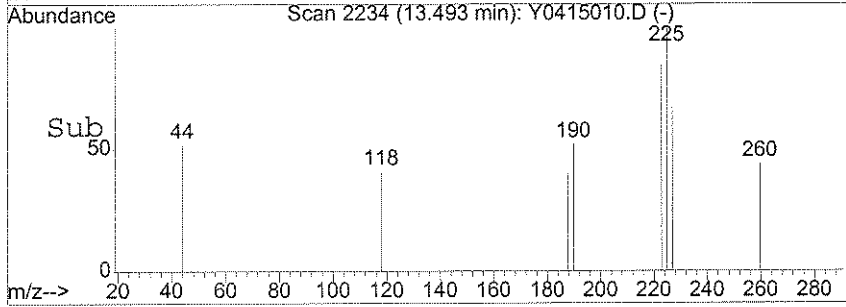
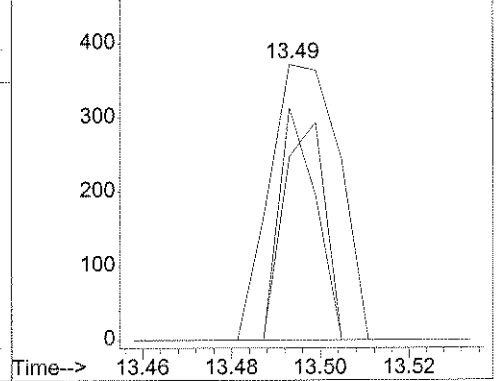


#90  
 Hexachlorobutadiene  
 Concen: 1.48 ug/l  
 RT: 13.49 min Scan# 2234  
 Delta R.T. -0.01 min  
 Lab File: Y0415010.D  
 Acq: 15 Apr 2007 15:00

Tgt Ion	Ratio	Lower	Upper
225	100		
223	44.6	50.2	75.2#
227	47.3	50.2	75.4#



Abundance Ion 224.80 (224.50 to 225.50): Y041501  
 Ion 222.80 (222.50 to 223.50): Y041501  
 Ion 226.80 (226.50 to 227.50): Y041501



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 17:27

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 17:27

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-006  
 Lab File ID: Y0415016.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 17:27  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

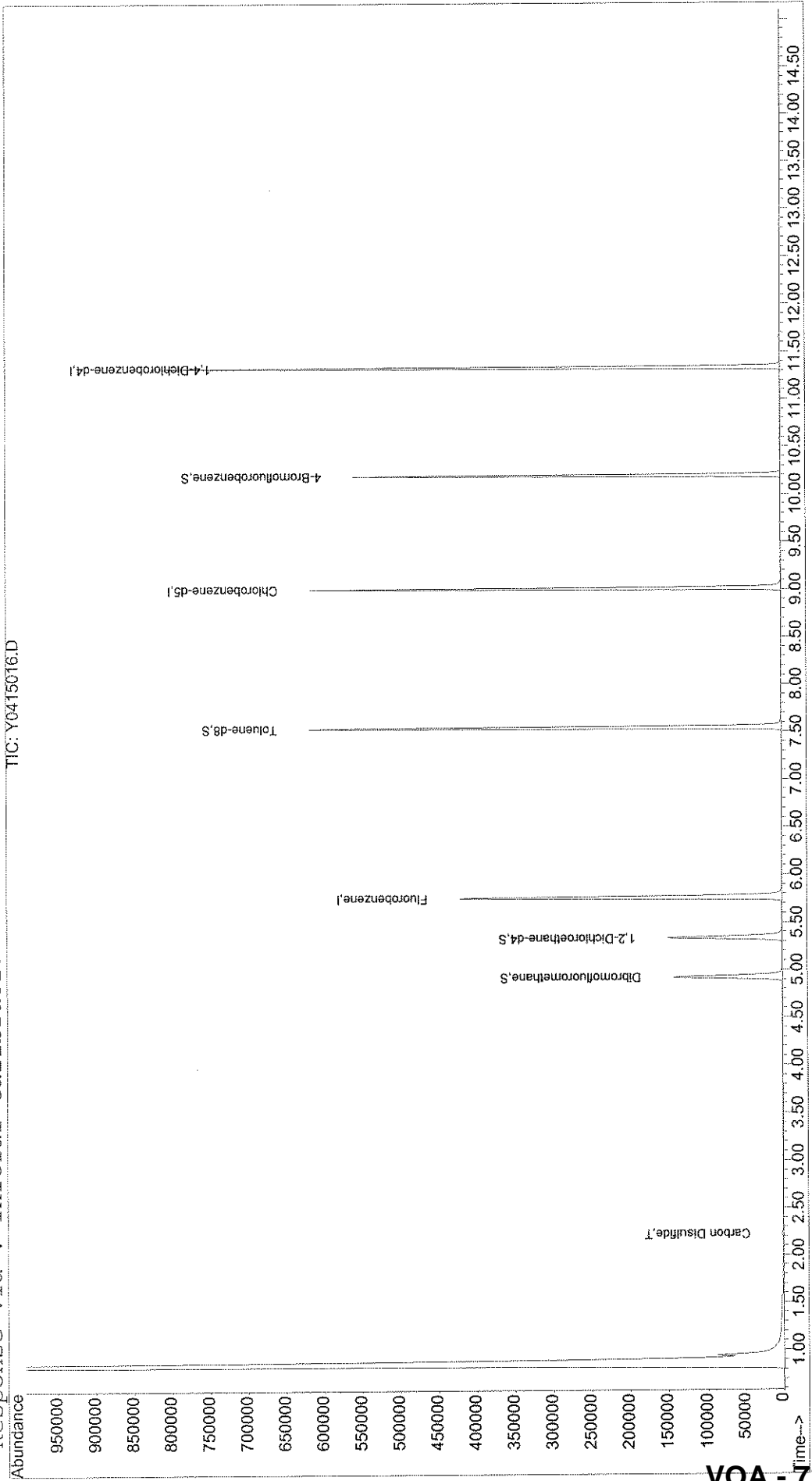
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415016.D  
Acq On : 15 Apr 2007 17:27  
Sample : JPL31-006  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:41 2007  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415016.D  
 Acq On : 15 Apr 2007 17:27  
 Sample : JPL31-006  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Vial: 29  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	384540	50.00	ug/l	0.00 113.19%
50) Chlorobenzene-d5	9.01	82	154052	50.00	ug/l	0.00 95.58%
70) 1,4-Dichlorobenzene-d4	11.34	152	181260	50.00	ug/l	0.00 80.43%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	103049	44.29	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	88.58%	
37) 1,2-Dichloroethane-d4	5.34	65	117753	47.71	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	95.42%	
51) Toluene-d8	7.54	98	360887	59.03	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	118.06%	
72) 4-Bromofluorobenzene	10.20	95	144911	55.34	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	4317	0.55	ug/l	100
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

100  
 4/16/07

(#) = qualifier out of range (m) = manual integration  
 Y0415016.D 8260B.M Mon Apr 16 13:41:51 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415016.D  
 Acq On : 15 Apr 2007 17:27  
 Sample : JPL31-006  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Vial: 29  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	274		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	489		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	514		N.D.	
65) m,p-Xylene	9.29	106	276		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	9.70	104	905		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415016.D 8260B.M Mon Apr 16 13:41:51 2007

Quantitation Report

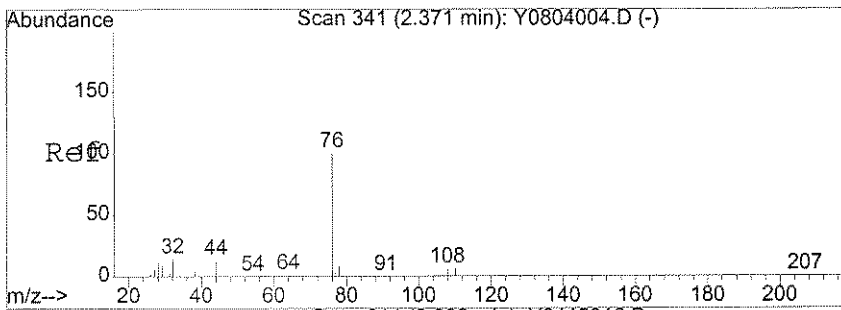
Data File : X:\MSVOA\YODA\041507\Y0415016.D  
 Acq On : 15 Apr 2007 17:27  
 Sample : JPL31-006  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:41 2007

Vial: 29  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

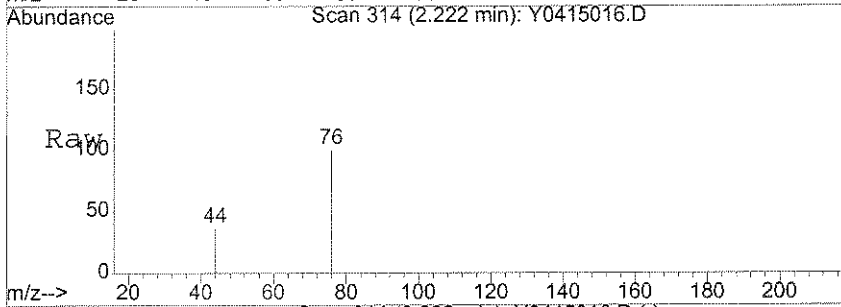
Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

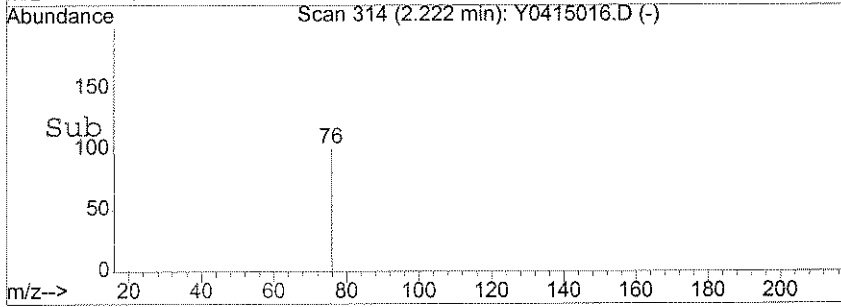
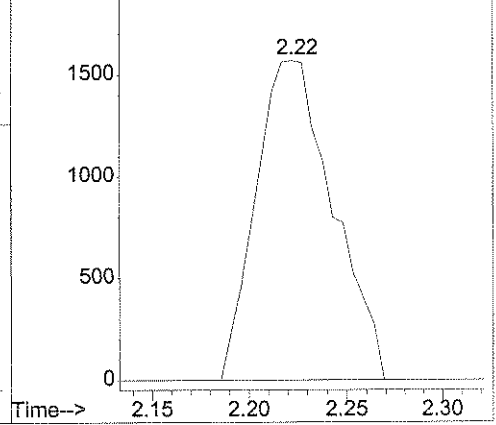
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.19	105	60		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.53	91	61		N.D.	
78) 4-Chlorotoluene	10.53	91	61		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	116		N.D.	
82) sec-butylbenzene	11.18	105	114		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	55		N.D.	
84) 4-Isopropyltoluene	11.34	119	385		N.D.	
85) 1,4-Dichlorobenzene	11.27	146	55		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	377		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



#14  
 Carbon Disulfide  
 Concen: 0.55 ug/l  
 RT: 2.22 min Scan# 314  
 Delta R.T. 0.01 min  
 Lab File: Y0415016.D  
 Acq: 15 Apr 2007 17:27  
 Tgt Ion: 76 Resp: 4317



Abundance on 76.00 (75.70 to 76.70): Y0415016.D



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-007  
 Lab File ID: Y0415017.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 17:51  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 17:51

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-007  
 Lab File ID: Y0415017.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 17:51  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

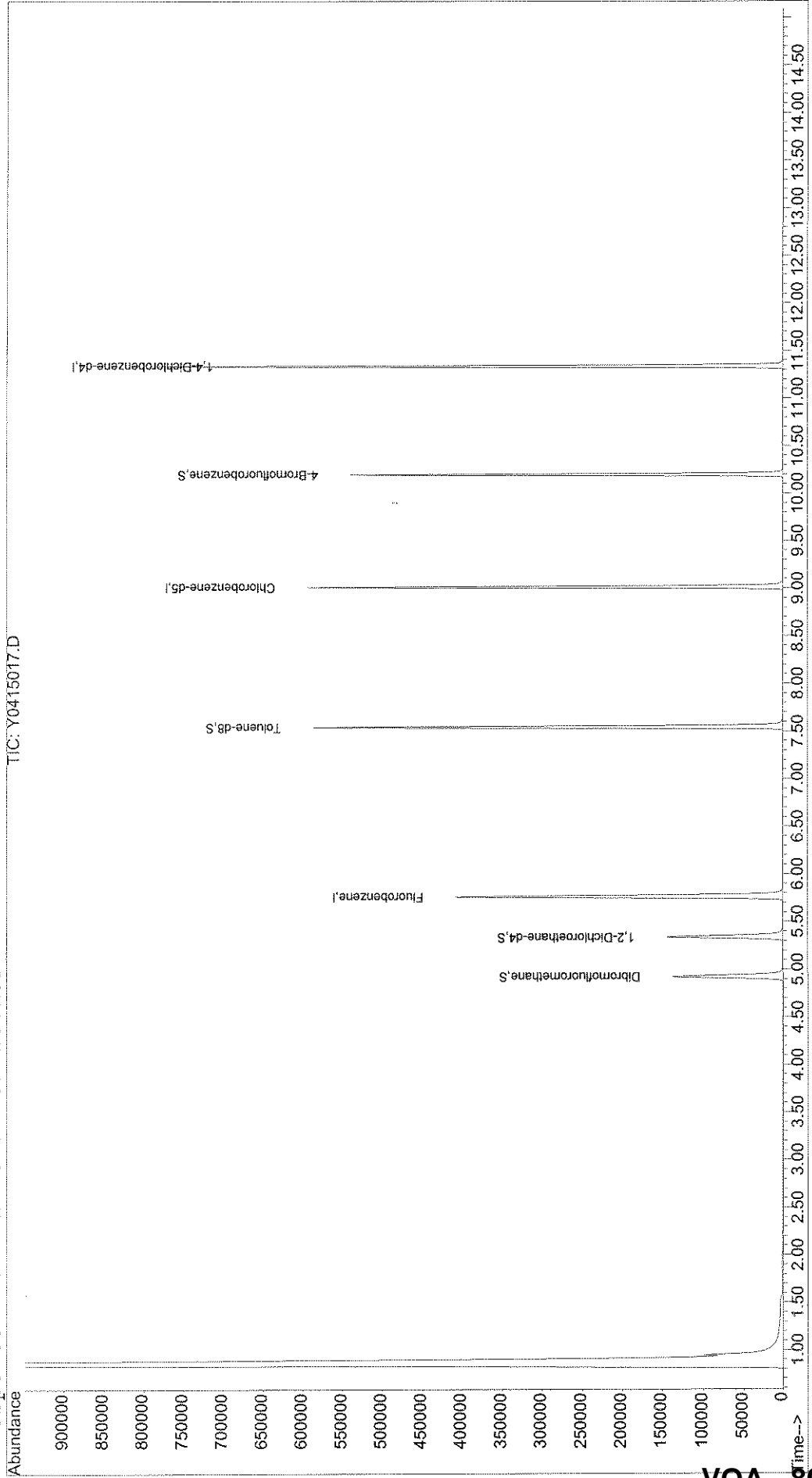
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415017.D Vial: 30  
Acq On : 15 Apr 2007 17:51 Operator: LH  
Sample : JPL31-007 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:42 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415017.D  
 Acq On : 15 Apr 2007 17:51  
 Sample : JPL31-007  
 Misc : 5mL+IS/SS #1

Vial: 30  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:42 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

Response via : Initial Calibration

DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	367209	50.00	ug/l	0.00 108.09%
50) Chlorobenzene-d5	9.01	82	148353	50.00	ug/l	0.00 92.04%
70) 1,4-Dichlorobenzene-d4	11.34	152	175584	50.00	ug/l	0.00 77.92%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	100448	45.21	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	90.42%
37) 1,2-Dichloroethane-d4	5.34	65	112151	47.58	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.16%
51) Toluene-d8	7.55	98	348616	59.21	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	118.42%
72) 4-Bromofluorobenzene	10.20	95	138016	54.41	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	1825	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415017.D  
 Acq On : 15 Apr 2007 17:51  
 Sample : JPL31-007  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:42 2007

Vial: 30  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	425		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	54		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	486		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	142		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415017.D 8260B.M Mon Apr 16 13:43:55 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415017.D  
 Acq On : 15 Apr 2007 17:51  
 Sample : JPL31-007  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:42 2007

Vial: 30  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	150		N.D.	
78) 4-Chlorotoluene	10.47	91	150		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.19	105	170		N.D.	
82) sec-butylbenzene	11.19	105	170		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	59		N.D.	
84) 4-Isopropyltoluene	11.33	119	325		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	59		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.75	91	256		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415017.D 8260B.M Mon Apr 16 13:43:56 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-008  
 Lab File ID: Y0415018.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 18:15  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-008  
 Lab File ID: Y0415018.D  
 Date Collected: 04/03/2007  
 Date/Time Analyzed: 04/15/2007 18:15  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



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VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 18:15

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

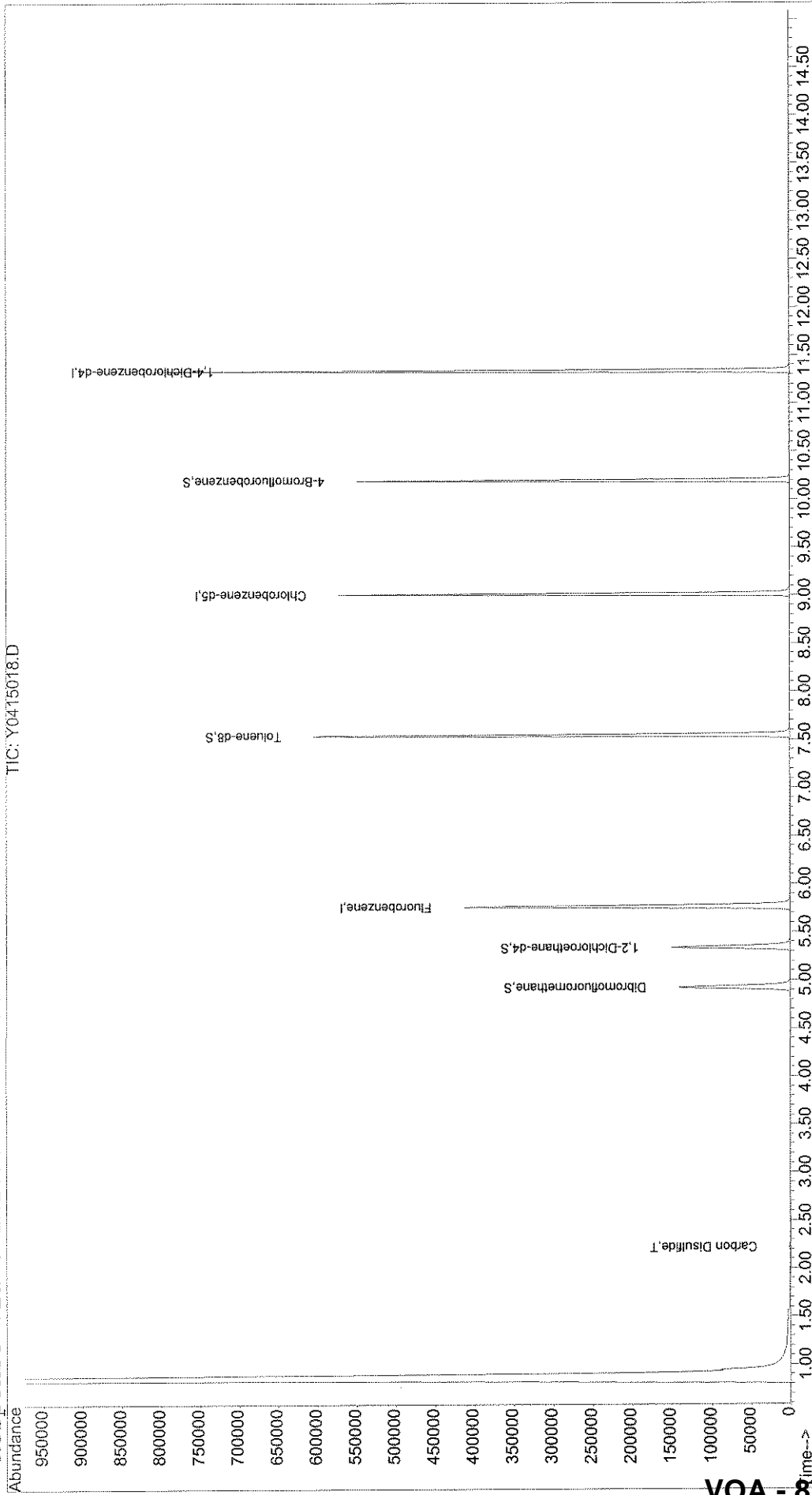
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415018.D Vial: 31  
Acq On : 15 Apr 2007 18:15 Operator: LH  
Sample : JPL31-008 Inst : Yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:45 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415018.D  
 Acq On : 15 Apr 2007 18:15  
 Sample : JPL31-008  
 Misc : 5mL+IS/SS #1

Vial: 31  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:45 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	374097	50.00	ug/l	0.00 110.11%
50) Chlorobenzene-d5	9.01	82	149144	50.00	ug/l	0.00 92.53%
70) 1,4-Dichlorobenzene-d4	11.34	152	179631	50.00	ug/l	0.00 79.71%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	101810	44.98	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	89.96%	
37) 1,2-Dichloroethane-d4	5.34	65	113979	47.47	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	94.94%	
51) Toluene-d8	7.54	98	351676	59.42	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	118.84%	
72) 4-Bromofluorobenzene	10.20	95	140470	54.13	ug/l	0.00

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	2799	0.37 ug/l	100	
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.		
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0415018.D 8260B.M Mon Apr 16 13:45:54 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415018.D  
 Acq On : 15 Apr 2007 18:15  
 Sample : JPL31-008  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:45 2007

Vial: 31  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.40	78	53		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	78		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	422		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	146		N.D.	
65) m,p-Xylene	9.29	106	75		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415018.D 8260B.M Mon Apr 16 13:45:54 2007

Quantitation Report

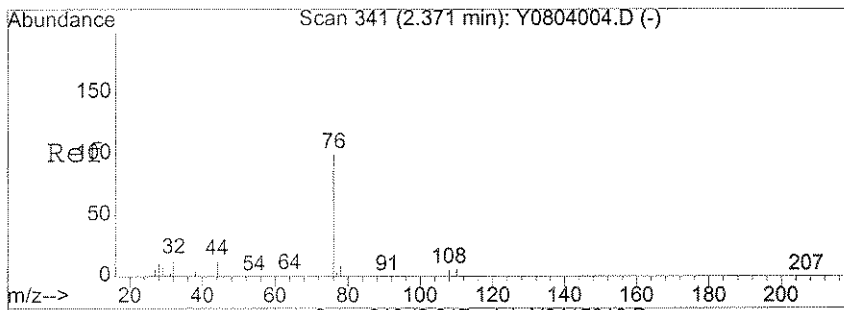
Data File : X:\MSVOA\YODA\041507\Y0415018.D  
 Acq On : 15 Apr 2007 18:15  
 Sample : JPL31-008  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:45 2007

Vial: 31  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

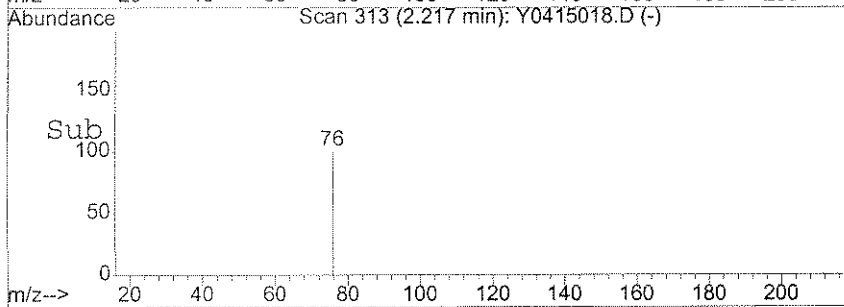
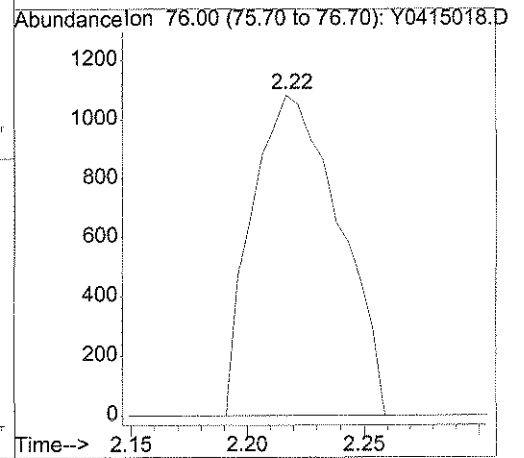
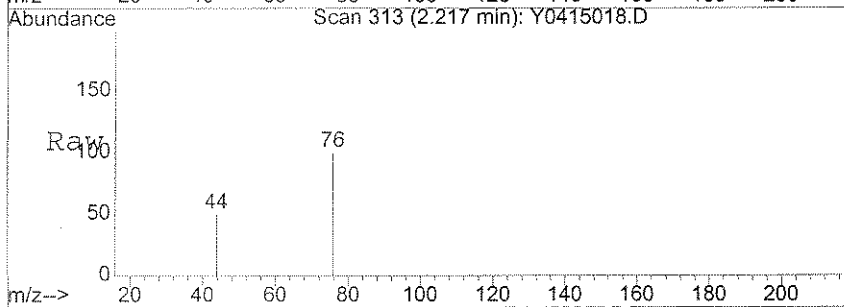
Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.19	105	54		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	122		N.D.	
78) 4-Chlorotoluene	10.47	91	122		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.02	105	55		N.D.	
82) sec-butylbenzene	11.18	105	56		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.33	119	257		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	236		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



#14  
 Carbon Disulfide  
 Concen: 0.37 ug/l  
 RT: 2.22 min Scan# 313  
 Delta R.T. 0.01 min  
 Lab File: Y0415018.D  
 Acq: 15 Apr 2007 18:15  
 Tgt Ion: 76 Resp: 2799



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 18:40

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 18:40

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



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VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-11-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/03/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 18:40

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

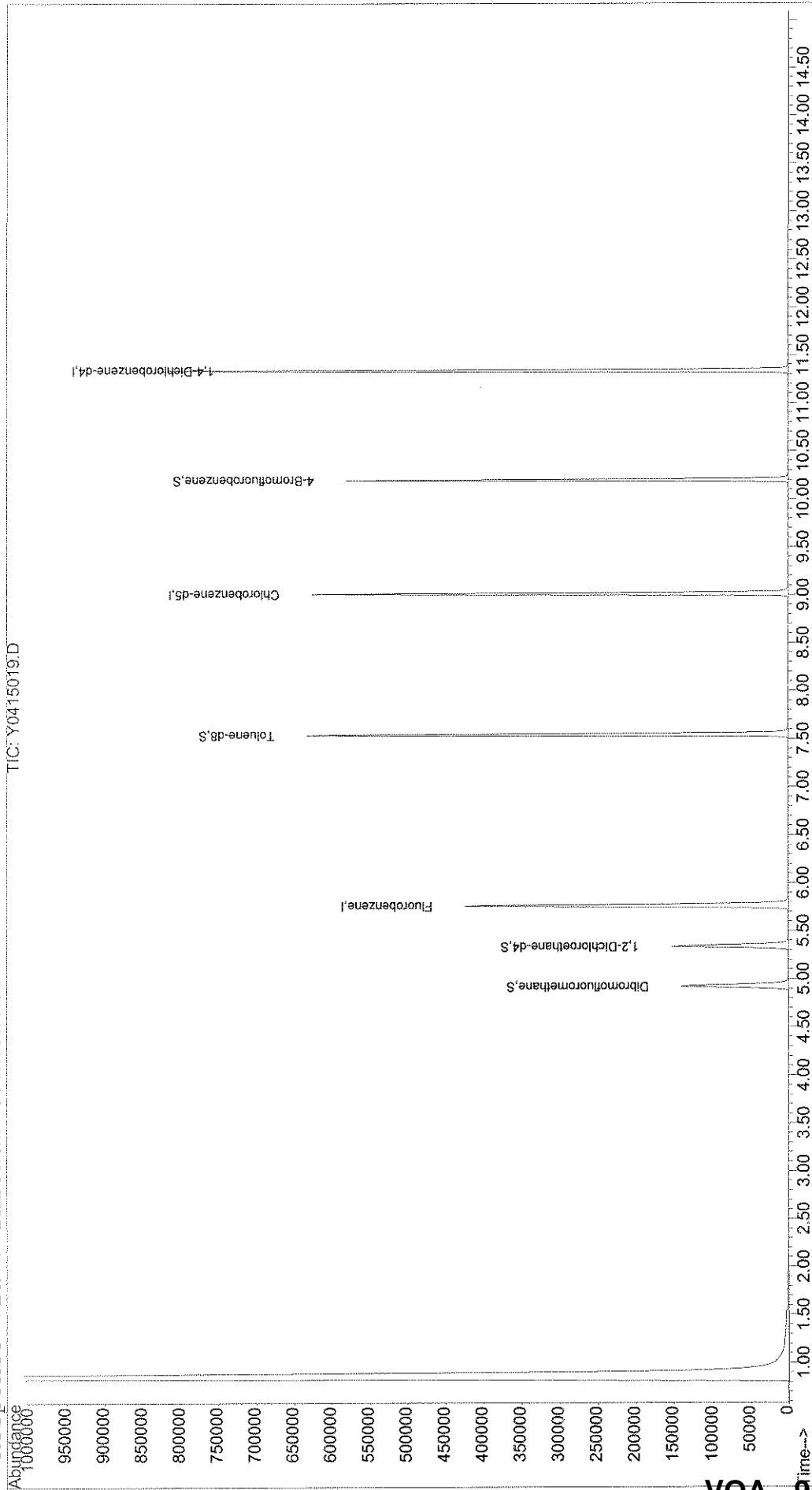
Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415019.D  
Acq On : 15 Apr 2007 18:40  
Sample : JPL31-009  
Misc : 5mL+IS/SS #4  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:46 2007

Vial: 32  
Operator: LH  
Inst : Yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415019.D  
 Acq On : 15 Apr 2007 18:40  
 Sample : JPL31-009  
 Misc : 5mL+IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:46 2007

Vial: 32  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	382611	50.00	ug/l	0.00 112.62%
50) Chlorobenzene-d5	9.01	82	156328	50.00	ug/l	0.00 96.99%
70) 1,4-Dichlorobenzene-d4	11.34	152	182977	50.00	ug/l	0.00 81.20%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	103125	44.55	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	89.10%	
37) 1,2-Dichloroethane-d4	5.34	65	116948	47.62	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	95.24%	
51) Toluene-d8	7.54	98	366214	59.03	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	118.06%	
72) 4-Bromofluorobenzene	10.20	95	145802	55.15	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	2.22	76	541	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		
23) Vinyl acetate	0.00	43	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 Y0415019.D 8260B.M Mon Apr 16 13:46:58 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415019.D  
 Acq On : 15 Apr 2007 18:40  
 Sample : JPL31-009  
 Misc : 5mL+IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:46 2007

Vial: 32  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	53		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	485		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	212		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415019.D 8260B.M Mon Apr 16 13:46:58 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415019.D  
 Acq On : 15 Apr 2007 18:40  
 Sample : JPL31-009  
 Misc : 5mL+IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:46 2007

Vial: 32  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	69		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	65		N.D.	
82) sec-butylbenzene	11.18	105	65		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	208		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	253		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	54		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-011  
 Lab File ID: Y0415020.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/15/2007 19:05  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 19:05

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-011  
 Lab File ID: Y0415020.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/15/2007 19:05  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

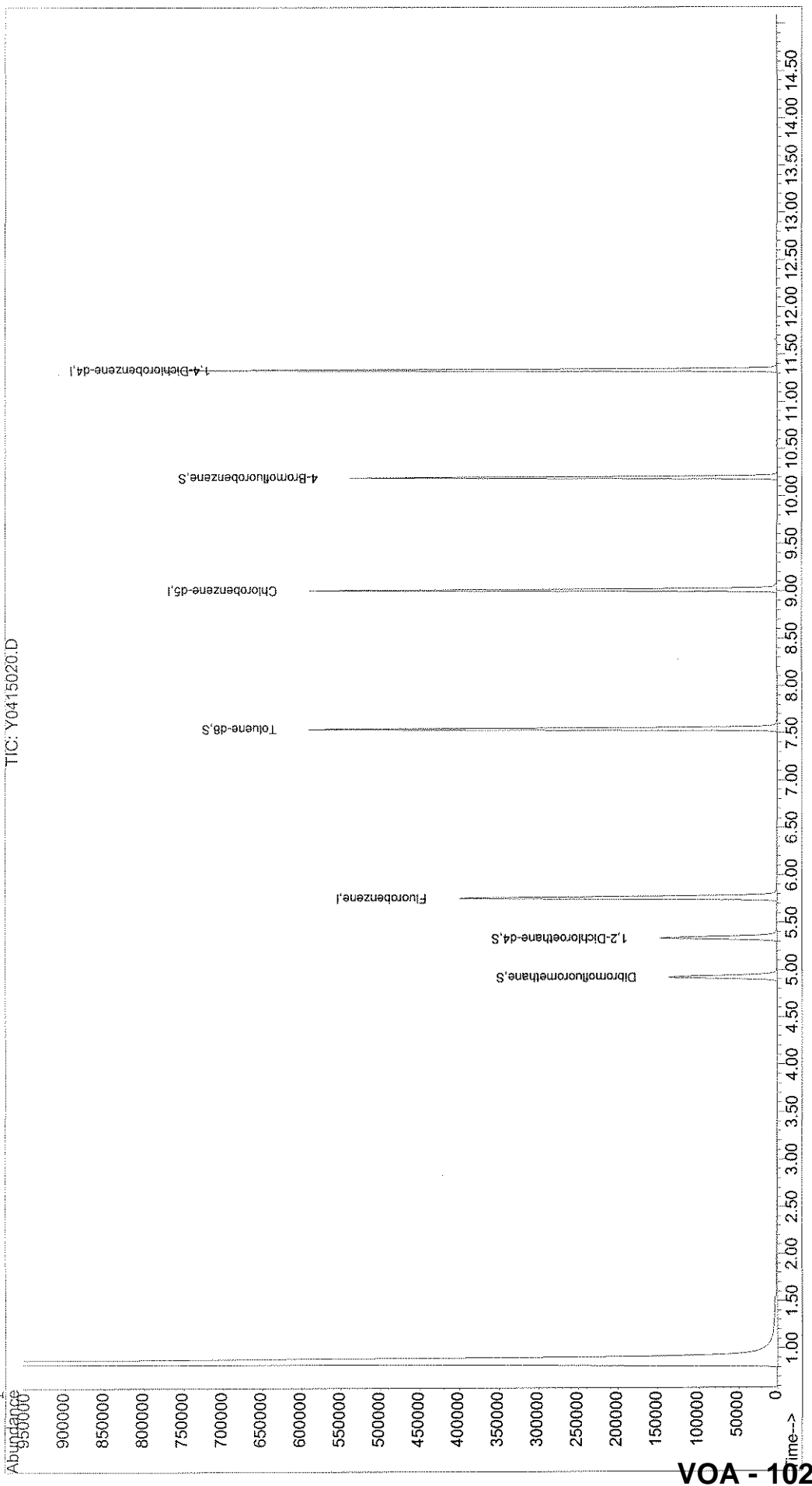
Comments:



Data File : X:\MSVOA\YODA\041507\Y0415020.D  
Acq On : 15 Apr 2007 19:05  
Sample : JPL31-011  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:48 2007

Vial: 33  
Operator: LH  
Inst : yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415020.D  
 Acq On : 15 Apr 2007 19:05  
 Sample : JPL31-011  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:48 2007

Vial: 33  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
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 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	5.76	96	370267	50.00	ug/l	0.00 108.99%
50) Chlorobenzene-d5	9.01	82	150647	50.00	ug/l	0.00 93.47%
70) 1,4-Dichlorobenzene-d4	11.34	152	173859	50.00	ug/l	0.00 77.15%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	100839	45.01	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	90.02%
37) 1,2-Dichloroethane-d4	5.34	65	113515	47.76	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.52%
51) Toluene-d8	7.55	98	352820	59.02	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	118.04%
72) 4-Bromofluorobenzene	10.20	95	138627	55.19	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.21	76	536		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415020.D 8260B.M Mon Apr 16 13:48:29 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415020.D  
 Acq On : 15 Apr 2007 19:05  
 Sample : JPL31-011  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:48 2007

Vial: 33  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	110		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	57		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	369		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	53		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415020.D 8260B.M Mon Apr 16 13:48:30 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415020.D  
 Acq On : 15 Apr 2007 19:05  
 Sample : JPL31-011  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:48 2007

Vial: 33  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.18	105	60		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	65		N.D.	
78) 4-Chlorotoluene	10.47	91	65		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	80		N.D.	
82) sec-butylbenzene	11.18	105	80		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	57		N.D.	
84) 4-Isopropyltoluene	11.34	119	252		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	68		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	203		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:31

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-012  
 Lab File ID: Y0417024.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 16:31  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	0.50		U
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	0.50		U
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-012  
 Lab File ID: Y0417024.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 16:31  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

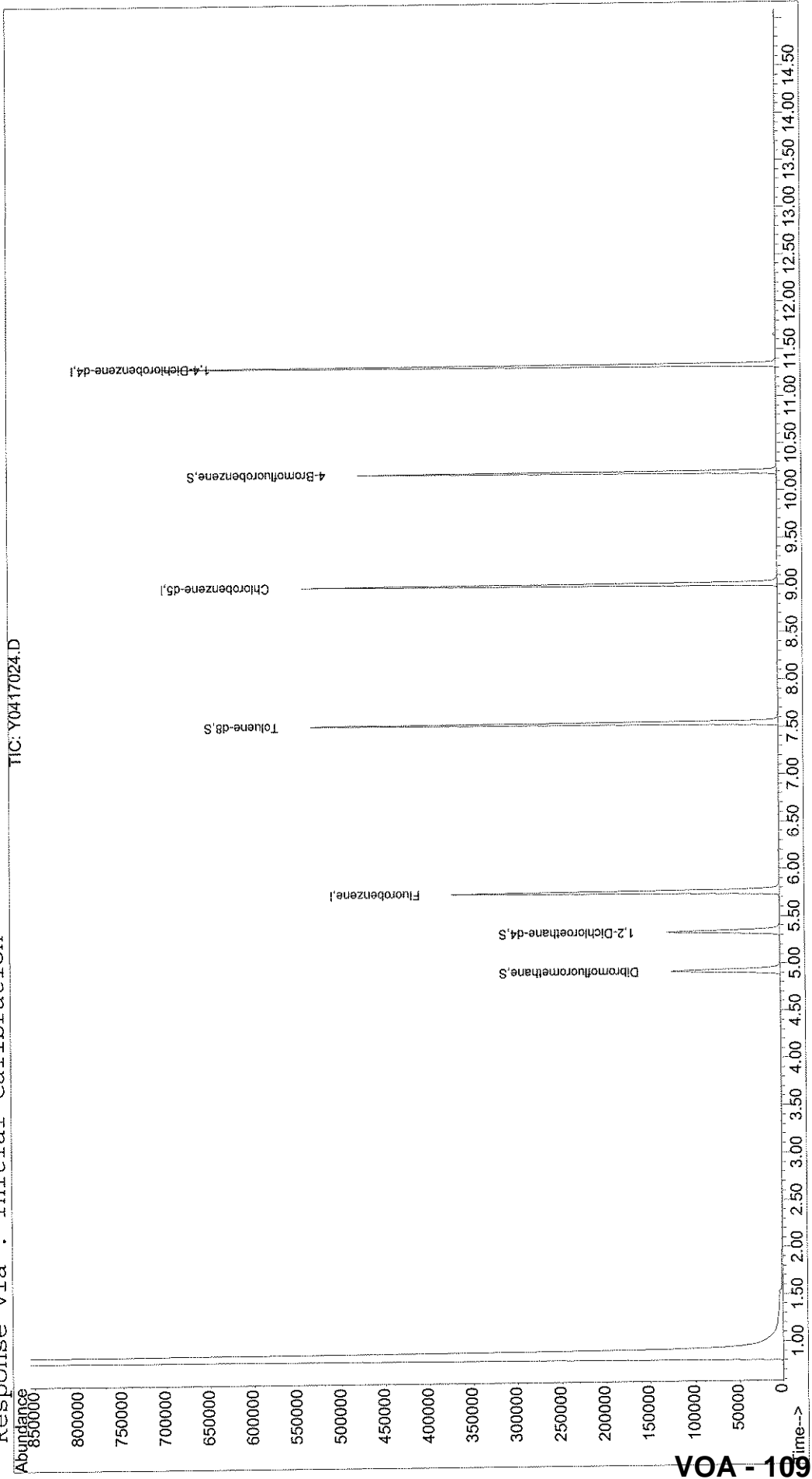
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417024.D  
Acq On : 17 Apr 2007 16:31  
Sample : JPL31-012  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:23 2007  
Vial: 30  
Operator: LNH  
Inst : yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 109



Data File : X:\MSVOA\YODA\041707\Y0417024.D  
 Acq On : 17 Apr 2007 16:31  
 Sample : JPL31-012  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:23 2007

Vial: 30  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	5.76	96	325926	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	135395	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	157820	50.00	ug/l	0.00

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	89926	45.60	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	91.20%
37) 1,2-Dichloroethane-d4	5.34	65	99187	47.41	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	94.82%
51) Toluene-d8	7.54	98	311493	57.97	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	115.94%
72) 4-Bromofluorobenzene	10.20	95	124823	54.74	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.09	101	284		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.21	76	323		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	2.39	40	64		N.D.	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	3.32	63	407		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417024.D 8260B.M Wed Apr 18 07:23:36 2007

Data File : X:\MSVOA\YODA\041707\Y0417024.D  
 Acq On : 17 Apr 2007 16:31  
 Sample : JPL31-012  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:23 2007

Vial: 30  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.70	83	939		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	490		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	442		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	376		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	232		N.D.	
65) m,p-Xylene	9.29	106	136		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417024.D 8260B.M Wed Apr 18 07:23:37 2007

Data File : X:\MSVOA\YODA\041707\Y0417024.D  
Acq On : 17 Apr 2007 16:31  
Sample : JPL31-012  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:23 2007

Vial: 30  
Operator: LNH  
Inst : yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration  
DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.46	91	137		N.D.	
78) 4-Chlorotoluene	10.46	91	137		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.19	105	54		N.D.	
82) sec-butylbenzene	11.19	105	54		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	426		N.D.	
84) 4-Isopropyltoluene	11.33	119	242		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	170		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	253		N.D.	
87) n-Butylbenzene	11.74	91	269		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.31	180	135		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	221		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:55

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
75-71-8	Dichlorodifluoromethane	0.50		U
74-87-3	Chloromethane	0.50		U
75-01-4	Vinyl chloride	0.50		U
74-83-9	Bromomethane	0.50		U
75-00-3	Chloroethane	0.50		U
75-69-4	Trichlorofluoromethane	0.50		U
75-35-4	1,1-Dichloroethene	0.50		U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50		U
75-09-2	Methylene chloride	0.50		U
1634-04-4	Methyl tert-butyl ether	0.50		U
156-60-5	trans-1,2-Dichloroethene	0.50		U
75-34-3	1,1-Dichloroethane	0.50		U
594-20-7	2,2-Dichloropropane	0.50		U
156-59-2	cis-1,2-Dichloroethene	0.50		U
78-93-3	2-Butanone	5.0		U
74-97-5	Bromochloromethane	0.50		U
67-66-3	Chloroform	0.50		U
71-55-6	1,1,1-Trichloroethane	0.50		U
56-23-5	Carbon tetrachloride	0.50		U
563-58-6	1,1-Dichloropropene	0.50		U
71-43-2	Benzene	0.50		U
107-06-2	1,2-Dichloroethane	0.50		U
79-01-6	Trichloroethene	1.4		
78-87-5	1,2-Dichloropropane	0.50		U
74-95-3	Dibromomethane	0.50		U
75-27-4	Bromodichloromethane	0.50		U
10061-01-	cis-1,3-Dichloropropene	0.50		U
108-10-1	4-Methyl-2-pentanone	5.0		U
108-88-3	Toluene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-013  
 Lab File ID: Y0417025.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 16:55  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.61	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-23-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:55

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

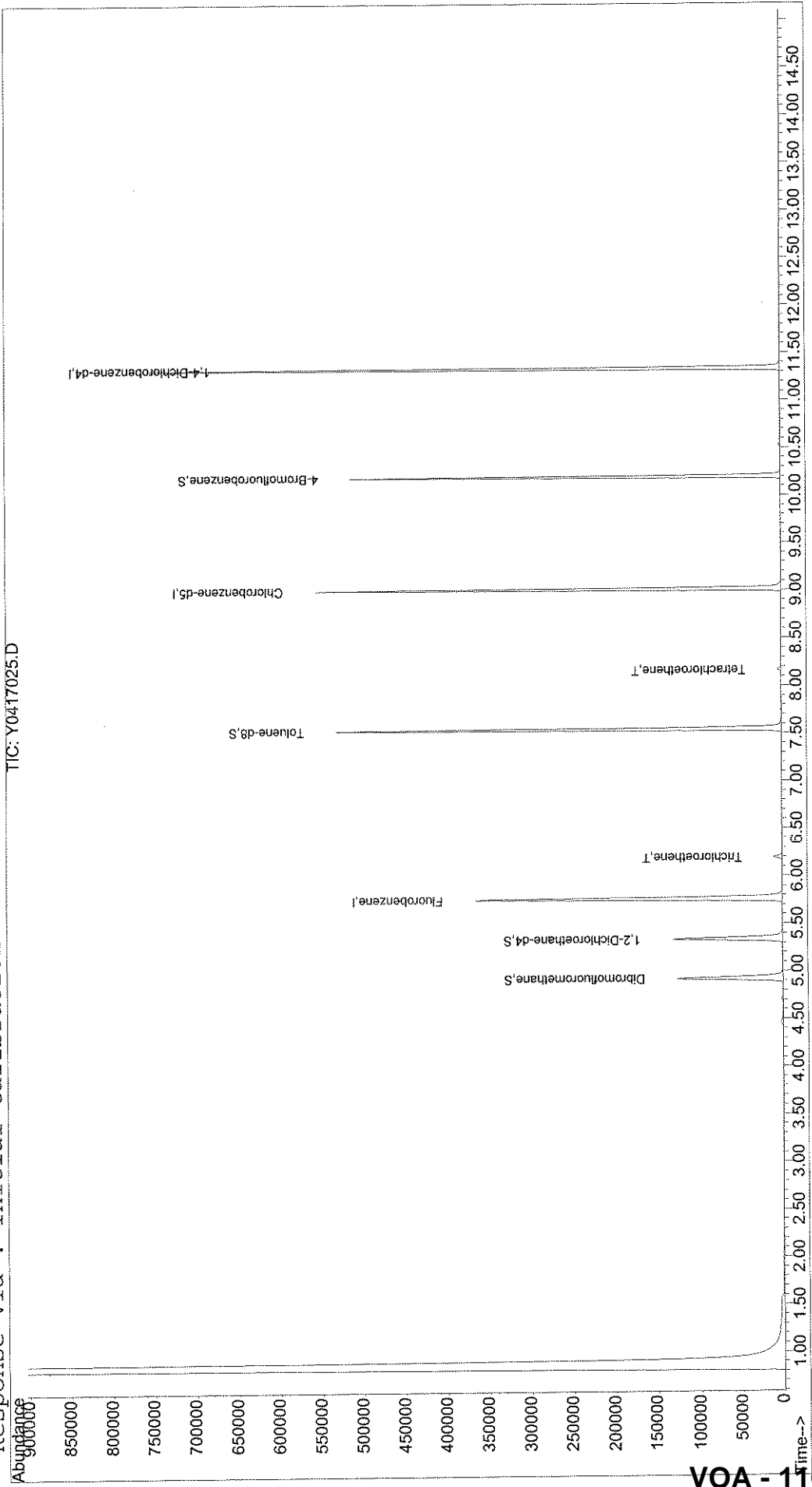
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417025.D  
Acq On : 17 Apr 2007 16:55  
Sample : JPL31-013  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:24 2007  
Vial: 31  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Data File : X:\MSVOA\YODA\041707\Y0417025.D  
 Acq On : 17 Apr 2007 16:55  
 Sample : JPL31-013  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:24 2007

Vial: 31  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	5.76	96	326258	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	136714	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	166662	50.00	ug/l	0.00

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	90781	45.99	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	91.98%		
37) 1,2-Dichloroethane-d4	5.34	65	100827	48.15	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	96.30%		
51) Toluene-d8	7.54	98	312715	57.64	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	115.28%		
72) 4-Bromofluorobenzene	10.20	95	128620	53.42	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.09	101	381	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	173	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	3.33	63	688	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	

(#) = qualifier out of range (m) = manual integration



Data File : X:\MSVOA\YODA\041707\Y0417025.D  
 Acq On : 17 Apr 2007 16:55  
 Sample : JPL31-013  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:24 2007

Vial: 31  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.70	83	1687	N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	0.00	117	0	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	5.40	78	63	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	6.20	130	3236	1.41 ug/l ✓	92
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	6.78	83	110	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	7.62	92	63	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	8.17	166	1160	0.61 ug/l # ✓	90
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	0.00	112	0	N.D.	
62) 1-Chlorohexane	9.02	91	310	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.29	91	208	N.D.	
65) m,p-Xylene	0.00	106	0	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	0.00	104	0	N.D.	
68) Bromoform	0.00	173	0	N.D.	
69) Isopropylbenzene	10.20	105	59	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	0.00	156	0	N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0	N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417025.D 8260B.M Wed Apr 18 07:24:17 2007

Data File : X:\MSVOA\YODA\041707\Y0417025.D

Vial: 31

Acq On : 17 Apr 2007 16:55

Operator: LNH

Sample : JPL31-013

Inst : yoda

Misc : 5mL+IS/SS #1

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 8260B.RES

Quant Time: Apr 18 7:24 2007

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

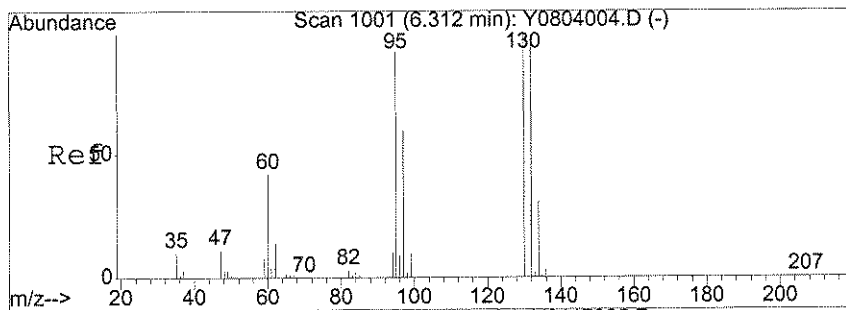
Response via : Initial Calibration

DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.47	91	59		N.D.	
78) 4-Chlorotoluene	10.47	91	59		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.18	105	55		N.D.	
82) sec-butylbenzene	11.18	105	55		N.D.	
83) 1,3-Dichlorobenzene	11.36	146	129		N.D.	
84) 4-Isopropyltoluene	11.34	119	265		N.D.	
85) 1,4-Dichlorobenzene	11.36	146	129		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	87		N.D.	
87) n-Butylbenzene	11.74	91	258		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	138		N.D.	

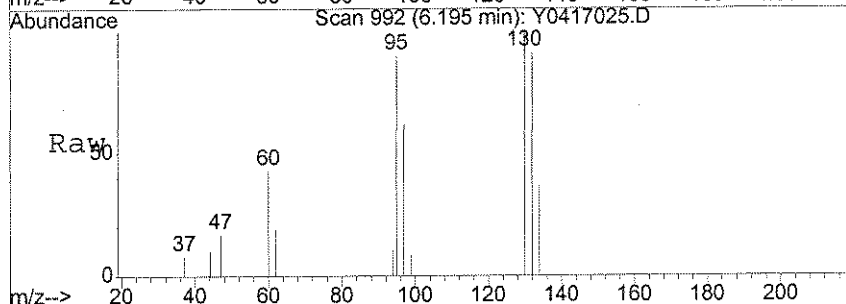
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(#) = qualifier out of range (m) = manual integration

Y0417025.D 8260B.M Wed Apr 18 07:24:18 2007

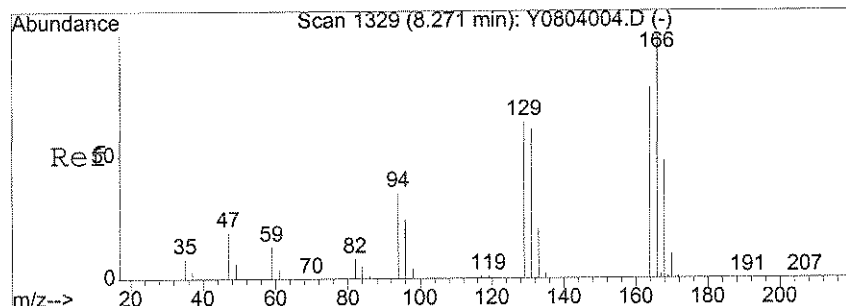
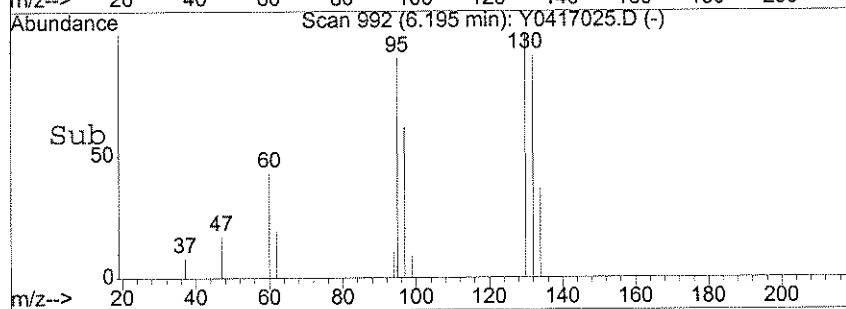
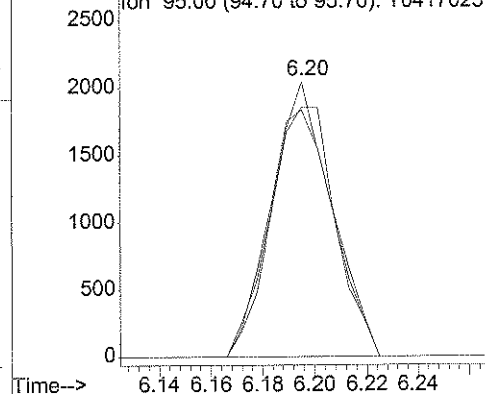


#41  
 Trichloroethene  
 Concen: 1.41 ug/l  
 RT: 6.20 min Scan# 992  
 Delta R.T. 0.01 min  
 Lab File: Y0417025.D  
 Acq: 17 Apr 2007 16:55

Tgt Ion	Resp	Lower	Upper
130	3236		
130	100		
132	100.4	75.0	115.0
95	98.4	69.4	109.4

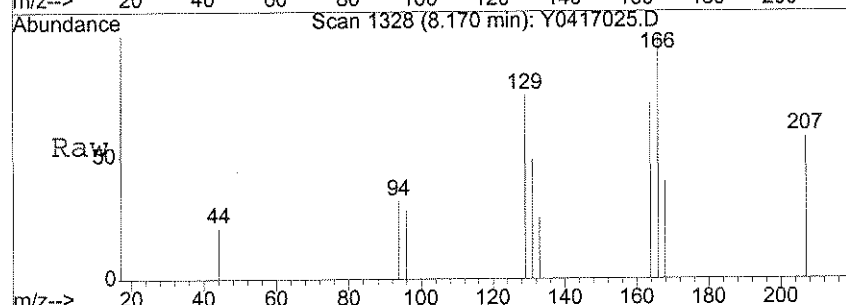


Abundance Ion 130.00 (129.70 to 130.70): Y041702  
 Ion 132.00 (131.70 to 132.70): Y041702  
 Ion 95.00 (94.70 to 95.70): Y0417025.D

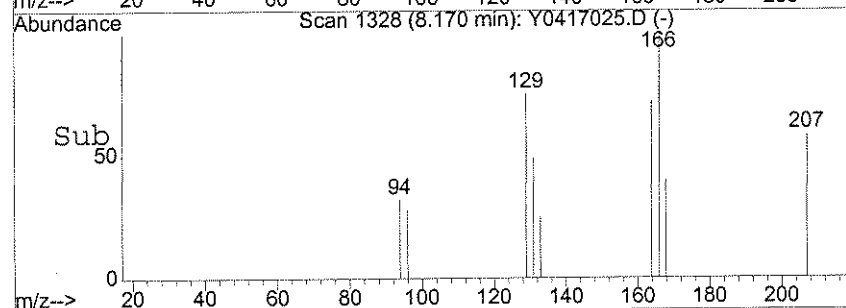
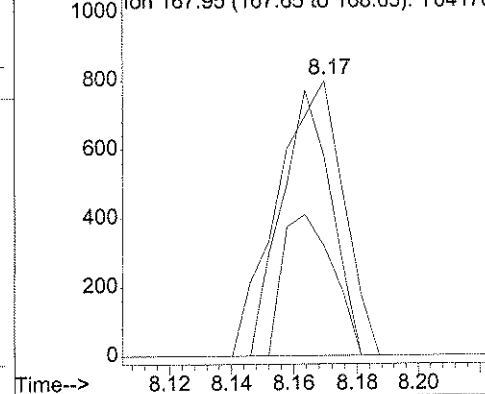


#56  
 Tetrachloroethene  
 Concen: 0.61 ug/l  
 RT: 8.17 min Scan# 1328  
 Delta R.T. 0.01 min  
 Lab File: Y0417025.D  
 Acq: 17 Apr 2007 16:55

Tgt Ion	Resp	Lower	Upper
166	1160		
166	100		
164	73.1	63.3	94.9
168	39.1	39.6	59.4#



Abundance Ion 165.95 (165.65 to 166.65): Y041702  
 Ion 163.95 (163.65 to 164.65): Y041702  
 Ion 167.95 (167.65 to 168.65): Y041702



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:20

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-014  
 Lab File ID: Y0417026.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 17:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,1,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-014  
 Lab File ID: Y0417026.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 17:20  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

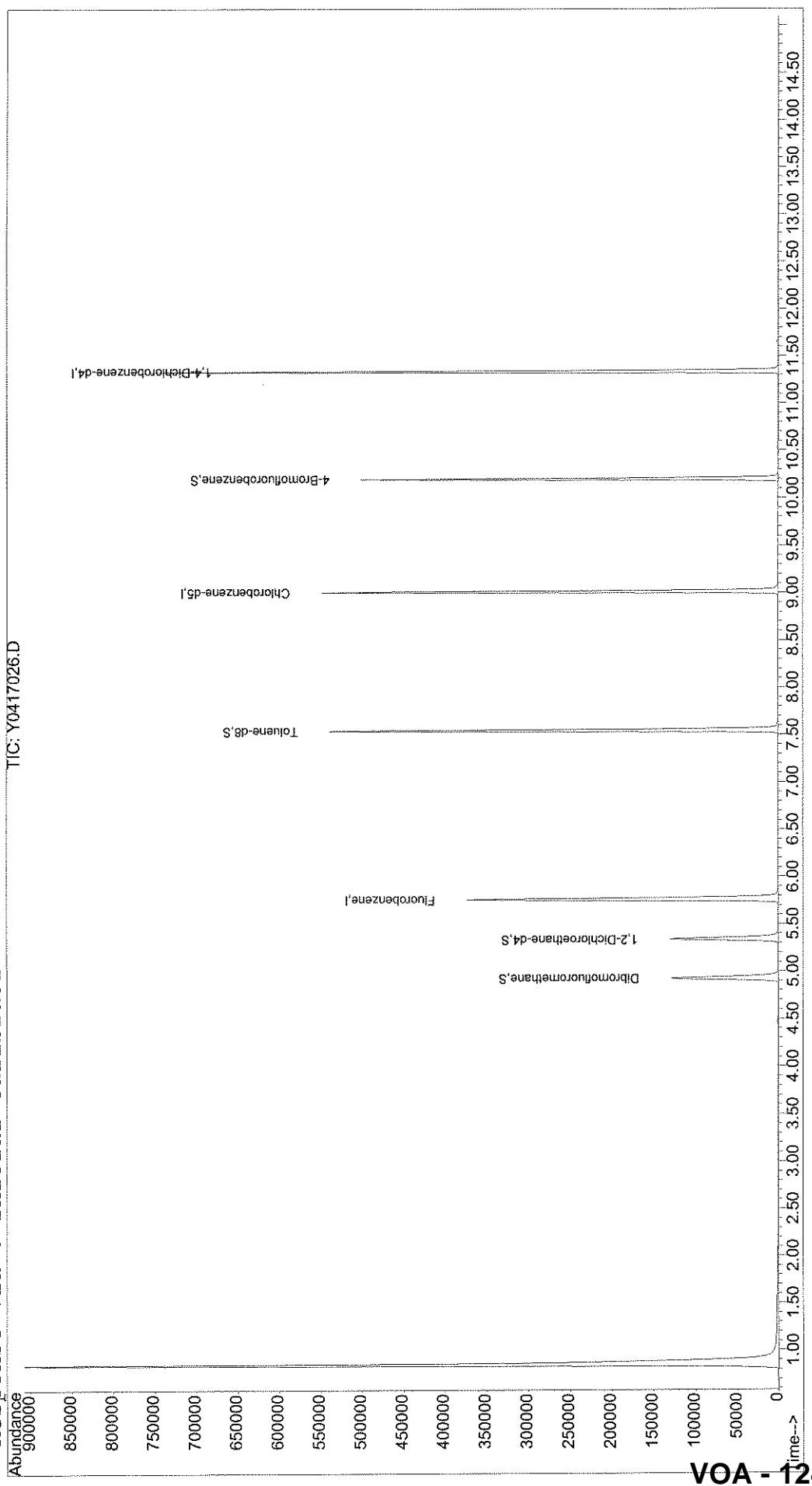
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417026.D Vial: 32  
Acq On : 17 Apr 2007 17:20 Operator: LNH  
Sample : JPL31-014 Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:25 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Data File : X:\MSVOA\YODA\041707\Y0417026.D

Vial: 32

Acq On : 17 Apr 2007 17:20

Operator: LNH

Sample : JPL31-014

Inst : yoda

Misc : 5mL+IS/SS #2

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 8260B.RES

Quant Time: Apr 18 7:25 2007

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

Response via : Initial Calibration

DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	5.76	96	331284	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	138350	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	163512	50.00	ug/l	0.00

## System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	92271	46.04	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	92.08%
37) 1,2-Dichloroethane-d4	5.34	65	101989	47.96	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.92%
51) Toluene-d8	7.54	98	314863	57.35	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	114.70%
72) 4-Bromofluorobenzene	10.20	95	127680	54.05	ug/l	0.00

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.09	101	169		N.D.	
11) Acetone	0.00	43	0		N.D.	d
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	0.00	76	0		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	

(#) = qualifier out of range (m) = manual integration

Y0417026.D 8260B.M Wed Apr 18 07:25:28 2007



Data File : X:\MSVOA\YODA\041707\Y0417026.D  
 Acq On : 17 Apr 2007 17:20  
 Sample : JPL31-014  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:25 2007

Vial: 32  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	140		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	437		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.01	91	437		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417026.D 8260B.M Wed Apr 18 07:25:28 2007

Data File : X:\MSVOA\YODA\041707\Y0417026.D  
Acq On : 17 Apr 2007 17:20  
Sample : JPL31-014  
Misc : 5mL+IS/SS #2  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:25 2007

Vial: 32  
Operator: LNH  
Inst : yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration  
DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.46	91	57		N.D.	
78) 4-Chlorotoluene	10.46	91	57		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	219		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	229		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-015  
 Lab File ID: Y0415011.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/15/2007 15:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-015

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 15:25

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-015

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/15/2007 15:25

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

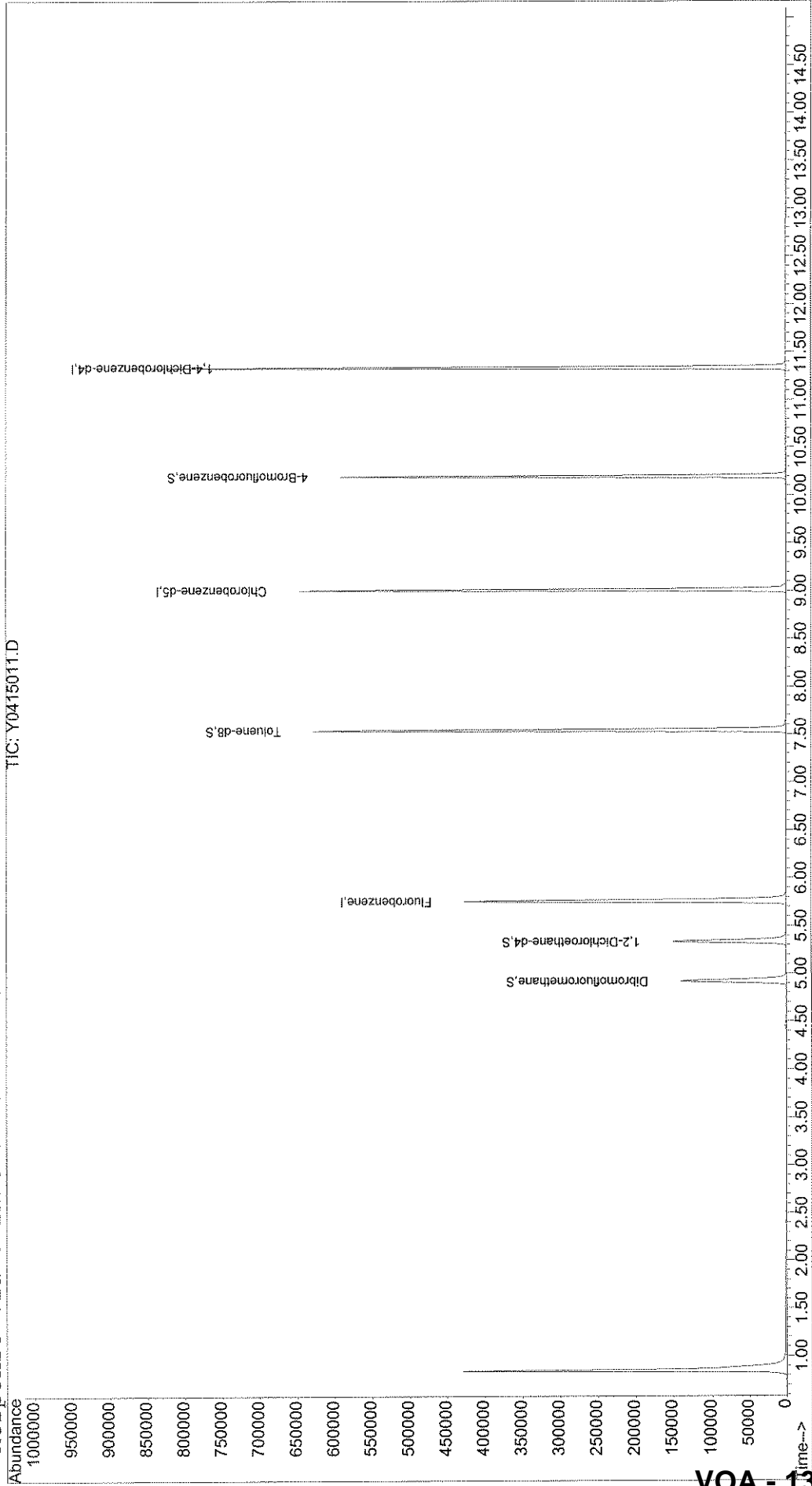
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D Vial: 24  
Acq On : 15 Apr 2007 15:25 Operator: LH  
Sample : JPL31-015 TB Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 16 13:36 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 131

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D  
 Acq On : 15 Apr 2007 15:25  
 Sample : JPL31-015 TB  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:36 2007

Vial: 24  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : X:\MSVOA\YODA\011707\Y0117007.D (17 Jan 2007 11:52)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	390484	50.00	ug/l	0.00 114.94%
50) Chlorobenzene-d5	9.01	82	160220	50.00	ug/l	0.00 99.41%
70) 1,4-Dichlorobenzene-d4	11.34	152	186356	50.00	ug/l	0.00 82.70%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	104152	44.09	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	88.18%	
37) 1,2-Dichloroethane-d4	5.34	65	117159	46.74	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	93.48%	
51) Toluene-d8	7.54	98	371079	58.36	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	116.72%	
72) 4-Bromofluorobenzene	10.20	95	153412	56.98	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	d
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	282		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	2.52	84	340		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415011.D 8260B.M Mon Apr 16 13:36:54 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D  
 Acq On : 15 Apr 2007 15:25  
 Sample : JPL31-015 TB  
 Misc : 5mL+IS/SS #2

Vial: 24  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:36 2007

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	55		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	6.38	83	60		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	109		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	68		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	539		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.16	91	142		N.D.	
65) m,p-Xylene	9.29	106	69		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415011.D 8260B.M Mon Apr 16 13:36:54 2007



Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D  
 Acq On : 15 Apr 2007 15:25  
 Sample : JPL31-015 TB  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:36 2007

Vial: 24  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	5.38	78	55		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	6.38	83	60		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	109		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.16	166	68		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	539		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.16	91	142		N.D.	
65) m,p-Xylene	9.29	106	69		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0415011.D 8260B.M Mon Apr 16 13:36:54 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D  
 Acq On : 15 Apr 2007 15:25  
 Sample : JPL31-015 TB  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 16 13:36 2007

Vial: 24  
 Operator: LH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.06	105	60		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.53	91	74		N.D.	
78) 4-Chlorotoluene	10.64	91	362		N.D.	
79) 1,3,5-Trimethylbenzene	10.65	105	150		N.D.	
80) tert-Butylbenzene	10.97	119	57		N.D.	
81) 1,2,4-Trimethylbenzene	11.01	105	237		N.D.	
82) sec-butylbenzene	11.18	105	458		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	204		N.D.	
84) 4-Isopropyltoluene	11.34	119	685		N.D.	
85) 1,4-Dichlorobenzene	11.35	146	258		N.D.	
86) 1,2-Dichlorobenzene	11.72	146	116		N.D.	
87) n-Butylbenzene	11.74	91	762		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	13.30	180	236		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	d
91) Naphthalene	13.54	128	221		N.D.	
92) 1,2,3-Trichlorobenzene	13.79	180	221		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-016  
 Lab File ID: Y0417027.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 17:44  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-016

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417027.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:44

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-016  
 Lab File ID: Y0417027.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 17:44  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

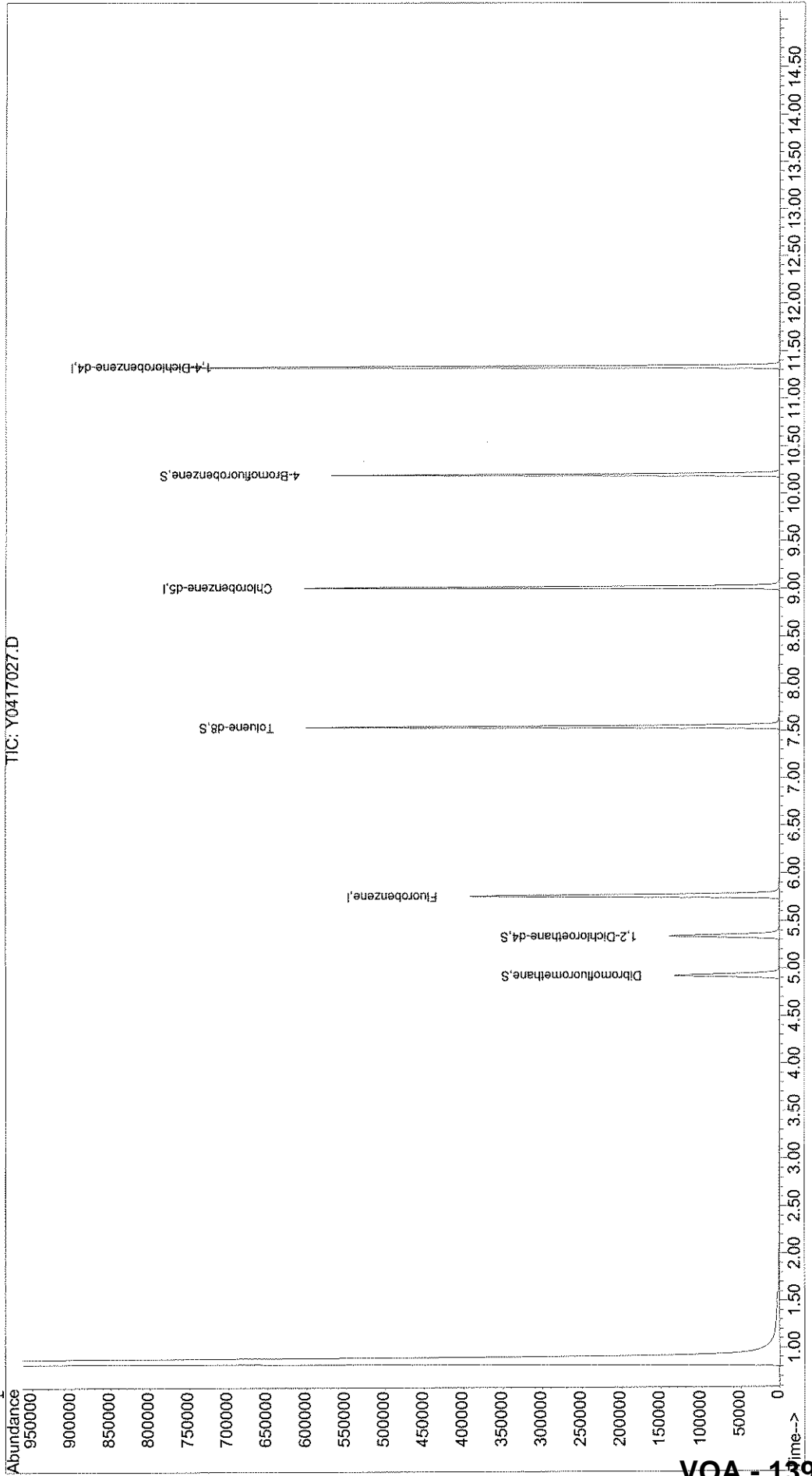
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417027.D  
Acq On : 17 Apr 2007 17:44 Vial: 33  
Sample : JPL31-016 Operator: LNH  
Misc : 5mL+IS/SS #2 Inst : Yoda  
MS Integration Params: rteint.p Multiplr: 1.00  
Quant Time: Apr 18 7:25 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Data File : X:\MSVOA\YODA\041707\Y0417027.D  
 Acq On : 17 Apr 2007 17:44  
 Sample : JPL31-016  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:25 2007

Vial: 33  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	5.76	96	354128	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	152423	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	175918	50.00	ug/l	0.00

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	96866	45.21	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	90.42%
37) 1,2-Dichloroethane-d4	5.34	65	109085	47.99	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.98%
51) Toluene-d8	7.54	98	346833	57.34	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	114.68%
72) 4-Bromofluorobenzene	10.20	95	143944	56.63	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.08	101	83	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	492	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	2.40	40	62	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417027.D 8260B.M Wed Apr 18 07:26:03 2007

Data File : X:\MSVOA\YODA\041707\Y0417027.D

Vial: 33

Acq On : 17 Apr 2007 17:44

Operator: LNH

Sample : JPL31-016

Inst : yoda

Misc : 5mL+IS/SS #2

Multiplr: 1.00

MS Integration Params: rteint.p

Quant Results File: 8260B.RES

Quant Time: Apr 18 7:25 2007

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973Y

Last Update : Tue Mar 20 10:56:50 2007

Response via : Initial Calibration

DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	72		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.02	91	455		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	145		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	0.00	105	0		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	

(#)=qualifier out of range (m)=manual integration

Y0417027.D 8260B.M Wed Apr 18 07:26:04 2007



Data File : X:\MSVOA\YODA\041707\Y0417027.D  
 Acq On : 17 Apr 2007 17:44  
 Sample : JPL31-016  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:25 2007

Vial: 33  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.47	91	64		N.D.	
78) 4-Chlorotoluene	10.47	91	64		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	11.27	146	58		N.D.	
84) 4-Isopropyltoluene	11.33	119	236		N.D.	
85) 1,4-Dichlorobenzene	11.27	146	58		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	110		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:09

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.57	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	1.4	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-017  
 Lab File ID: Y0417028.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 18:09  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-4

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:09

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

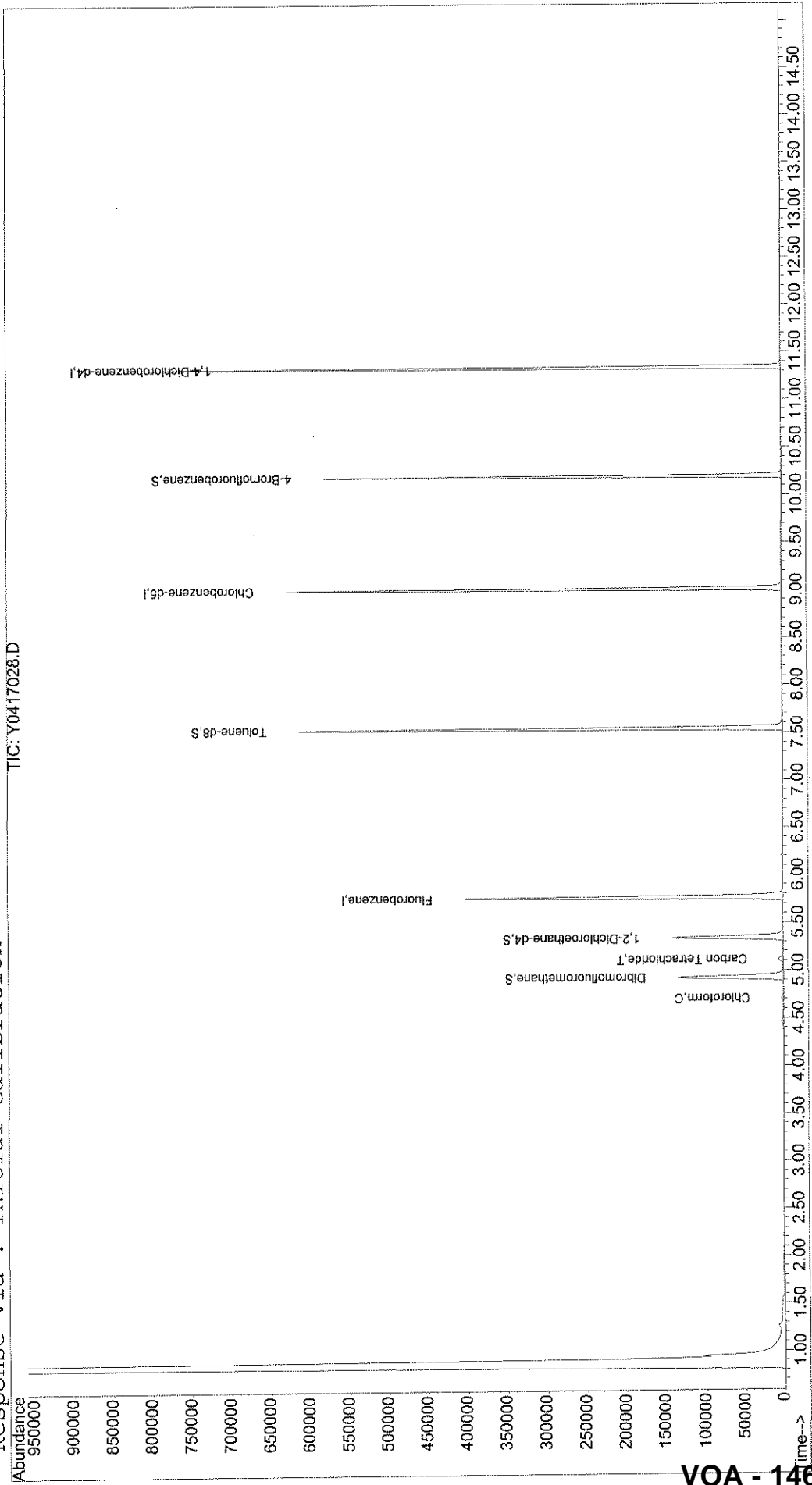
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417028.D  
Acq On : 17 Apr 2007 18:09  
Sample : JPL31-017  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:26 2007  
Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Data File : X:\MSVOA\YODA\041707\Y0417028.D  
 Acq On : 17 Apr 2007 18:09  
 Sample : JPL31-017  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:26 2007

Vial: 34  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	5.76	96	361542	50.00	ug/l	0.00
50) Chlorobenzene-d5	9.01	82	159092	50.00	ug/l	0.00
70) 1,4-Dichlorobenzene-d4	11.34	152	174100	50.00	ug/l	0.00

#### System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	97604	44.62	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	89.24%
37) 1,2-Dichloroethane-d4	5.34	65	110152	47.47	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	94.94%
51) Toluene-d8	7.55	98	356096	56.40	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.80%
72) 4-Bromofluorobenzene	10.20	95	147340	58.58	ug/l	0.00

#### Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.09	101	135	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	810	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	2.35	40	53	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	0.00	43	0	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Y0417028.D 8260B.M Wed Apr 18 07:26:36 2007

Data File : X:\MSVOA\YODA\041707\Y0417028.D  
 Acq On : 17 Apr 2007 18:09  
 Sample : JPL31-017  
 Misc : 5mL+IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:26 2007

Vial: 34  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30) Methacrylonitrile	0.00	41	0	N.D.	
31) Chloroform	4.71	83	2655	0.57 ug/l ✓	99
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	5.11	117	4509	1.39 ug/l ✓	98
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	0.00	78	0	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	6.20	130	952	N.D.	
42) Methylcyclohexane	0.00	83	0	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	0.00	83	0	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	7.61	92	347	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.	
56) Tetrachloroethene	0.00	166	0	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D.	
59) Dibromochloromethane	0.00	129	0	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.04	112	53	N.D.	
62) 1-Chlorohexane	9.01	91	578	N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.17	91	111	N.D.	
65) m,p-Xylene	9.30	106	59	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	9.70	104	138	N.D.	
68) Bromoform	0.00	173	0	N.D.	
69) Isopropylbenzene	10.19	105	78	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	0.00	156	0	N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0	N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0	N.D.	
76) n-Propylbenzene	0.00	120	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417028.D 8260B.M Wed Apr 18 07:26:37 2007

Data File : X:\MSVOA\YODA\041707\Y0417028.D  
Acq On : 17 Apr 2007 18:09  
Sample : JPL31-017  
Misc : 5mL+IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:26 2007

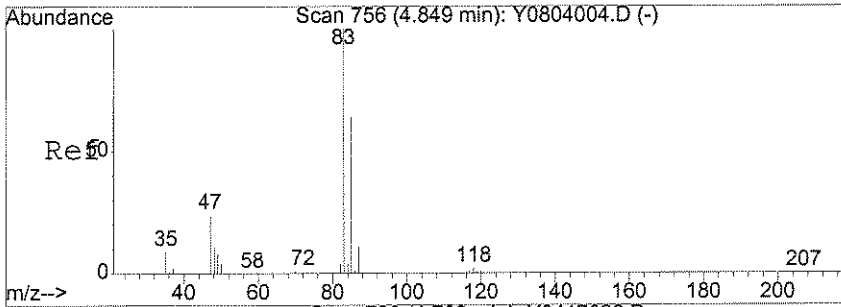
Vial: 34  
Operator: LNH  
Inst : yoda  
Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration  
DataAcq Meth : 8260B

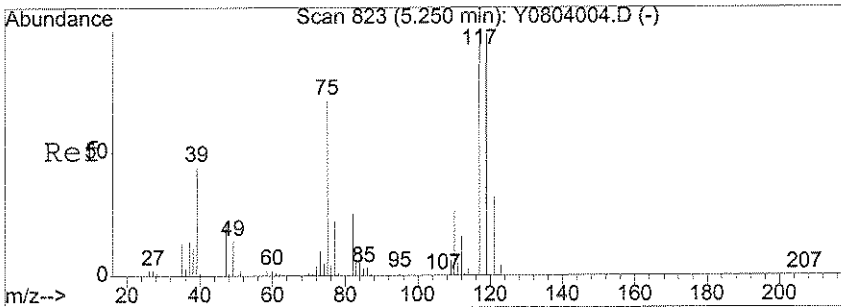
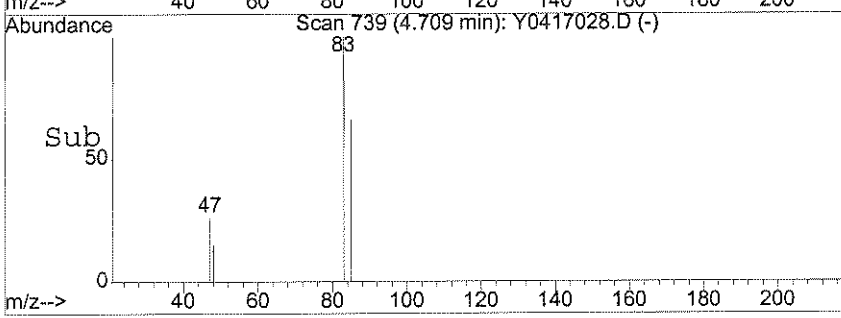
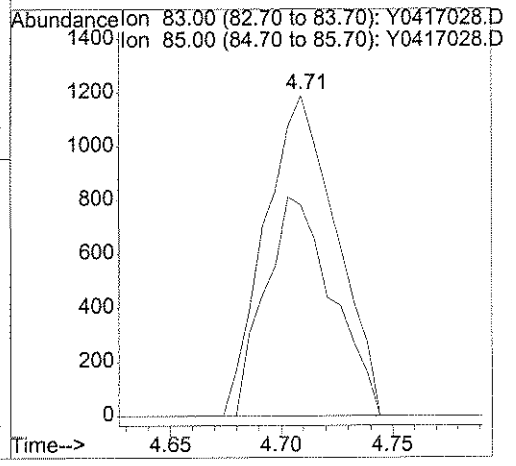
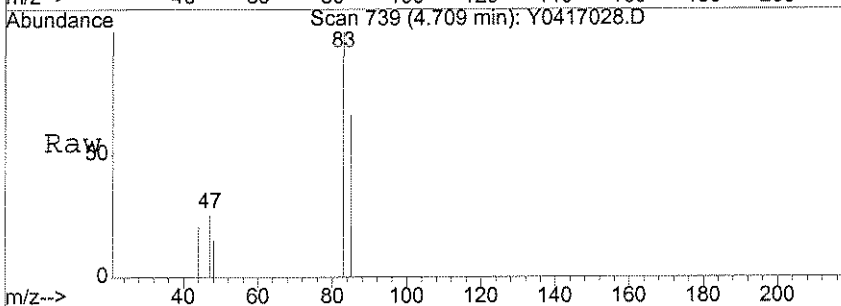
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77) 2-Chlorotoluene	10.46	91	70		N.D.	
78) 4-Chlorotoluene	10.46	91	70		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	136		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	53		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	





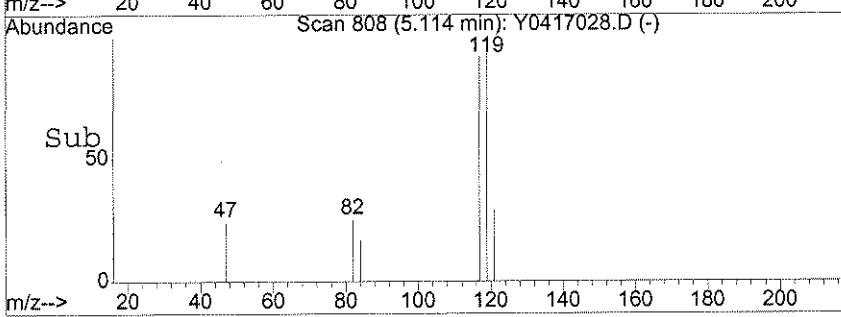
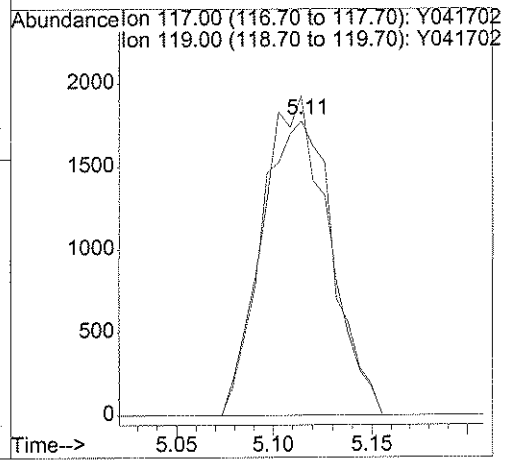
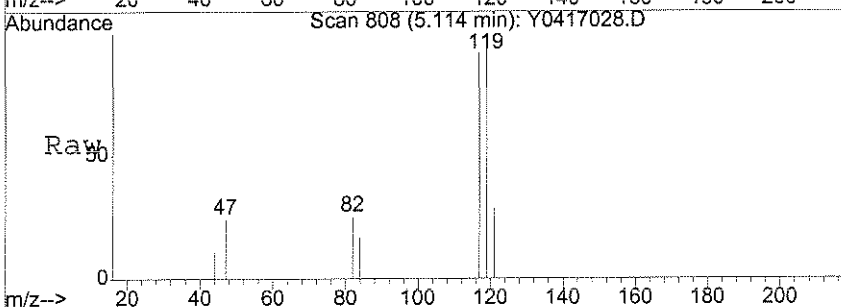
#31  
 Chloroform  
 Concen: 0.57 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0417028.D  
 Acq: 17 Apr 2007 18:09

Tgt Ion:	83	Resp:	2655
Ion Ratio	Lower	Upper	
83	100		
85	64.3	43.3	83.3



#35  
 Carbon Tetrachloride  
 Concen: 1.39 ug/l  
 RT: 5.11 min Scan# 808  
 Delta R.T. 0.01 min  
 Lab File: Y0417028.D  
 Acq: 17 Apr 2007 18:09

Tgt Ion:	117	Resp:	4509
Ion Ratio	Lower	Upper	
117	100		
119	100.5	78.2	118.2



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417029.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:33

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	2.6	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	2.0	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-018  
 Lab File ID: Y0417029.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 18:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-018  
 Lab File ID: Y0417029.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 18:33  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

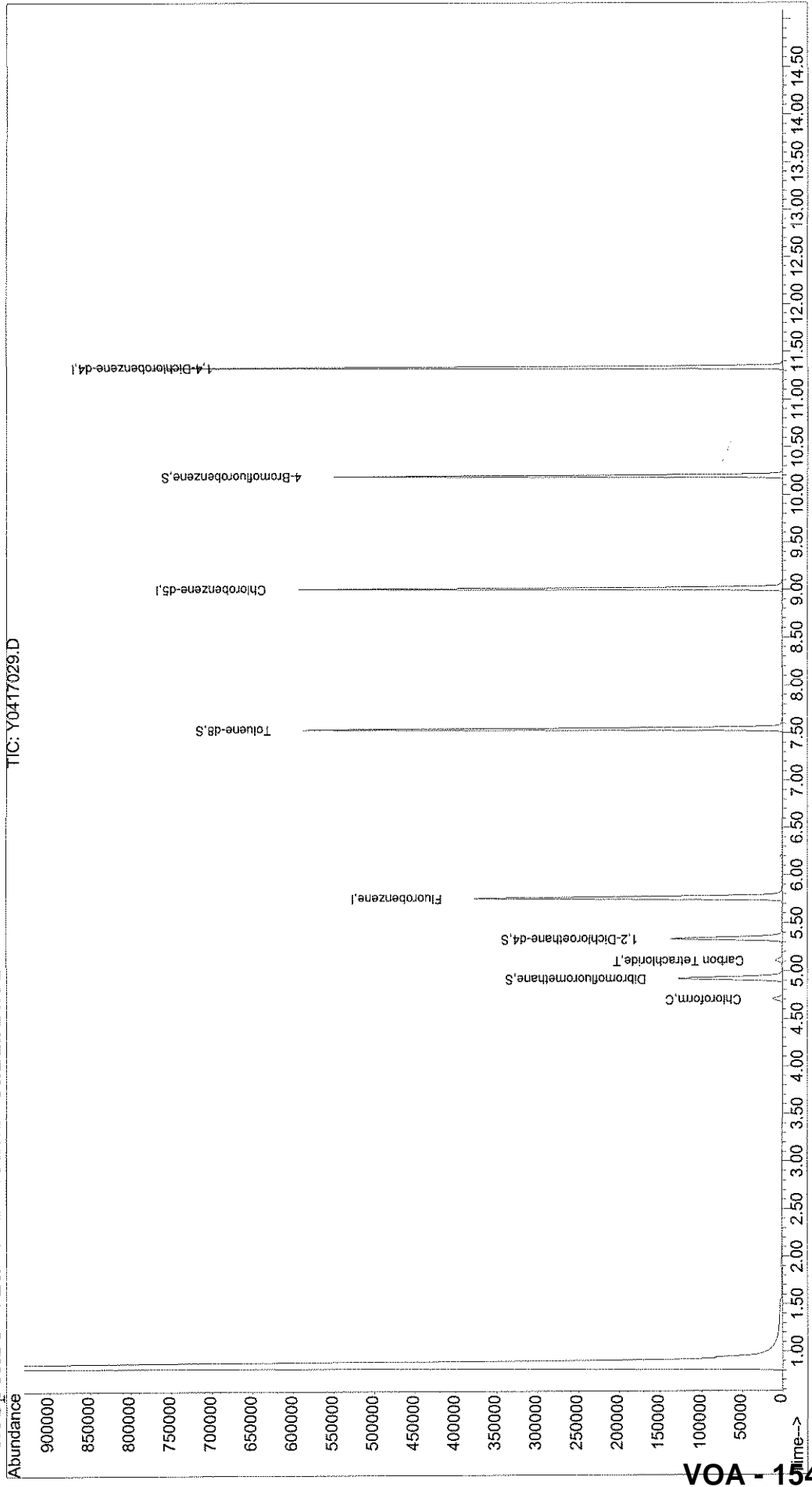
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417029.D Vial: 35  
Acq On : 17 Apr 2007 18:33 Operator: LNH  
Sample : JPL31-018 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:27 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



VOA - 154

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417029.D  
 Acq On : 17 Apr 2007 18:33  
 Sample : JPL31-018  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	342333	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	150817	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	168546	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	94491	45.62	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	91.24%
37) 1,2-Dichloroethane-d4	5.34	65	104887	47.73	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.46%
51) Toluene-d8	7.54	98	336478	56.22	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	112.44%
72) 4-Bromofluorobenzene	10.20	95	141663	58.18	ug/l	0.00

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	2.08	101	137	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	2.22	76	580	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	2.39	40	55	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	
23) Vinyl acetate	0.00	43	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417029.D 8260B.M Wed Apr 18 07:55:28 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417029.D  
 Acq On : 17 Apr 2007 18:33  
 Sample : JPL31-018  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	11585	2.63	ug/l ✓	96
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	5.11	117	6294	2.04	ug/l ✓	98
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	6.20	130	1000		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	7.61	92	60		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.02	91	431		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	247		N.D.	
65) m,p-Xylene	9.29	106	110		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

*LNH 4/19/07*

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417029.D  
 Acq On : 17 Apr 2007 18:33  
 Sample : JPL31-018  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

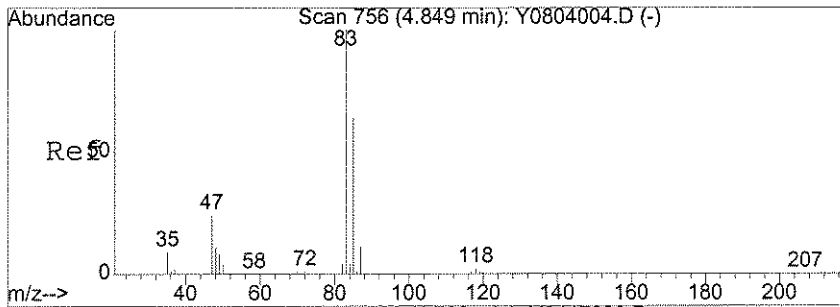
Vial: 35  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

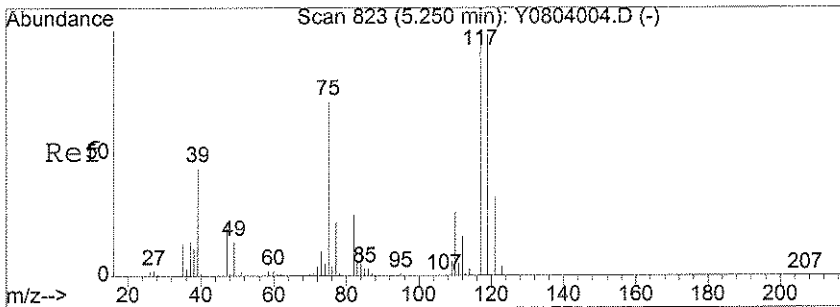
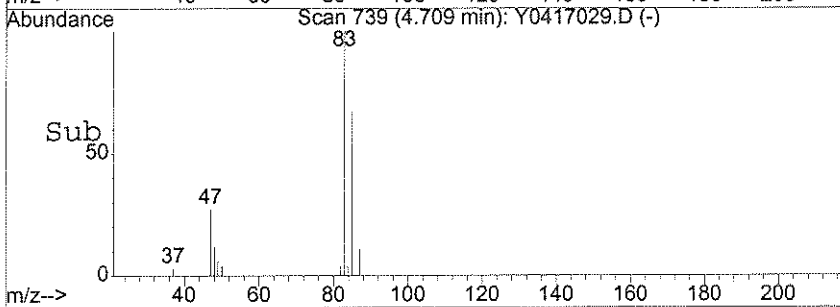
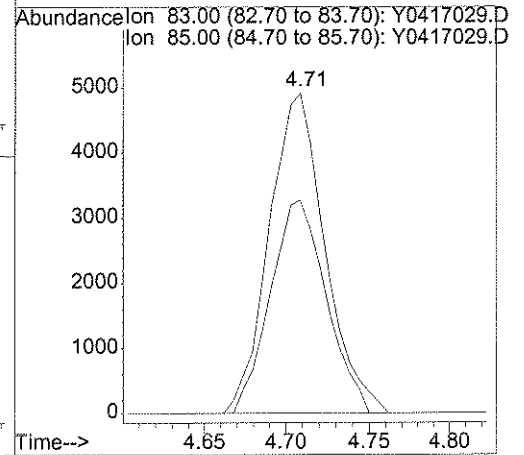
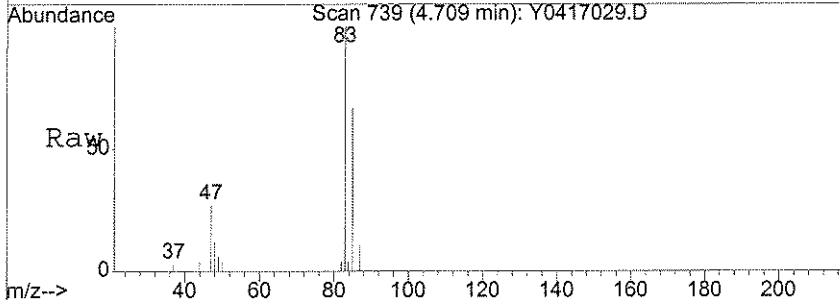
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.21	105	83		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	212		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	128		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	





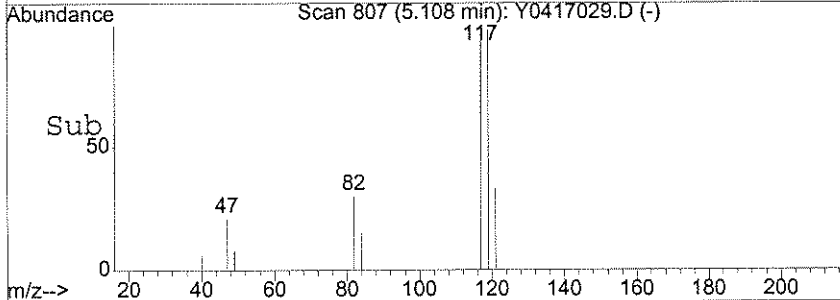
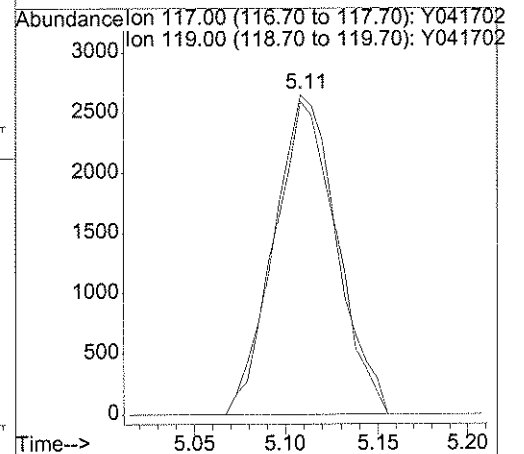
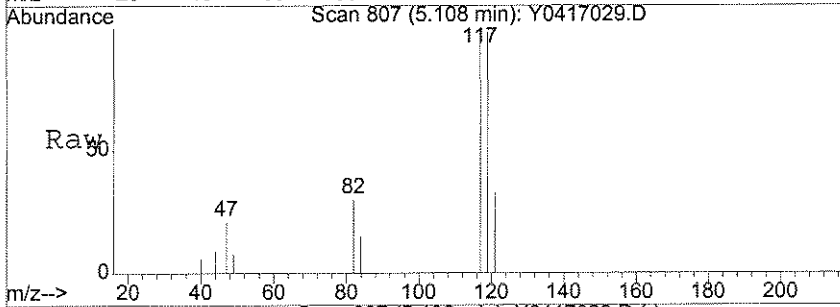
#31  
 Chloroform  
 Concen: 2.63 ug/l  
 RT: 4.71 min Scan# 739  
 Delta R.T. 0.01 min  
 Lab File: Y0417029.D  
 Acq: 17 Apr 2007 18:33

Tgt Ion: 83 Resp: 11585  
 Ion Ratio Lower Upper  
 83 100  
 85 66.7 43.3 83.3



#35  
 Carbon Tetrachloride  
 Concen: 2.04 ug/l  
 RT: 5.11 min Scan# 807  
 Delta R.T. -0.00 min  
 Lab File: Y0417029.D  
 Acq: 17 Apr 2007 18:33

Tgt Ion: 117 Resp: 6294  
 Ion Ratio Lower Upper  
 117 100  
 119 96.3 78.2 118.2



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-019  
 Lab File ID: Y0417030.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 18:58  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-019  
 Lab File ID: Y0417030.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 18:58  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL31-019

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417030.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:58

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

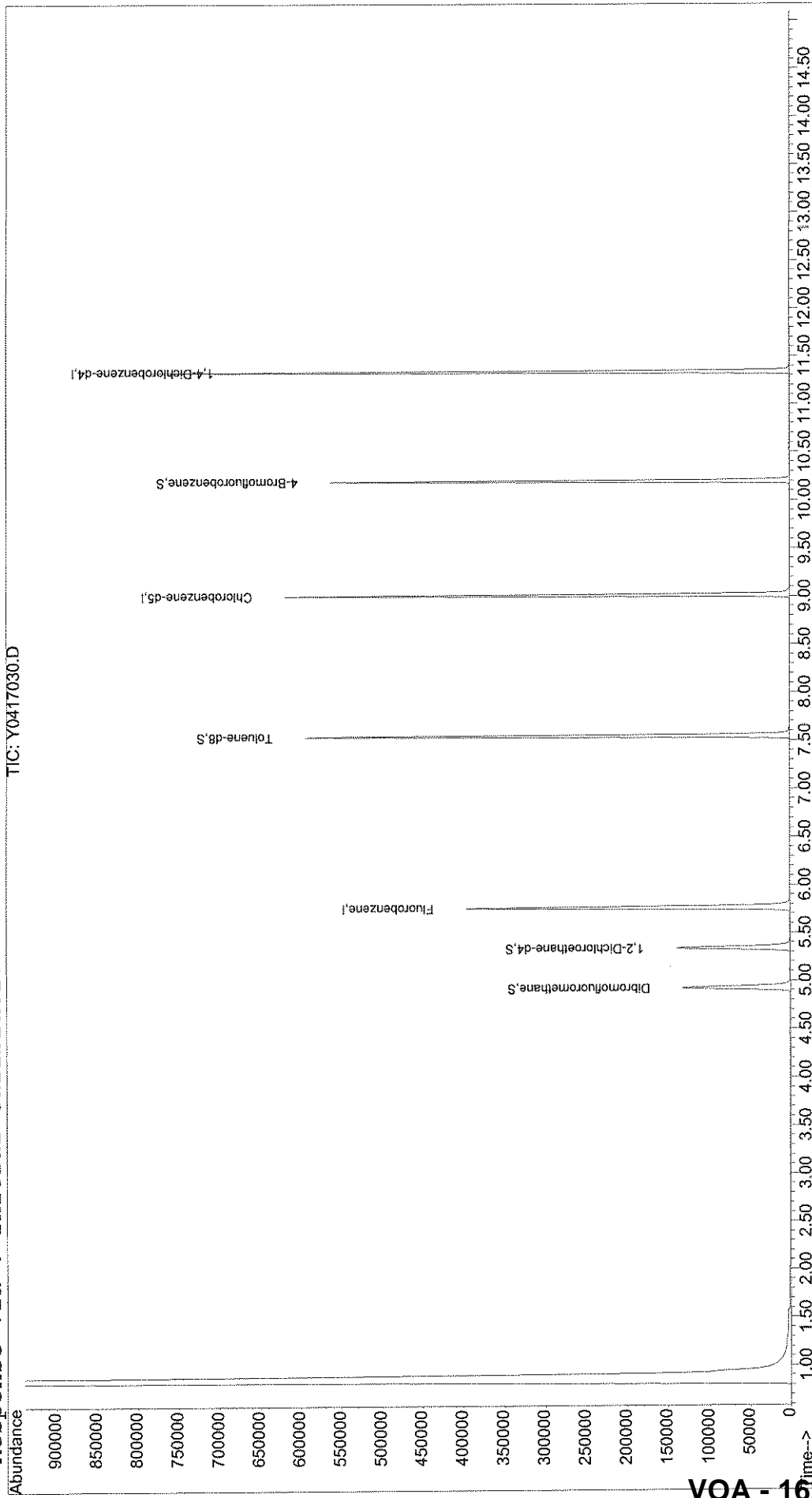
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417030.D Vial: 36  
Acq On : 17 Apr 2007 18:58 Operator: LNH  
Sample : JPL31-019 Inst : Yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:27 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417030.D  
 Acq On : 17 Apr 2007 18:58  
 Sample : JPL31-019  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.76	96	352260	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	152835	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	168349	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.92	111	97097	45.56	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	91.12%
37) 1,2-Dichloroethane-d4	5.34	65	108213	47.86	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	95.72%
51) Toluene-d8	7.54	98	346636	57.15	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	114.30%
72) 4-Bromofluorobenzene	10.20	95	145407	59.78	ug/l	0.00

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	495		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	2.36	40	59		N.D.	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417030.D 8260B.M Wed Apr 18 07:56:04 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417030.D  
 Acq On : 17 Apr 2007 18:58  
 Sample : JPL31-019  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	4.71	83	313		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	5.11	117	777		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	7.61	92	168		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	372		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.17	91	88		N.D.	
65) m,p-Xylene	9.30	106	121		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417030.D 8260B.M Wed Apr 18 07:56:05 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417030.D  
 Acq On : 17 Apr 2007 18:58  
 Sample : JPL31-019  
 Misc : 5mL+IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:27 2007

Vial: 36  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	71		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	10.47	91	57		N.D.	
78) 4-Chlorotoluene	10.47	91	57		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	196		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	135		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-020  
 Lab File ID: Y0417031.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 19:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-020  
 Lab File ID: Y0417031.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 19:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-12-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL31  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-020  
 Lab File ID: Y0417031.D  
 Date Collected: 04/04/2007  
 Date/Time Analyzed: 04/17/2007 19:22  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

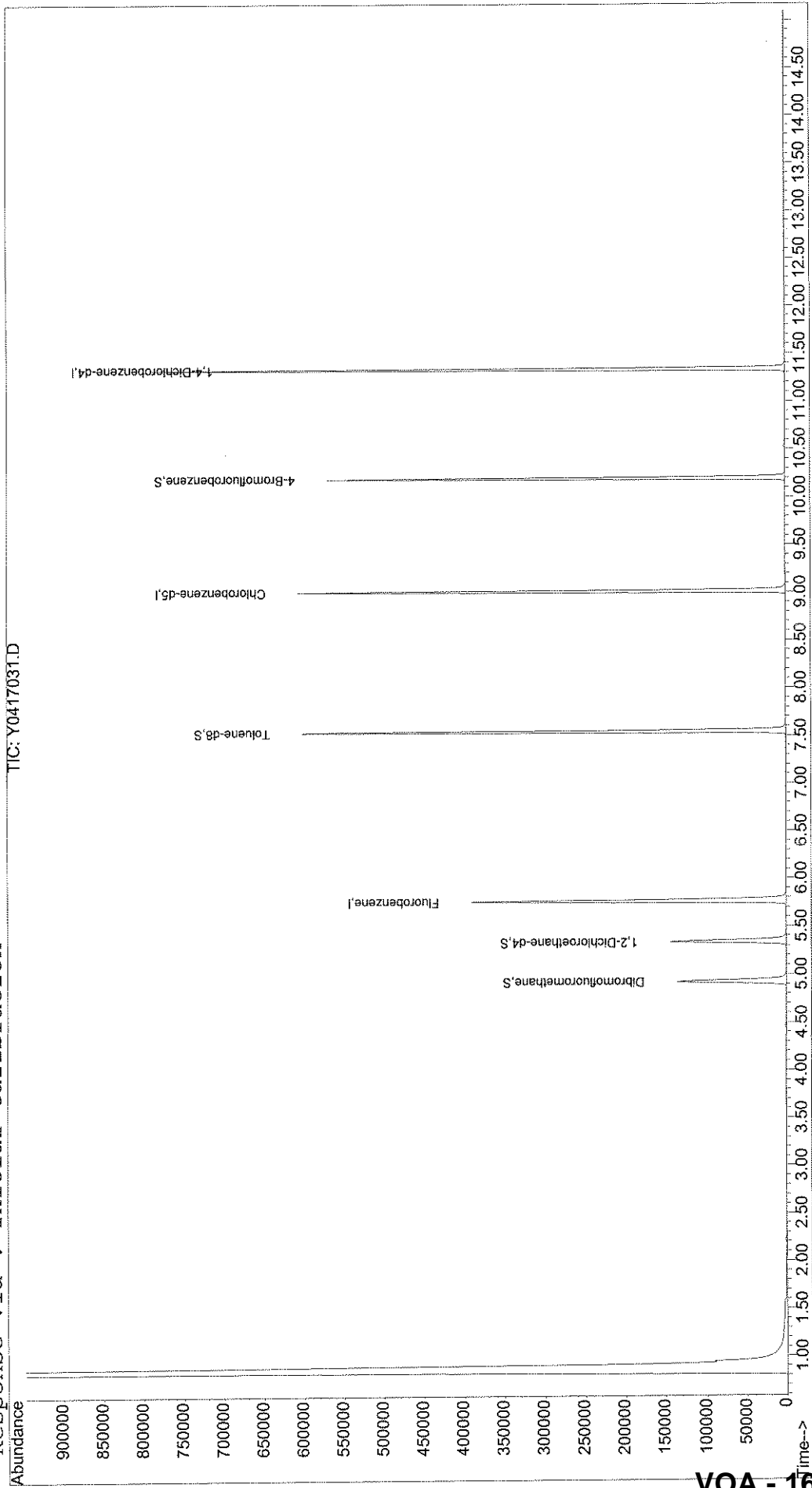
Comments:

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417031.D  
Acq On : 17 Apr 2007 19:22  
Sample : JPL31-020  
Misc : 5mL+IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 18 7:28 2007  
Quant Results File: 8260B.RES

Vial: 37  
Operator: LNH  
Inst : Yoda  
Multiplr: 1.00

Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973Y  
Last Update : Tue Mar 20 10:56:50 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417031.D  
 Acq On : 17 Apr 2007 19:22  
 Sample : JPL31-020  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:28 2007

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B  
 IS QA File : File name for IS QA unknown

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	5.77	96	358518	50.00	ug/l	0.00 NA%
50) Chlorobenzene-d5	9.01	82	155830	50.00	ug/l	0.00 NA%
70) 1,4-Dichlorobenzene-d4	11.34	152	175473	50.00	ug/l	0.00 NA%

System Monitoring Compounds

33) Dibromofluoromethane	4.93	111	98328	45.33	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	90.66%
37) 1,2-Dichloroethane-d4	5.34	65	109289	47.49	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	94.98%
51) Toluene-d8	7.54	98	351791	56.89	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	113.78%
72) 4-Bromofluorobenzene	10.20	95	149871	59.12	ug/l	0.00

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	0.00	43	0		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	2.22	76	2221		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	2.36	40	57		N.D.	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	
23) Vinyl acetate	0.00	43	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417031.D 8260B.M Wed Apr 18 07:56:12 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417031.D  
 Acq On : 17 Apr 2007 19:22  
 Sample : JPL31-020  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:28 2007

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	0.00	78	0		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	0.00	92	0		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	0.00	112	0		N.D.	
62) 1-Chlorohexane	9.01	91	476		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.29	91	244		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	
68) Bromoform	0.00	173	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 Y0417031.D 8260B.M Wed Apr 18 07:56:12 2007

Quantitation Report

Data File : X:\MSVOA\YODA\041707\Y0417031.D  
 Acq On : 17 Apr 2007 19:22  
 Sample : JPL31-020  
 Misc : 5mL+IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 18 7:28 2007

Vial: 37  
 Operator: LNH  
 Inst : yoda  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973Y  
 Last Update : Tue Mar 20 10:56:50 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260B

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
69) Isopropylbenzene	10.20	105	174		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	0.00	83	0		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	0.00	91	0		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	0.00	105	0		N.D.	
82) sec-butylbenzene	0.00	105	0		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	11.34	119	207		N.D.	
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	11.74	91	146		N.D.	
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

**TICS SUMMARY**

SDG JPL31

VOLATILES ANALYSIS



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041507MVOWY1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: B041507MVOWY1  
 Lab File ID: Y0415009.D  
 Date Collected: \_\_\_\_\_  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415009.D Vial: 22  
Acq On : 15 Apr 2007 13:48 Operator: LH  
Sample : B041507MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-1) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415009.D 8260B.M Mon Apr 16 15:53:22 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041707MVOWY1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: B041707MVOWY1  
 Lab File ID: Y0417016.D  
 Date Collected: \_\_\_\_\_  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417016.D Vial: 23  
Acq On : 17 Apr 2007 13:14 Operator: LNH  
Sample : B041707MVOWY1 Inst : yoda  
Misc : 5mLpfw+IS/SS(MV8-38-11) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417016.D 8260B.M Thu Apr 19 18:48:41 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-4-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-001  
 Lab File ID: Y0415012.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415012.D Vial: 25  
Acq On : 15 Apr 2007 15:49 Operator: LH  
Sample : JPL31-001 Inst : yoda  
Misc : 5mL+IS/SS #6 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415012.D 8260B.M Mon Apr 16 15:53:54 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-4-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-002  
 Lab File ID: Y0415013.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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04					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415013.D Vial: 26  
Acq On : 15 Apr 2007 16:14 Operator: LH  
Sample : JPL31-002 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415013.D 8260B.M Mon Apr 16 15:54:04 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-4-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-003  
 Lab File ID: Y0415014.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415014.D Vial: 27  
Acq On : 15 Apr 2007 16:38 Operator: LH  
Sample : JPL31-003 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415014.D 8260B.M Mon Apr 16 15:54:24 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-004  
 Lab File ID: Y0415015.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415015.D Vial: 28  
Acq On : 15 Apr 2007 17:02 Operator: LH  
Sample : JPL31-004 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415015.D 8260B.M Mon Apr 16 15:54:32 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-6-4/3/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec.  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-005  
 Lab File ID: Y0415010.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415010.D Vial: 23  
Acq On : 15 Apr 2007 15:00 Operator: LH  
Sample : JPL31-005 TB Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415010.D 8260B.M Mon Apr 16 15:53:35 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-11-4

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/15/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415016.D Vial: 29  
Acq On : 15 Apr 2007 17:27 Operator: LH  
Sample : JPL31-006 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415016.D 8260B.M Mon Apr 16 15:55:03 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-11-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-007  
 Lab File ID: Y0415017.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415017.D Vial: 30  
Acq On : 15 Apr 2007 17:51 Operator: LH  
Sample : JPL31-007 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415017.D 8260B.M Mon Apr 16 15:55:17 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-11-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/04/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/15/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415018.D Vial: 31  
Acq On : 15 Apr 2007 18:15 Operator: LH  
Sample : JPL31-008 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415018.D 8260B.M Mon Apr 16 15:55:38 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-11-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016855  
 Lab Sample ID: JPL31-009  
 Lab File ID: Y0415019.D  
 Date Collected: 04/04/2007  
 Date Analyzed: 04/15/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
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29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415019.D Vial: 32  
Acq On : 15 Apr 2007 18:40 Operator: LH  
Sample : JPL31-009 Inst : yoda  
Misc : 5mL+IS/SS #4 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415019.D 8260B.M Mon Apr 16 15:55:50 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-23-3

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/15/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
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29					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415020.D Vial: 33  
Acq On : 15 Apr 2007 19:05 Operator: LH  
Sample : JPL31-011 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415020.D 8260B.M Mon Apr 16 15:56:12 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-23-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-012  
 Lab File ID: Y0417024.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417024.D Vial: 30  
Acq On : 17 Apr 2007 16:31 Operator: LNH  
Sample : JPL31-012 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417024.D 8260B.M Wed Apr 18 07:52:40 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-23-1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/17/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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29					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417025.D Vial: 31  
Acq On : 17 Apr 2007 16:55 Operator: LNH  
Sample : JPL31-013 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417025.D 8260B.M Wed Apr 18 07:52:47 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-014  
 Lab File ID: Y0417026.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417026.D Vial: 32  
Acq On : 17 Apr 2007 17:20 Operator: LNH  
Sample : JPL31-014 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417026.D 8260B.M Wed Apr 18 07:52:58 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-7-4/4/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016855

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-015

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0415011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/15/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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04					
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041507\Y0415011.D Vial: 24  
Acq On : 15 Apr 2007 15:25 Operator: LH  
Sample : JPL31-015 TB Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0415011.D 8260B.M Mon Apr 16 15:53:45 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-12-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-016  
 Lab File ID: Y0417027.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417027.D Vial: 33  
Acq On : 17 Apr 2007 17:44 Operator: LNH  
Sample : JPL31-016 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417027.D 8260B.M Wed Apr 18 07:53:04 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-12-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-017  
 Lab File ID: Y0417028.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417028.D Vial: 34  
Acq On : 17 Apr 2007 18:09 Operator: LNH  
Sample : JPL31-017 Inst : yoda  
Misc : 5mL+IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417028.D 8260B.M Wed Apr 18 07:53:11 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-12-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-018  
 Lab File ID: Y0417029.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
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28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417029.D Vial: 35  
Acq On : 17 Apr 2007 18:33 Operator: LNH  
Sample : JPL31-018 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417029.D 8260B.M Wed Apr 18 07:53:18 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-12-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL31  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: DB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016916  
 Lab Sample ID: JPL31-019  
 Lab File ID: Y0417030.D  
 Date Collected: 04/05/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417030.D Vial: 36  
Acq On : 17 Apr 2007 18:58 Operator: LNH  
Sample : JPL31-019 Inst : yoda  
Misc : 5mL+IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417030.D 8260B.M Wed Apr 18 07:53:26 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-12-1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL31

Run Sequence: R016916

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL31-020

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: Y0417031.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/17/2007

GC Column: DB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\YODA\041707\Y0417031.D Vial: 37  
Acq On : 17 Apr 2007 19:22 Operator: LNH  
Sample : JPL31-020 Inst : yoda  
Misc : 5mL+IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\YODA\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973Y  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
Y0417031.D 8260B.M Wed Apr 18 07:53:33 2007

**Metals Data**

**JPL31**

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-4-3	JPL31-001
MW-4-3D	JPL31-001D
MW-4-3MS	JPL31-001MS
MW-4-2	JPL31-002
MW-4-1	JPL31-003
EB-6-4/3/07	JPL31-004
MW-11-3	JPL31-007
MW-11-3MS	JPL31-007MS
MW-11-3MSD	JPL31-007MSD
MW-11-2	JPL31-008
MW-11-1	JPL31-009
MW-23-4	JPL31-010
MW-23-3	JPL31-011
MW-23-3MS	JPL31-011MS
MW-23-3MSD	JPL31-011MSD
MW-23-2	JPL31-012
MW-23-1	JPL31-013
EB-7-4/4/07	JPL31-014
EB-7-4/4/07MS	JPL31-014MS
EB-7-4/4/07MSD	JPL31-014MSD
MW-12-3	JPL31-018

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

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I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher

Name: Bill Ambacher

Date: 4/27/07

Title: Inorganics/Metals Manager

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

SOW No.: \_\_\_\_\_

Sample No.  
MW-12-2  
MW-12-1

Lab Sample ID  
JPL31-019  
JPL31-020

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

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I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher

Name: Bill Ambacher

Date: 4/27/07

Title: Inorganics/Metals Manager

## **Metals Analysis Data Sheets**

## INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-4-3

Lab Name: Laucks Laboratories Contract: JPL Groundwater Monitorin  
 Lab Code: LAUCKS SDG No.: JPL31  
 Matrix (soil/water): Water Lab Sample ID: JPL31-001  
 Level (low/med): LOW Date Received: 04/04/2007  
 % Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.98			M	R016781

Color Before: Colorless Clarity Before: Clear Texture: \_\_\_\_\_  
 Color After: Colorless Clarity After: Clear Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-4-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-002

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.82			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-4-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-003

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.71			M	R016781

Color Before: Colorless Clarity Before: Clear Texture: \_\_\_\_\_

Color After: Colorless Clarity After: Clear Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-6-4/3/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-004

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.39			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-11-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-007

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-11-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-008

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-11-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-009

Level (low/med): LOW

Date Received: 04/04/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-23-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-010

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.47			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-23-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-011

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.13			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-23-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-012

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.02			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-23-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-013

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	2.77			M	R016781

Color Before: Colorless Clarity Before: Clear Texture: \_\_\_\_\_

Color After: Colorless Clarity After: Clear Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-7-4/4/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-014

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.60			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-12-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-018

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.00	U		M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-12-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-019

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.22			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-12-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL31

Matrix (soil/water): Water

Lab Sample ID: JPL31-020

Level (low/med): LOW

Date Received: 04/05/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.11			M	R016724

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL31**

**COVER PAGE - INORGANIC ANALYSES DATA PACKAGE**

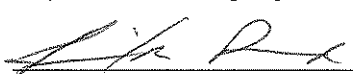
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL31

Client Identification	Lab Sample Work Order Number
MW-4-3	JPL31-001DL
MW-4-3MS	JPL31-001MS
MW-4-3MSD	JPL31-001MSD
MW-4-2	JPL31-002DL
MW-4-1	JPL31-003DL
EB-6-4/3/07	JPL31-004
MW-11-4	JPL31-006
MW-11-3	JPL31-007DL
MW-11-3MS	JPL31-007MS
MW-11-3MSD	JPL31-007MSD
MW-11-2	JPL31-008DL
MW-11-1	JPL31-009
MW-11-1MS	JPL31-009MS
MW-11-1MSD	JPL31-009MSD
MW-23-3	JPL31-011DL
MW-23-3MS	JPL31-011MS
MW-23-3MSD	JPL31-011MSD
MW-23-2	JPL31-012DL
MW-23-1	JPL31-013DL
EB-7-4/4/07	JPL31-014
EB-7-4/4/07MS	JPL31-014MS

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Penner

Date: 4-25-07

Title: Inorganics Lead

COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

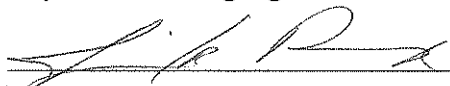
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL31

Client Identification	Lab Sample Work Order Number
EB-7-4/4/07MSD	JPL31-014MSD
MW-12-5	JPL31-016DL
MW-12-4	JPL31-017DL
MW-12-3	JPL31-018DL
MW-12-2	JPL31-019DL
MW-12-1	JPL31-020DL

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Penner

Date: 4-25-07

Title: Inorganics Lead



## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-4-3 **Date/Time Collected:** 04/03/2007 07:44  
**Lab Sample ID:** JPL31-001 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL31  
Sample Number: MW-4-2 Date/Time Collected: 04/03/2007 08:37  
Lab Sample ID: JPL31-002 Date/Time Received: 04/04/2007 09:00  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-4-1 **Date/Time Collected:** 04/03/2007 09:10  
**Lab Sample ID:** JPL31-003 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	280		4.0	0.56	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** EB-6-4/3/07 **Date/Time Collected:** 04/03/2007 08:59  
**Lab Sample ID:** JPL31-004 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-11-4 **Date/Time Collected:** 04/03/2007 10:19  
**Lab Sample ID:** JPL31-006 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-11-3 **Date/Time Collected:** 04/03/2007 10:50  
**Lab Sample ID:** JPL31-007 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-11-2 **Date/Time Collected:** 04/03/2007 11:38  
**Lab Sample ID:** JPL31-008 **Date/Time Received:** 04/04/2007 09:00  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894





Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-23-3 **Date/Time Collected:** 04/04/2007 08:22  
**Lab Sample ID:** JPL31-011 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL31  
Sample Number: MW-23-2 Date/Time Collected: 04/04/2007 09:10  
Lab Sample ID: JPL31-012 Date/Time Received: 04/05/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	7.9		4.0	0.56	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-23-1 **Date/Time Collected:** 04/04/2007 09:44  
**Lab Sample ID:** JPL31-013 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** EB-7-4/4/07 **Date/Time Collected:** 04/04/2007 09:33  
**Lab Sample ID:** JPL31-014 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-12-5 **Date/Time Collected:** 04/04/2007 11:02  
**Lab Sample ID:** JPL31-016 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-12-4 **Date/Time Collected:** 04/04/2007 11:22  
**Lab Sample ID:** JPL31-017 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	3.2		2.0	0.28	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-12-3 **Date/Time Collected:** 04/04/2007 11:46  
**Lab Sample ID:** JPL31-018 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894



Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-12-2 **Date/Time Collected:** 04/04/2007 12:13  
**Lab Sample ID:** JPL31-019 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	2.5	U	2.5	0.35	04/17/2007	04/18/2007	R016894

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL31  
**Sample Number:** MW-12-1 **Date/Time Collected:** 04/04/2007 13:03  
**Lab Sample ID:** JPL31-020 **Date/Time Received:** 04/05/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/17/2007	04/18/2007	R016894

**SAMPLE DATA**

SDG JPL32

VOLATILES ANALYSIS

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417010.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 11:12

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-002  
 Lab File ID: M0417010.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 11:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.95	J
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-002  
 Lab File ID: M0417010.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 11:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

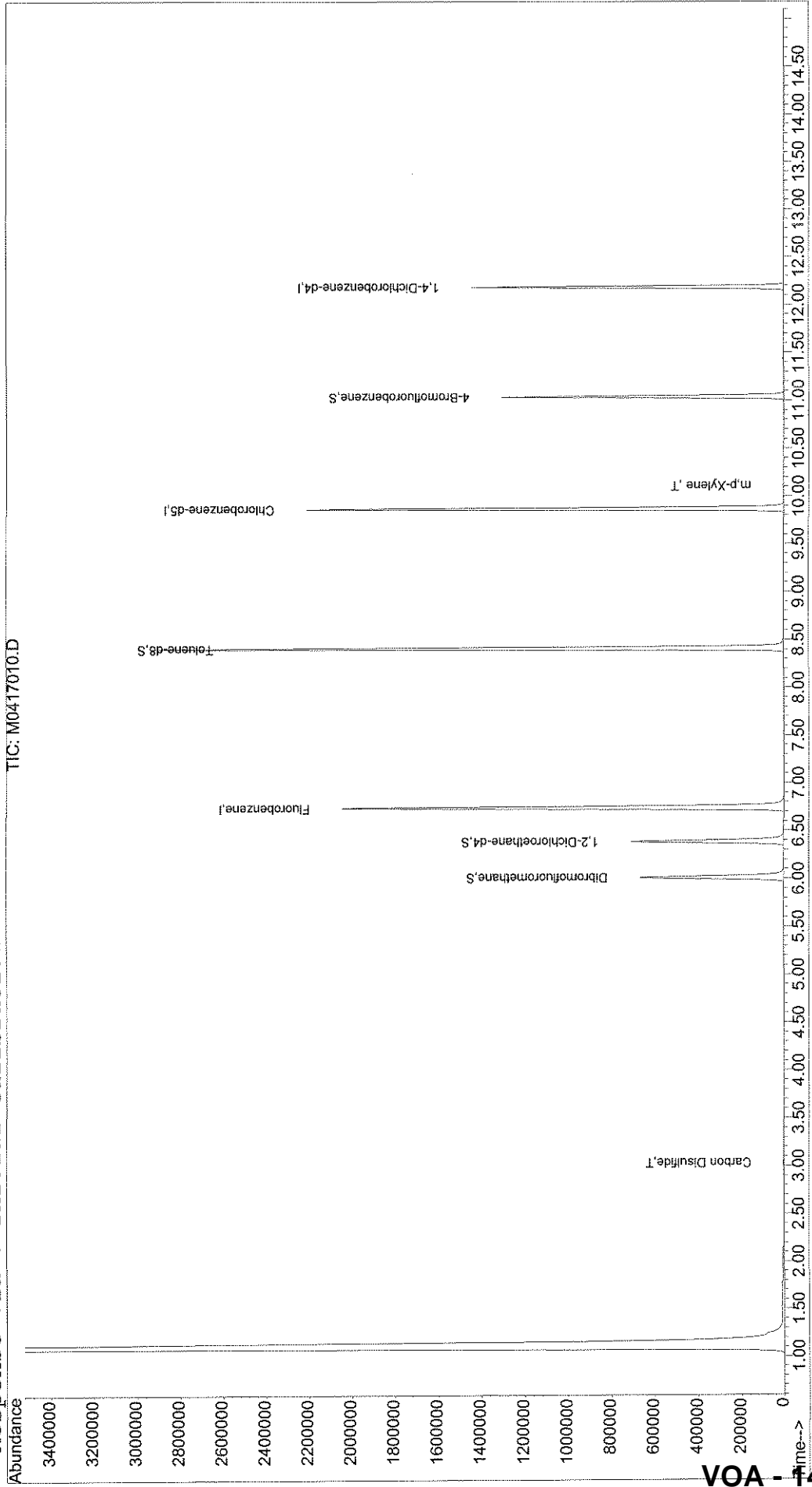
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments: \_\_\_\_\_

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417010.D Vial: 55  
Acq On : 17 Apr 2007 11:12 Operator: DGA  
Sample : JPL32-002 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:07 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA - 14

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417010.D  
 Acq On : 17 Apr 2007 11:12  
 Sample : JPL32-002  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:07 2007

Vial: 55  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar)
1) Fluorobenzene	6.74	96	1983740	50.00	ug/l	0.00	101.63%
50) Chlorobenzene-d5	9.86	82	651534	50.00	ug/l	0.00	92.64%
70) 1,4-Dichlorobenzene-d4	12.18	152	344928	50.00	ug/l	0.00	88.62%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	466100	49.80	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.60%	
37) 1,2-Dichloroethane-d4	6.38	65	528039	49.29	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.58%	
51) Toluene-d8	8.41	98	1912415	53.15	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.30%	
72) 4-Bromofluorobenzene	11.04	95	391156	54.98	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.96%	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.43	50	120	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.85	96	60	N.D.		
6) Chloroethane	1.95	64	60	N.D.		
7) Trichlorofluoromethane	2.20	101	68	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.98	43	295	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.03	76	7835	0.40	ug/l	100
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	3.33	43	59	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.87	96	142	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.60	63	57	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417010.D 8260B.M Wed Apr 25 09:07:59 2007

8 04/25/07  
 VOA-15 Page 1



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417010.D  
 Acq On : 17 Apr 2007 11:12  
 Sample : JPL32-002  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:07 2007

Vial: 55  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.67	43	55		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.37	96	121		N.D.	
27) 2-Butanone	5.54	43	196		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.78	41	61		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	5.94	97	62		N.D.	
34) Cyclohexane	6.01	56	61		N.D.	
35) Carbon Tetrachloride	6.14	117	69		N.D.	
36) 1,1-Dichloropropene	6.15	75	393		N.D.	
38) Benzene	6.43	78	1556		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.16	130	371		N.D.	
42) Methylcyclohexane	7.30	83	379		N.D.	
43) 1,2-Dichloropropane	7.38	63	58		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.50	41	55		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.17	75	57		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	1204		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.78	97	57		N.D.	
56) Tetrachloroethene	9.02	166	431		N.D.	
57) 1,3-Dichloropropane	9.09	76	59		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	797		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.10	131	93		N.D.	
64) Ethylbenzene	9.98	91	2003		N.D.	
65) m,p-Xylene	10.11	106	1588	0.95	ug/l	88
66) o-xylene	10.50	106	480		N.D.	
67) Styrene	10.53	104	558		N.D.	

Quantitation Report

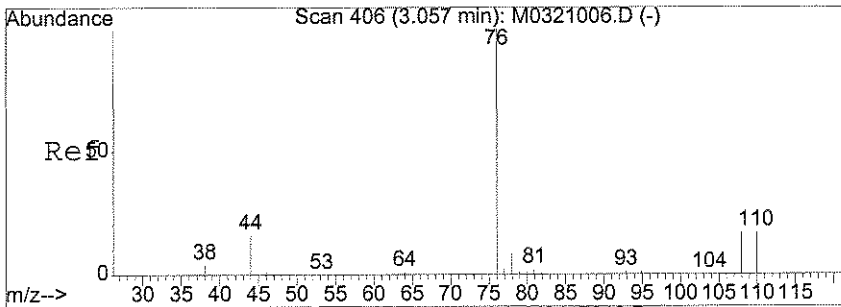
Data File : X:\MSVOA\MOBY\041707\M0417010.D  
 Acq On : 17 Apr 2007 11:12  
 Sample : JPL32-002  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:07 2007

Vial: 55  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

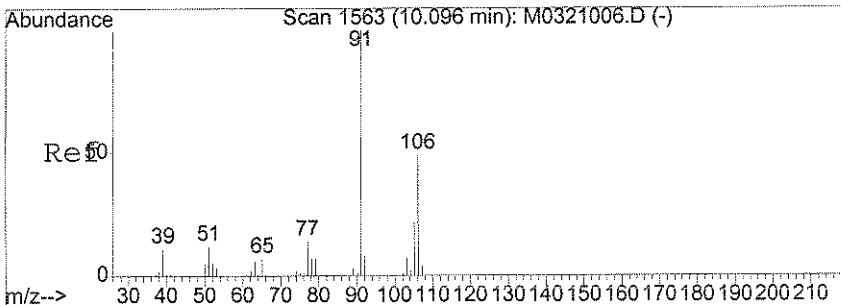
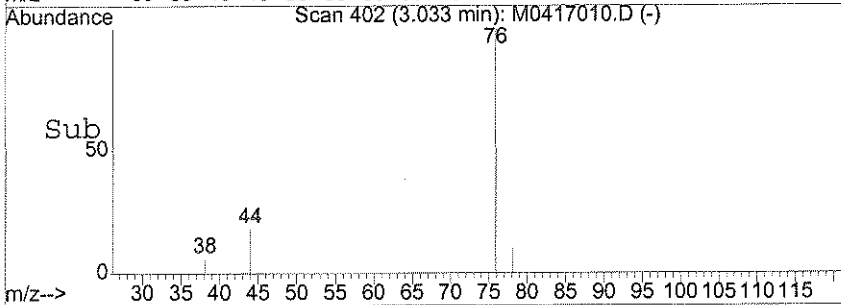
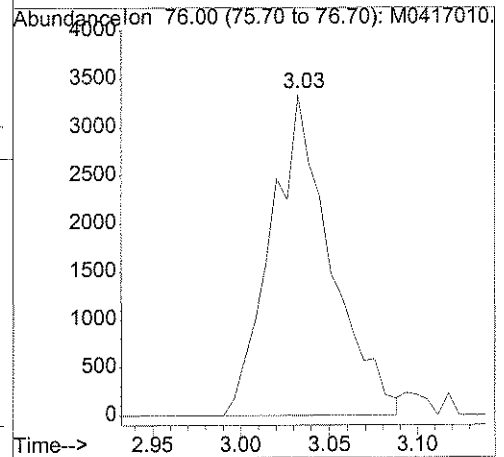
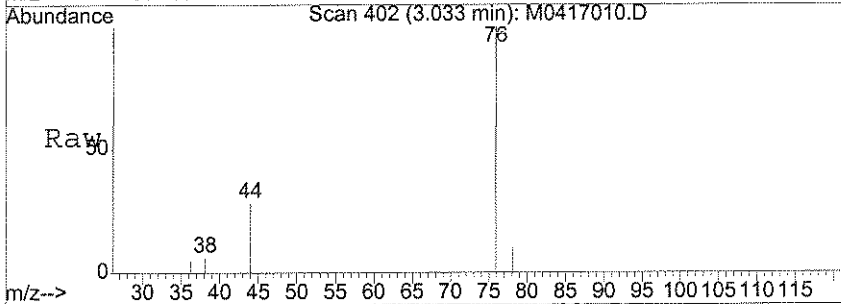
Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

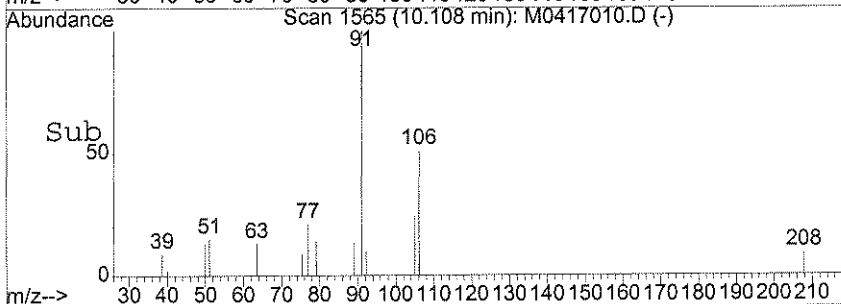
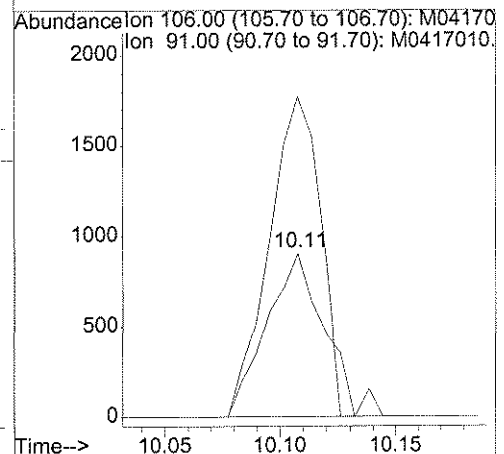
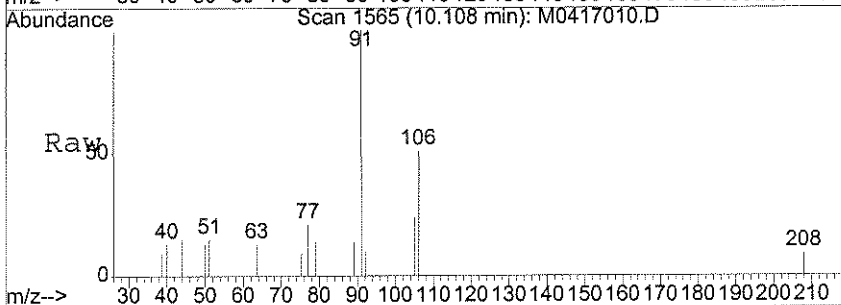
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	69		N.D.	
69) Isopropylbenzene	10.86	105	780		N.D.	
71) trans-1,4-Dichloro-2-buten	10.77	53	64		N.D.	
73) Bromobenzene	11.19	156	137		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.12	83	78		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	d
76) n-Propylbenzene	11.28	120	216		N.D.	
77) 2-Chlorotoluene	11.37	91	532		N.D.	
78) 4-Chlorotoluene	11.48	91	784		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	451		N.D.	
80) tert-Butylbenzene	11.77	119	610		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	685		N.D.	
82) sec-butylbenzene	11.98	105	1062		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	454		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	786		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	303		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.31	225	409		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



#14  
 Carbon Disulfide  
 Concen: 0.40 ug/l  
 RT: 3.03 min Scan# 402  
 Delta R.T. 0.01 min  
 Lab File: M0417010.D  
 Acq: 17 Apr 2007 11:12  
 Tgt Ion: 76 Resp: 7835



#65  
 m,p-Xylene  
 Concen: 0.95 ug/l  
 RT: 10.11 min Scan# 1565  
 Delta R.T. -0.00 min  
 Lab File: M0417010.D  
 Acq: 17 Apr 2007 11:12  
 Tgt Ion: 106 Resp: 1588  
 Ion Ratio Lower Upper  
 106 100  
 91 172.0 169.6 209.6



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 11:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.72	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	1.5	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.34	J
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 11:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417011.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 11:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

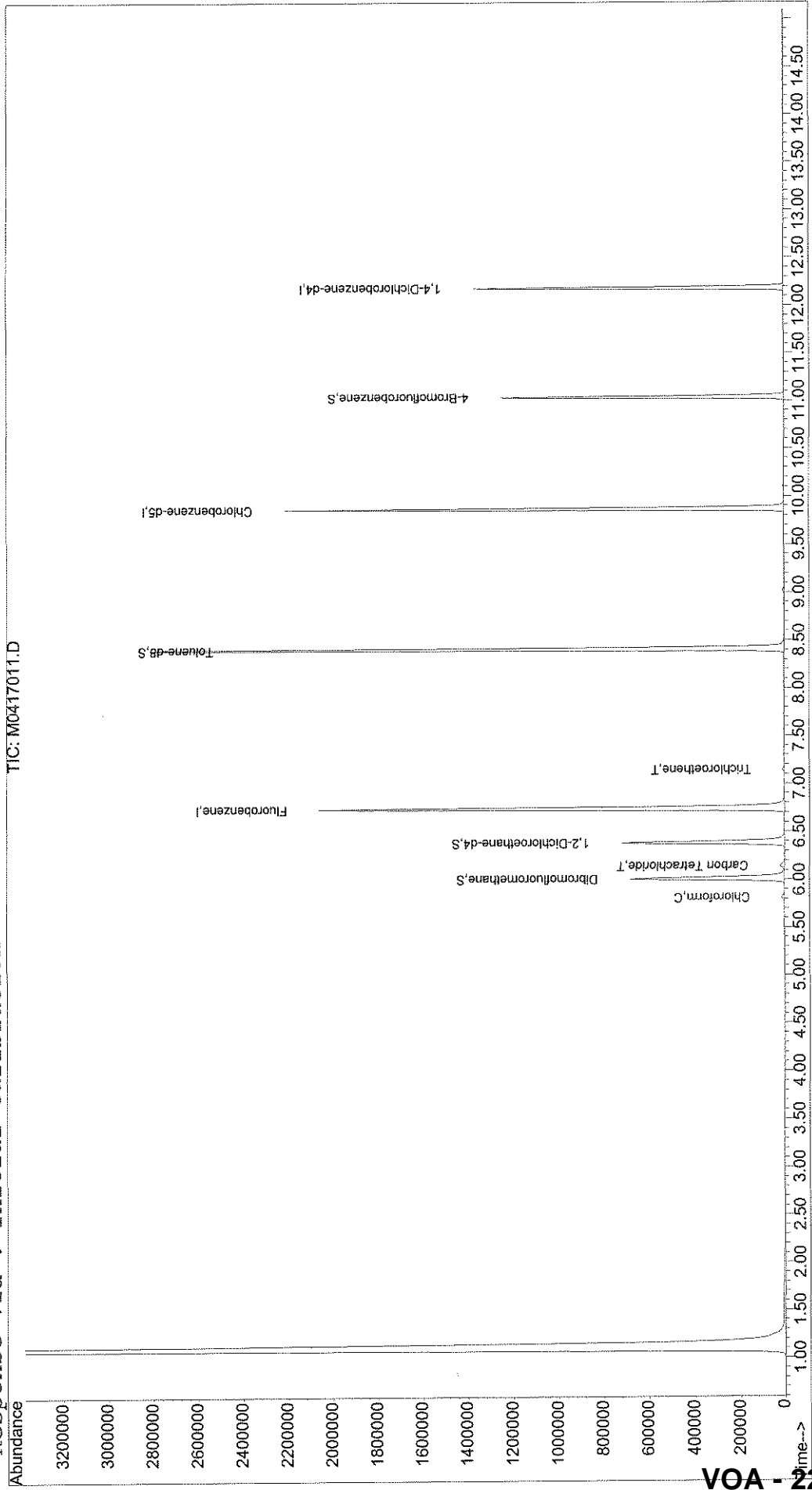
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417011.D  
Acq On : 17 Apr 2007 11:37  
Sample : JPL32-003  
Misc : #3 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 10:12 2007  
Vial: 56  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417011.D  
 Acq On : 17 Apr 2007 11:37  
 Sample : JPL32-003  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 10:12 2007

Vial: 56  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973M

Last Update : Tue Apr 24 12:05:10 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1969022	50.00	ug/l	0.00 100.88%
50) Chlorobenzene-d5	9.87	82	642952	50.00	ug/l	0.00 91.42%
70) 1,4-Dichlorobenzene-d4	12.18	152	342296	50.00	ug/l	0.00 87.94%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	466045	50.17	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.34%
37) 1,2-Dichloroethane-d4	6.39	65	534752	50.29	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.58%
51) Toluene-d8	8.41	98	1889064	53.20	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.40%
72) 4-Bromofluorobenzene	11.04	95	389726	55.20	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.40%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.81	96	95	N.D.	
6) Chloroethane	1.93	64	60	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	2.69	56	62	N.D.	
9) 1,1-Dichloroethene	2.79	96	495	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	2210	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.36	43	58	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	3.90	53	69	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	4.54	63	1302	N.D.	

(#) = qualifier out of range (m) = manual integration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417011.D  
 Acq On : 17 Apr 2007 11:37  
 Sample : JPL32-003  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 10:12 2007

Vial: 56  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	5.43	77	57		N.D.	
26) cis-1,2-Dichloroethene	5.32	96	57		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	
28) Propionitrile	5.51	54	60		N.D.	
29) Bromochloromethane	5.64	128	55		N.D.	
30) Methacrylonitrile	5.60	41	55		N.D.	
31) Chloroform	5.81	83	11449	0.72	ug/l	99
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	5.99	56	60		N.D.	
35) Carbon Tetrachloride	6.14	117	13991	1.49	ug/l	100
36) 1,1-Dichloropropene	6.16	75	61		N.D.	
38) Benzene	6.40	78	1304		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D. d	
41) Trichloroethene	7.15	130	3001	0.34	ug/l #	53
42) Methylcyclohexane	7.31	83	69		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.71	83	65		N.D.	
47) 2-Chloroethyl vinyl ether	7.86	63	66		N.D.	
48) cis-1,3-Dichloropropene	8.24	75	78		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.47	92	981		N.D.	
53) trans-1,3-Dichloropropene	8.69	75	58		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.89	97	73		N.D.	
56) Tetrachloroethene	9.03	166	1446		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	374		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	849		N.D.	
65) m,p-Xylene	0.00	106	0		N.D. d	
66) o-xylene	10.50	106	302		N.D.	
67) Styrene	10.52	104	262		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417011.D 8260B.M Wed Apr 25 10:12:34 2007

*J. O. J. 04/25/07*  
 VOA-24 Page 2

Quantitation Report

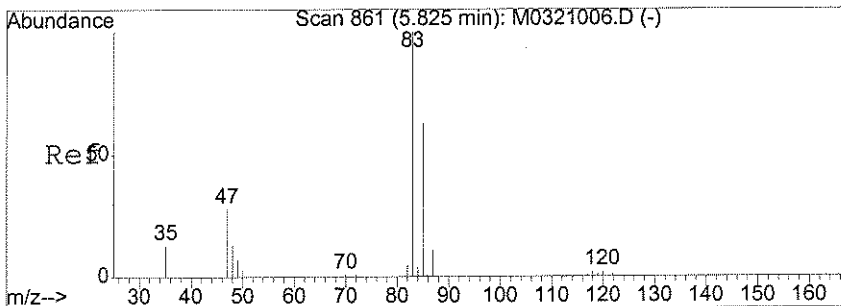
Data File : X:\MSVOA\MOBY\041707\M0417011.D  
 Acq On : 17 Apr 2007 11:37  
 Sample : JPL32-003  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 10:12 2007

Vial: 56  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

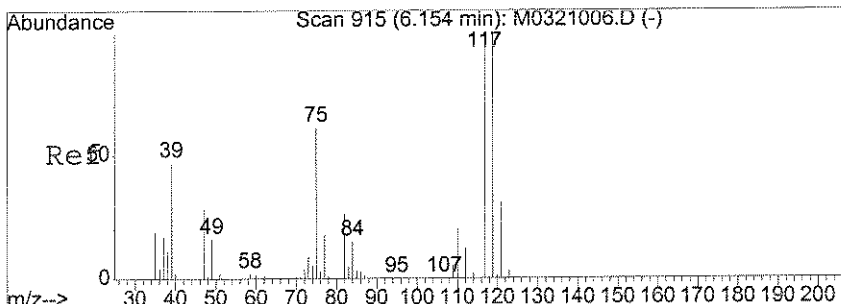
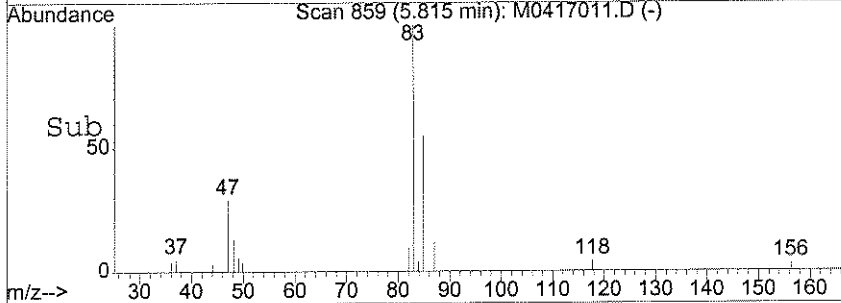
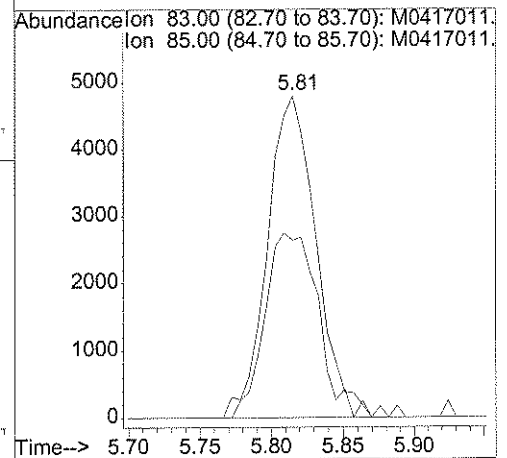
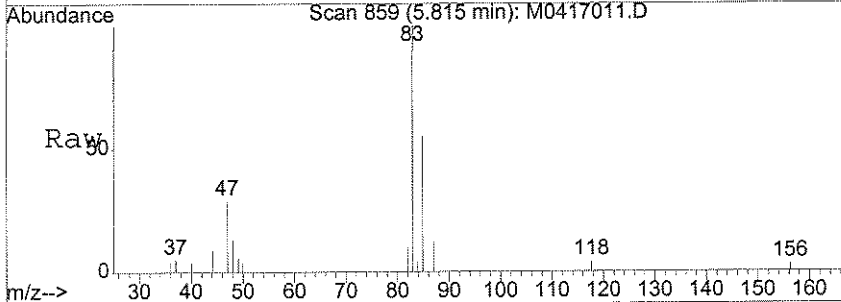
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	182		N.D.	
69) Isopropylbenzene	10.86	105	241		N.D.	
71) trans-1,4-Dichloro-2-buten	11.06	53	60		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	314		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	213		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.36	91	168		N.D.	
78) 4-Chlorotoluene	11.48	91	425		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	436		N.D.	
80) tert-Butylbenzene	11.77	119	68		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	525		N.D.	
82) sec-butylbenzene	11.97	105	619		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	750		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	771		N.D.	
86) 1,2-Dichlorobenzene	12.59	146	647		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



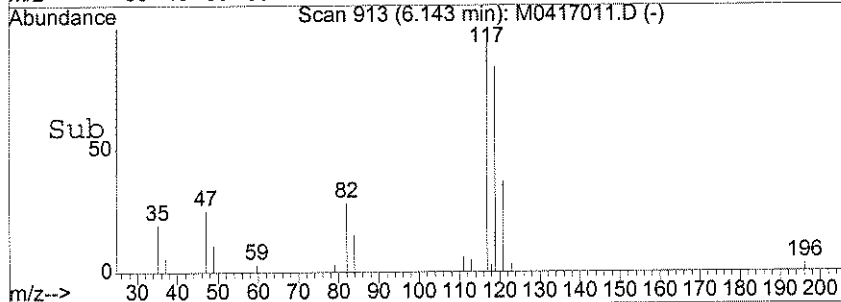
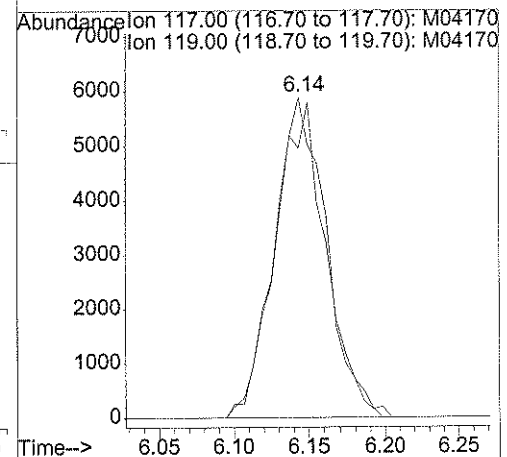
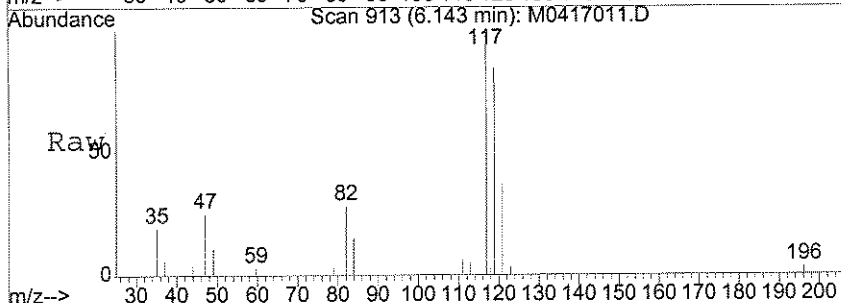
#31  
 Chloroform  
 Concen: 0.72 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0417011.D  
 Acq: 17 Apr 2007 11:37

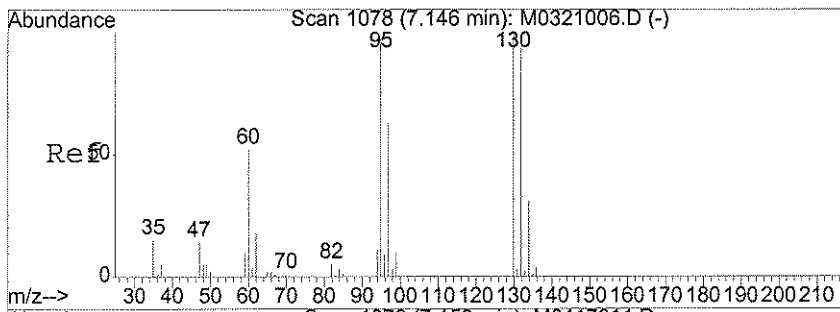
Tgt Ion	Resp	Lower	Upper
83	11449		
85	61.9	41.2	81.2



#35  
 Carbon Tetrachloride  
 Concen: 1.49 ug/l  
 RT: 6.14 min Scan# 913  
 Delta R.T. 0.00 min  
 Lab File: M0417011.D  
 Acq: 17 Apr 2007 11:37

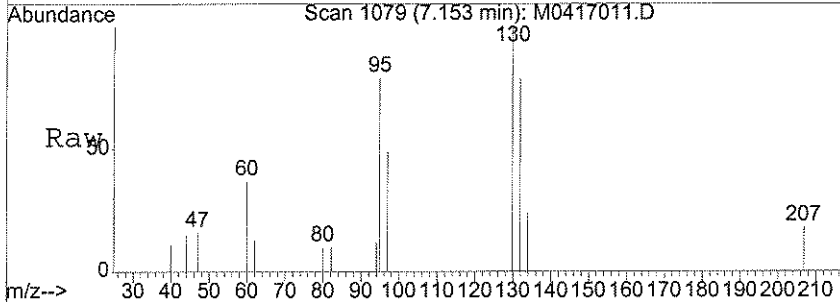
Tgt Ion	Resp	Lower	Upper
117	13991		
119	97.9	78.2	118.2



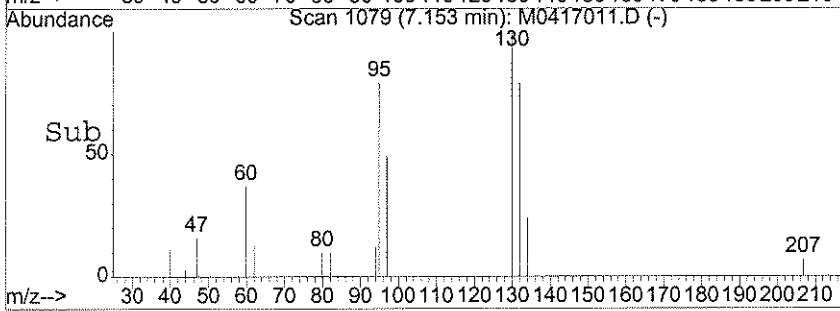
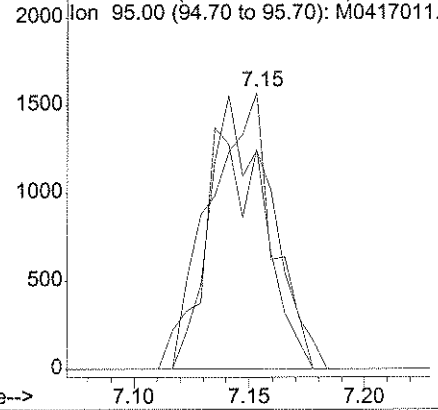


#41  
 Trichloroethene  
 Concen: 0.34 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0417011.D  
 Acq: 17 Apr 2007 11:37

Tgt Ion	Resp	Lower	Upper
130	3001		
132	92.8	75.0	115.0
95	0.0	69.4	109.4#



Abundance  
 Ion 130.00 (129.70 to 130.70): M04170  
 Ion 132.00 (131.70 to 132.70): M04170  
 Ion 95.00 (94.70 to 95.70): M0417011



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-004  
 Lab File ID: M0417012.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 12:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	1.7	
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	4.0	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	11	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-004  
 Lab File ID: M0417012.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 12:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	5.9	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-24-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-004  
 Lab File ID: M0417012.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 12:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

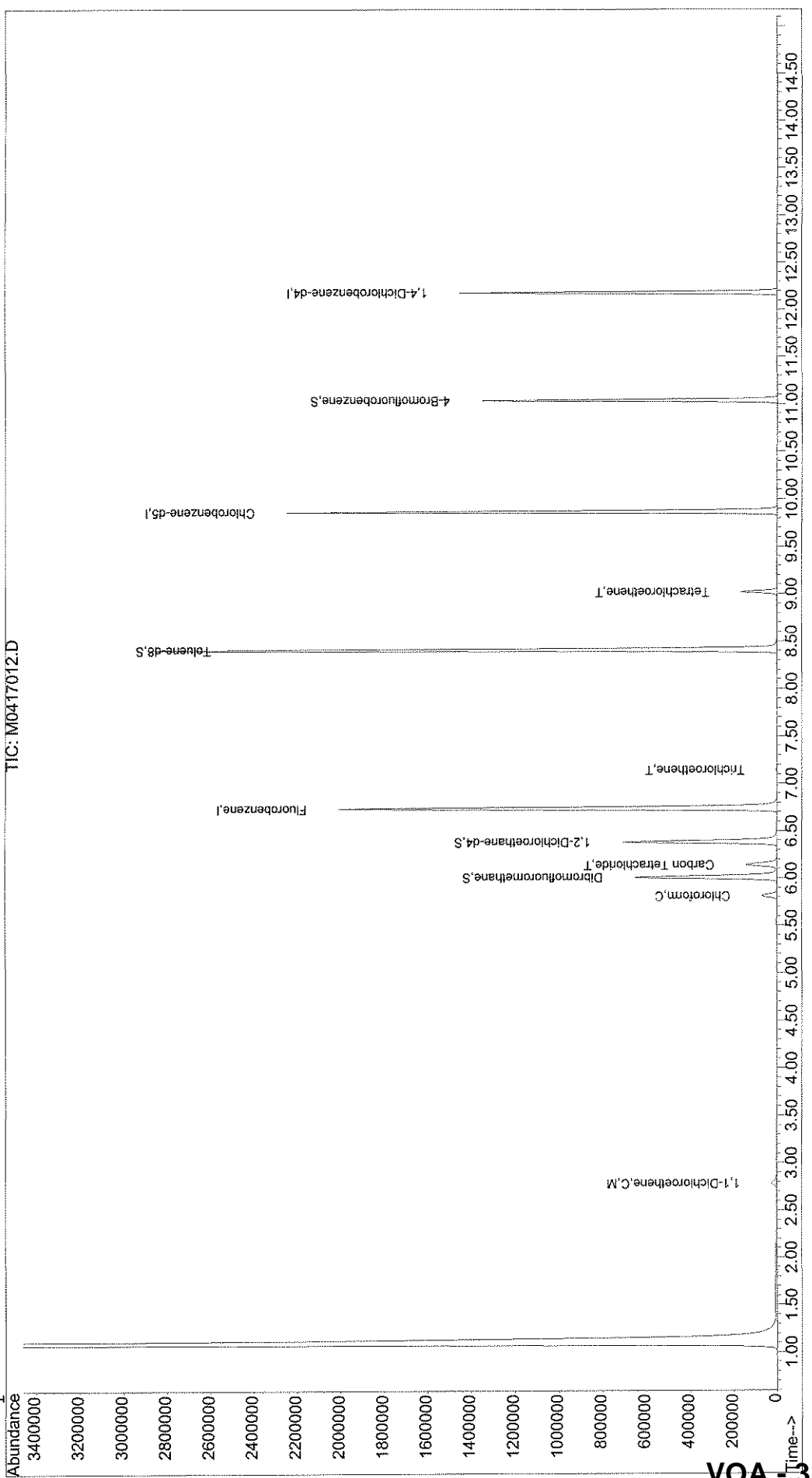
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417012.D  
Acq On : 17 Apr 2007 12:01  
Sample : JPL32-004  
Misc : #2 5ml +IS/SS(524)  
MS Integration Params: iteint.p  
Quant Time: Apr 25 9:11 2007  
Vial: 57  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration





Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417012.D  
 Acq On : 17 Apr 2007 12:01  
 Sample : JPL32-004  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:11 2007

Vial: 57  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	1929890	50.00	ug/l	0.00 98.87%
50) Chlorobenzene-d5	9.86	82	648721	50.00	ug/l	0.00 92.24%
70) 1,4-Dichlorobenzene-d4	12.18	152	343233	50.00	ug/l	0.00 88.18%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	459210	50.44	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.88%
37) 1,2-Dichloroethane-d4	6.38	65	524630	50.34	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.68%
51) Toluene-d8	8.41	98	1894749	52.88	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.76%
72) 4-Bromofluorobenzene	11.04	95	394932	55.78	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.56%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	1.97	64	59	N.D.		
7) Trichlorofluoromethane	2.18	101	55	N.D.		
8) Acrolein	2.75	56	57	N.D.		
9) 1,1-Dichloroethene	2.78	96	9943	1.69	ug/l	95
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	3.00	108	66	N.D.		
14) Carbon Disulfide	3.01	76	1156	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.54	63	215	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417012.D 8260B.M Wed Apr 25 11:36:02 2007

*[Handwritten Signature]*  
 VOA-32 Page 1

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417012.D  
 Acq On : 17 Apr 2007 12:01  
 Sample : JPL32-004  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:11 2007

Vial: 57  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.55	43	84	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	5.33	77	79	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.39	43	60	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	5.75	128	61	N.D.		
30) Methacrylonitrile	5.69	41	57	N.D.		
31) Chloroform	5.82	83	61829	3.99	ug/l	94
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	6.01	56	58	N.D.		
35) Carbon Tetrachloride	6.15	117	103716	11.26	ug/l	97
36) 1,1-Dichloropropene	6.10	75	65	N.D.		
38) Benzene	6.40	78	1232	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.	d	
41) Trichloroethene	7.15	130	2014	0.23	ug/l #	53
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	7.66	93	61	N.D.		
45) Methyl methacrylate	7.50	41	58	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	7.94	63	62	N.D.		
48) cis-1,3-Dichloropropene	8.06	75	78	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	8.47	92	1092	N.D.		
53) trans-1,3-Dichloropropene	8.76	75	60	N.D.		
54) Ethyl methacrylate	8.62	69	72	N.D.		
55) 1,1,2-Trichloroethane	9.01	97	375	N.D.		
56) Tetrachloroethene	9.02	166	46174	5.93	ug/l	97
57) 1,3-Dichloropropane	9.30	76	87	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.	d	
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	9.35	107	58	N.D.		
61) Chlorobenzene	9.90	112	358	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	10.00	91	864	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.	d	
66) o-xylene	10.50	106	66	N.D.		
67) Styrene	10.53	104	199	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417012.D 8260B.M Wed Apr 25 11:36:03 2007

*J. [Signature]*  
 Page 2  
 VOA-33

Quantitation Report

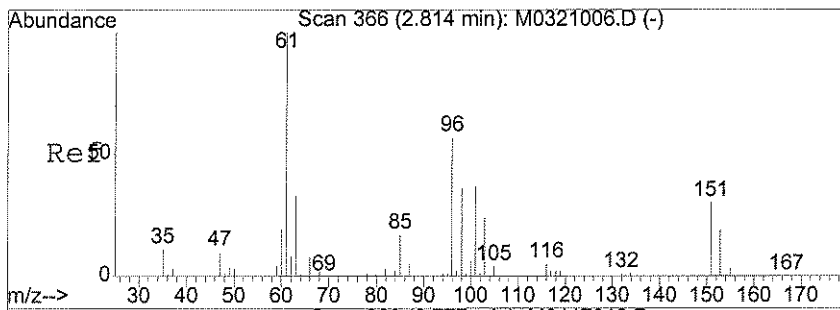
Data File : X:\MSVOA\MOBY\041707\M0417012.D  
 Acq On : 17 Apr 2007 12:01  
 Sample : JPL32-004  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:11 2007

Vial: 57  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

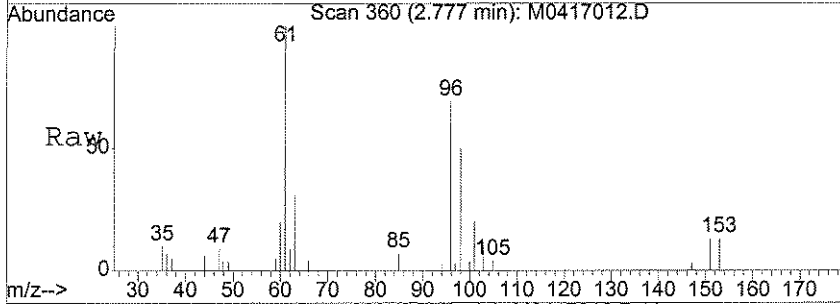
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	77		N.D.	
69) Isopropylbenzene	10.86	105	237		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.18	156	58		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.36	83	58		N.D.	
75) 1,2,3-Trichloropropane	11.39	110	56		N.D.	
76) n-Propylbenzene	11.26	120	68		N.D.	
77) 2-Chlorotoluene	11.48	91	174		N.D.	
78) 4-Chlorotoluene	11.48	91	174		N.D.	
79) 1,3,5-Trimethylbenzene	11.43	105	66		N.D.	
80) tert-Butylbenzene	11.74	119	57		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	335		N.D.	
82) sec-butylbenzene	11.98	105	242		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	73		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	320		N.D.	
86) 1,2-Dichlorobenzene	12.59	146	157		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.32	225	83		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

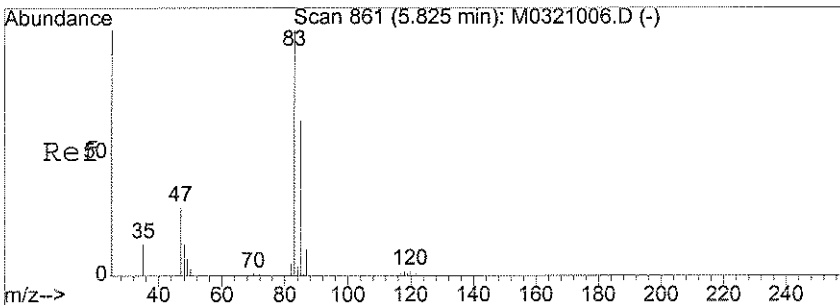
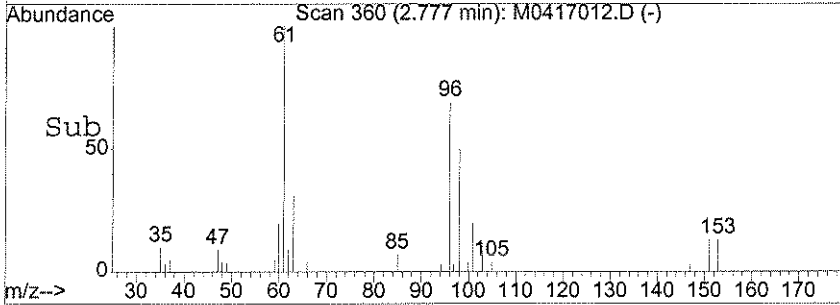
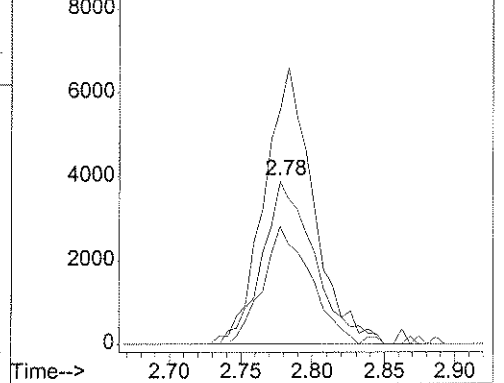


#9  
 1,1-Dichloroethene  
 Concen: 1.69 ug/l  
 RT: 2.78 min Scan# 360  
 Delta R.T. -0.00 min  
 Lab File: M0417012.D  
 Acq: 17 Apr 2007 12:01

Tgt Ion	Resp	Lower	Upper
96	9943		
61	159.5	138.6	178.6
98	65.9	32.2	72.2

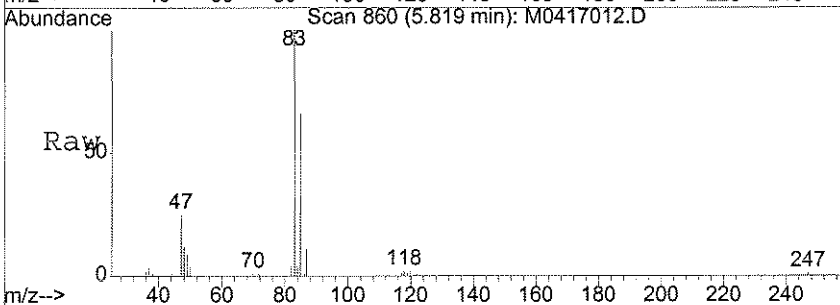


Abundance Ion 96.00 (95.70 to 96.70): M0417012  
 Ion 61.00 (60.70 to 61.70): M0417012  
 Ion 98.00 (97.70 to 98.70): M0417012

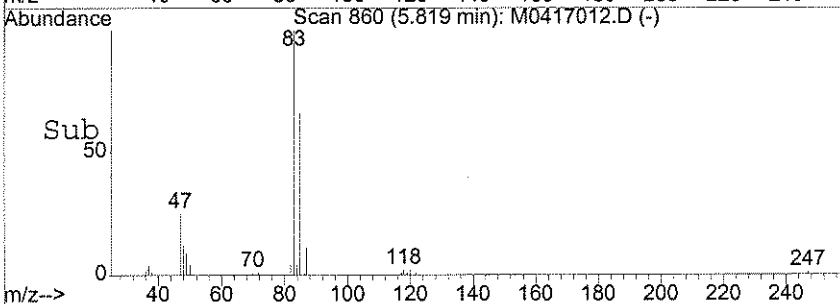
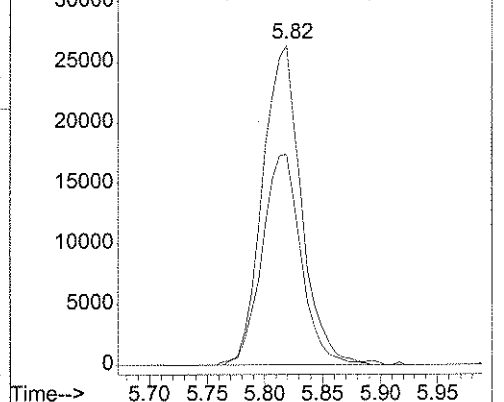


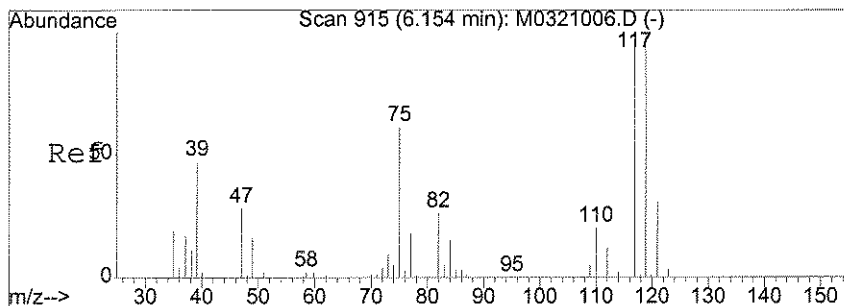
#31  
 Chloroform  
 Concen: 3.99 ug/l  
 RT: 5.82 min Scan# 860  
 Delta R.T. 0.01 min  
 Lab File: M0417012.D  
 Acq: 17 Apr 2007 12:01

Tgt Ion	Resp	Lower	Upper
83	61829		
85	66.1	41.2	81.2



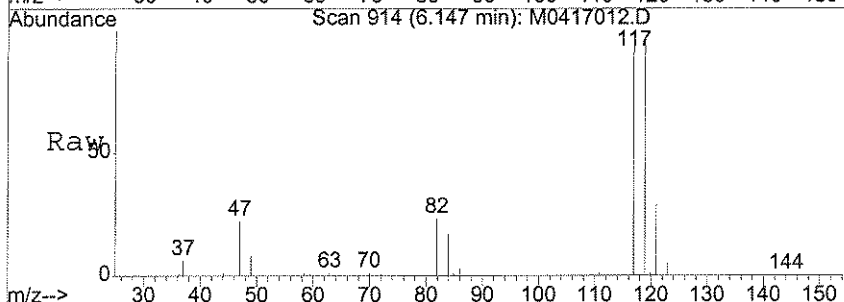
Abundance Ion 83.00 (82.70 to 83.70): M0417012  
 Ion 85.00 (84.70 to 85.70): M0417012



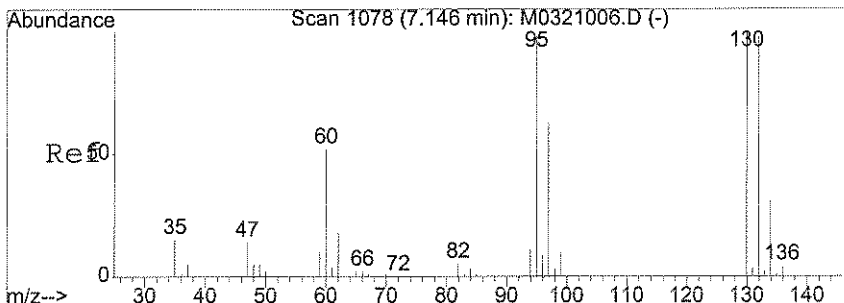
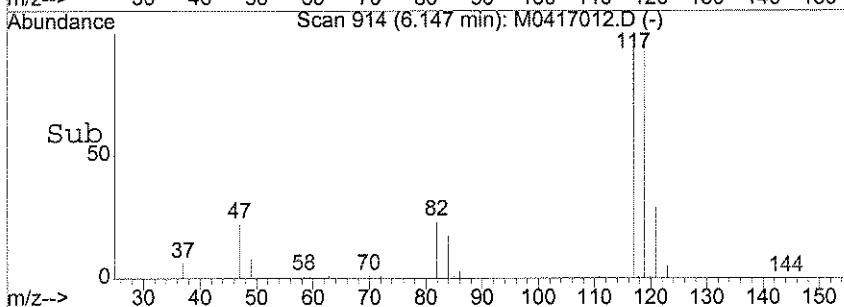
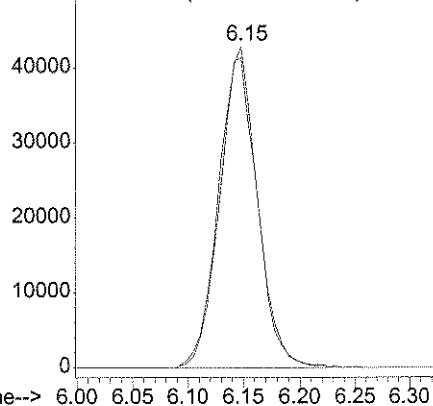


#35  
 Carbon Tetrachloride  
 Concen: 11.26 ug/l  
 RT: 6.15 min Scan# 914  
 Delta R.T. 0.01 min  
 Lab File: M0417012.D  
 Acq: 17 Apr 2007 12:01

Tgt Ion:117 Resp: 103716  
 Ion Ratio Lower Upper  
 117 100  
 119 94.8 78.2 118.2

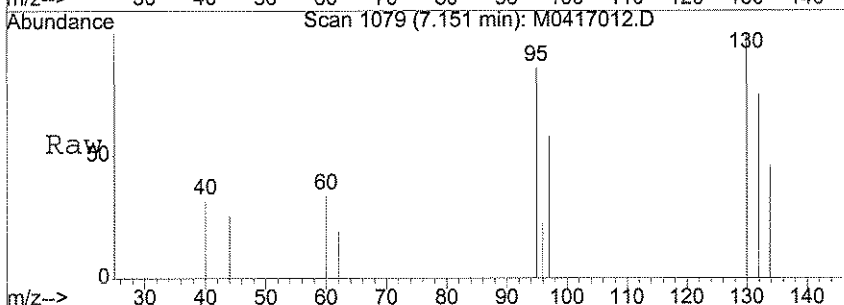


Abundance Ion 117.00 (116.70 to 117.70): M04170  
 50000 Ion 119.00 (118.70 to 119.70): M04170

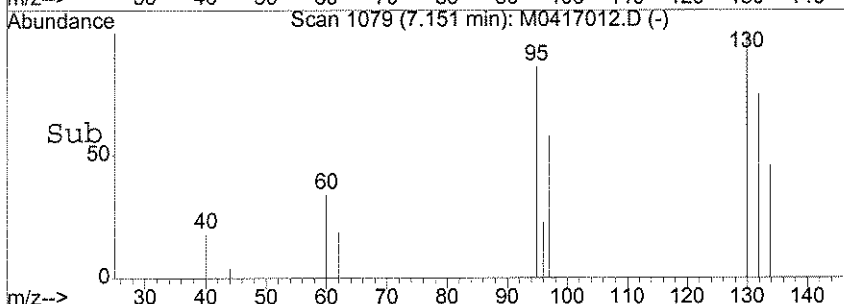
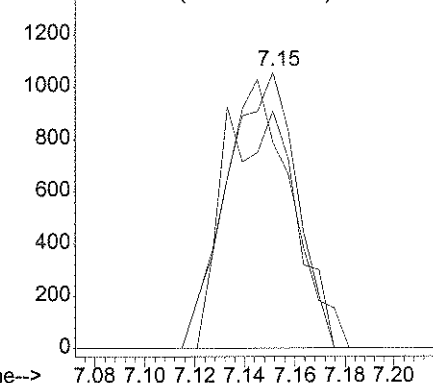


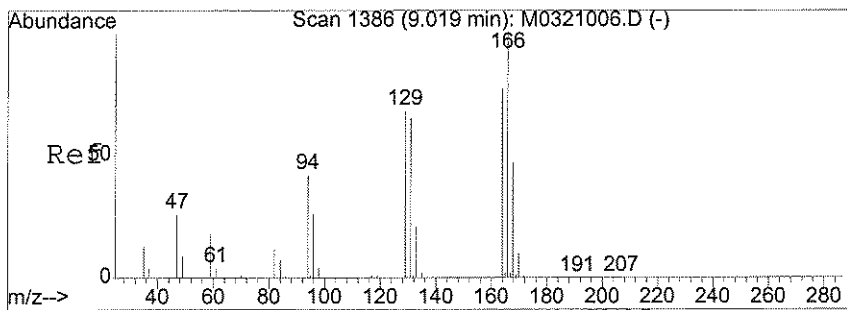
#41  
 Trichloroethene  
 Concen: 0.23 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0417012.D  
 Acq: 17 Apr 2007 12:01

Tgt Ion:130 Resp: 2014  
 Ion Ratio Lower Upper  
 130 100  
 132 96.5 75.0 115.0  
 95 0.0 69.4 109.4#



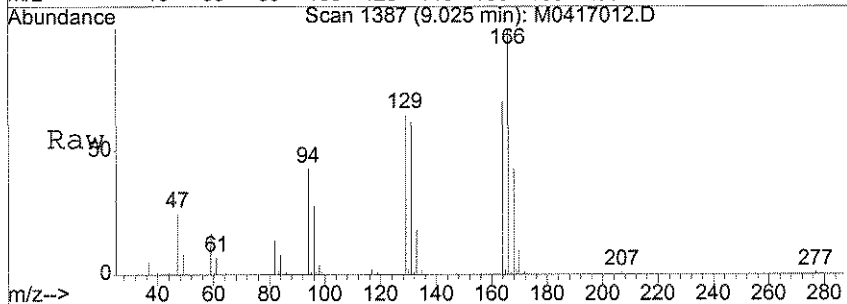
Abundance Ion 130.00 (129.70 to 130.70): M04170  
 1400 Ion 132.00 (131.70 to 132.70): M04170  
 Ion 95.00 (94.70 to 95.70): M0417012.



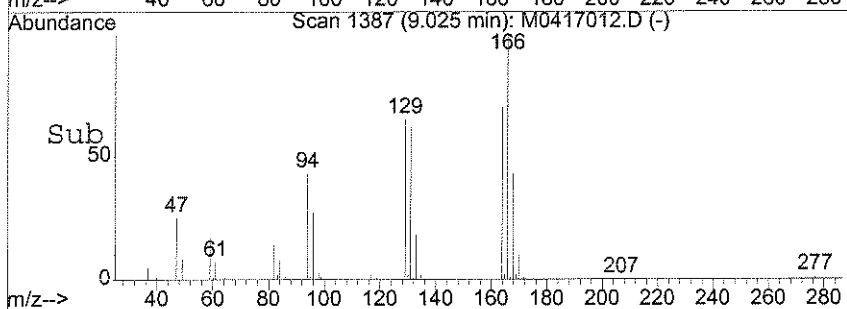
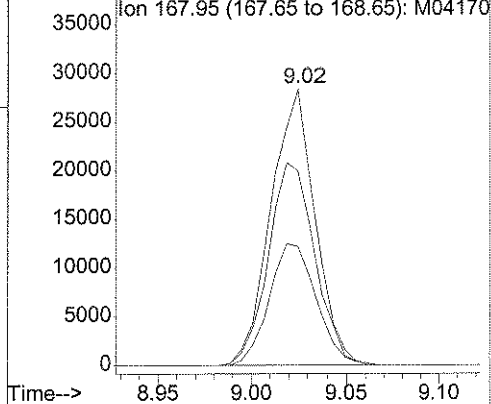


#56  
 Tetrachloroethene  
 Concen: 5.93 ug/l  
 RT: 9.02 min Scan# 1387  
 Delta R.T. -0.00 min  
 Lab File: M0417012.D  
 Acq: 17 Apr 2007 12:01

Tgt Ion	Resp	Lower	Upper
166	46174		
164	77.1	63.3	94.9
168	46.7	39.6	59.4



Abundance  
 Ion 165.95 (165.65 to 166.65): M04170  
 Ion 163.95 (163.65 to 164.65): M04170  
 Ion 167.95 (167.65 to 168.65): M04170



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-5-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-005  
 Lab File ID: M0417013.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 12:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
75-71-8	Dichlorodifluoromethane	0.50		U
74-87-3	Chloromethane	0.50		U
75-01-4	Vinyl chloride	0.50		U
74-83-9	Bromomethane	0.50		U
75-00-3	Chloroethane	0.50		U
75-69-4	Trichlorofluoromethane	0.50		U
75-35-4	1,1-Dichloroethene	1.9		
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50		U
75-09-2	Methylene chloride	0.50		U
1634-04-4	Methyl tert-butyl ether	0.50		U
156-60-5	trans-1,2-Dichloroethene	0.50		U
75-34-3	1,1-Dichloroethane	0.50		U
594-20-7	2,2-Dichloropropane	0.50		U
156-59-2	cis-1,2-Dichloroethene	0.50		U
78-93-3	2-Butanone	5.0		U
74-97-5	Bromochloromethane	0.50		U
67-66-3	Chloroform	4.8		
71-55-6	1,1,1-Trichloroethane	0.50		U
56-23-5	Carbon tetrachloride	14		
563-58-6	1,1-Dichloropropene	0.50		U
71-43-2	Benzene	0.50		U
107-06-2	1,2-Dichloroethane	0.50		U
79-01-6	Trichloroethene	0.26		J
78-87-5	1,2-Dichloropropane	0.50		U
74-95-3	Dibromomethane	0.50		U
75-27-4	Bromodichloromethane	0.50		U
10061-01-	cis-1,3-Dichloropropene	0.50		U
108-10-1	4-Methyl-2-pentanone	5.0		U
108-88-3	Toluene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-5-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 12:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	7.4	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-5-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-005

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 12:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

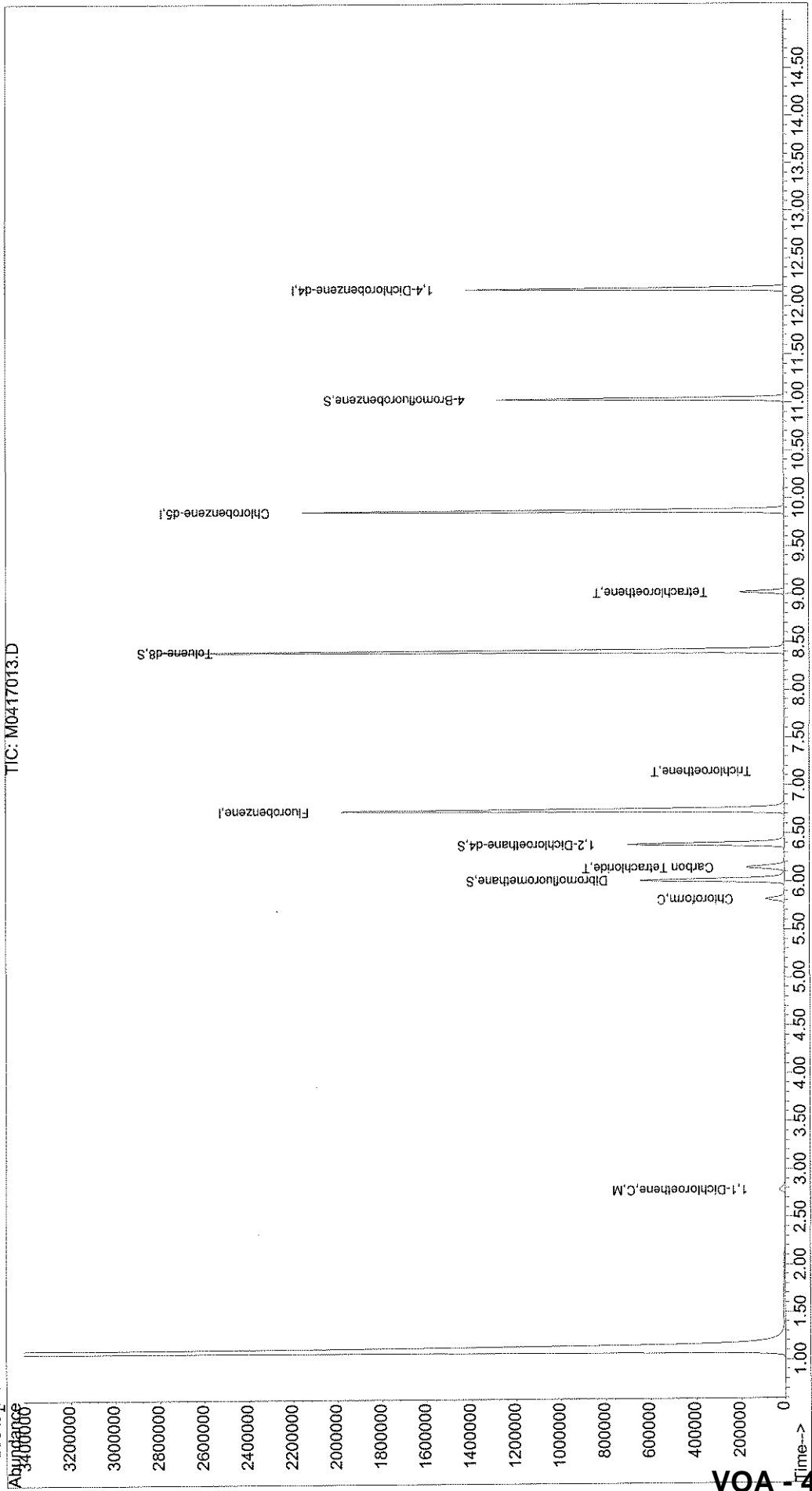
Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417013.D  
Acq On : 17 Apr 2007 12:25  
Sample : JPL32-005  
Misc : #2 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:12 2007

Vial: 58  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00

Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA-41

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417013.D  
 Acq On : 17 Apr 2007 12:25  
 Sample : JPL32-005  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:12 2007

Vial: 58  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1907297	50.00	ug/l	0.00 97.71%
50) Chlorobenzene-d5	9.87	82	639634	50.00	ug/l	0.00 90.95%
70) 1,4-Dichlorobenzene-d4	12.18	152	339016	50.00	ug/l	0.00 87.10%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	449342	49.94	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.88%
37) 1,2-Dichloroethane-d4	6.39	65	521134	50.59	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.18%
51) Toluene-d8	8.41	98	1854956	52.51	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.02%
72) 4-Bromofluorobenzene	11.04	95	388951	55.62	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.24%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.83	96	65	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.78	96	11299	1.94	ug/l	98
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.97	43	200	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	3.03	108	57	N.D.		
14) Carbon Disulfide	3.03	76	1957	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.88	96	68	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.55	63	245	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417013.D 8260B.M Wed Apr 25 09:13:09 2007

*jo 04/25/07*  
 VOA-42 Page 1

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417013.D  
 Acq On : 17 Apr 2007 12:25  
 Sample : JPL32-005  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:12 2007

Vial: 58  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	5.41	77	57	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.43	43	66	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	5.58	128	56	N.D.		
30) Methacrylonitrile	5.81	41	59	N.D.		
31) Chloroform	5.81	83	73779	4.82	ug/l	94
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	5.96	56	61	N.D.		
35) Carbon Tetrachloride	6.14	117	122850	13.50	ug/l	99
36) 1,1-Dichloropropene	6.06	75	61	N.D.		
38) Benzene	6.41	78	1253	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.	d	
41) Trichloroethene	7.15	130	2226	0.26	ug/l #	51
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	7.56	63	55	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	7.67	41	61	N.D.		
46) Bromodichloromethane	7.71	83	134	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	8.05	75	66	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	8.48	92	745	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	8.93	97	86	N.D.		
56) Tetrachloroethene	9.03	166	56849	7.40	ug/l	98
57) 1,3-Dichloropropane	8.89	76	56	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.	d	
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	9.89	112	134	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.99	91	921	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.	d	
66) o-xylene	10.52	106	141	N.D.		
67) Styrene	10.53	104	213	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417013.D 8260B.M Wed Apr 25 09:13:09 2007

*[Handwritten Signature]*  
 Page 2  
**VOA-43**

Quantitation Report

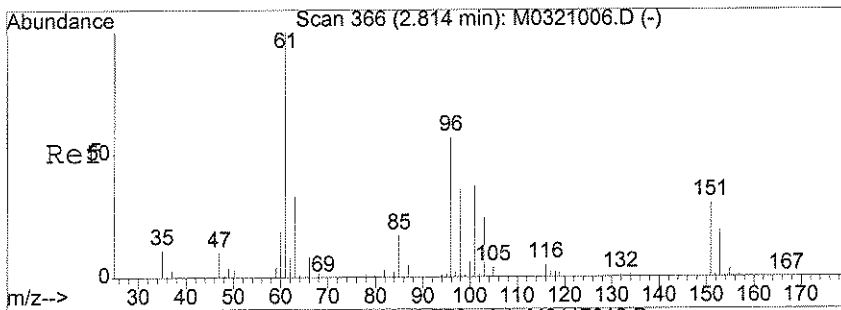
Data File : X:\MSVOA\MOBY\041707\M0417013.D  
 Acq On : 17 Apr 2007 12:25  
 Sample : JPL32-005  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:12 2007

Vial: 58  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

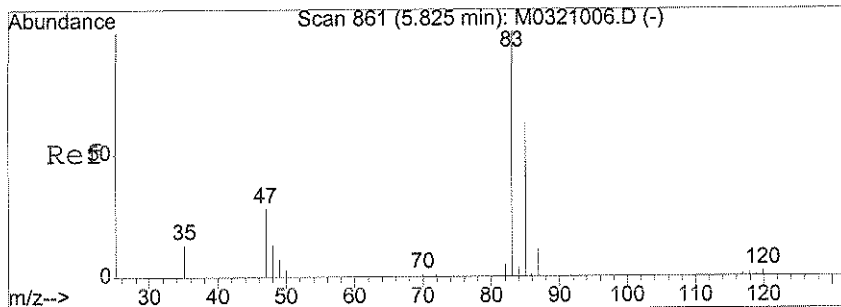
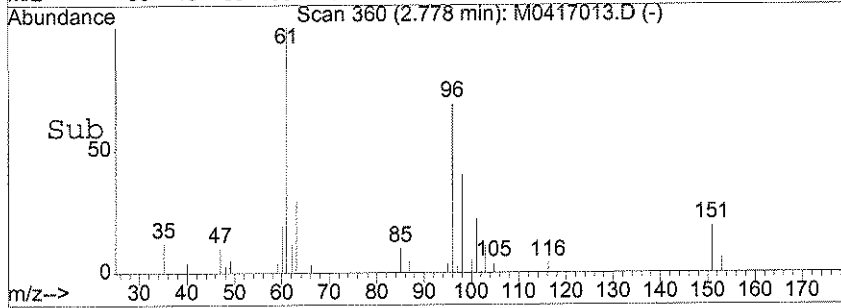
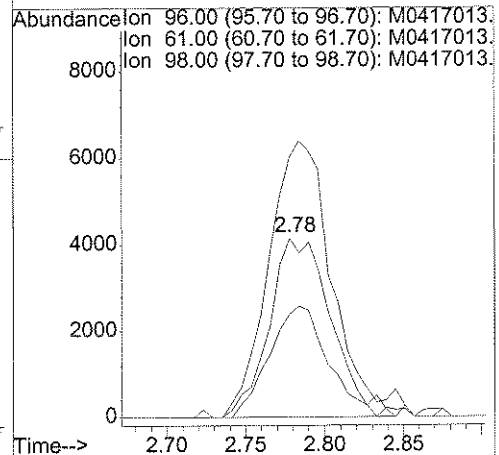
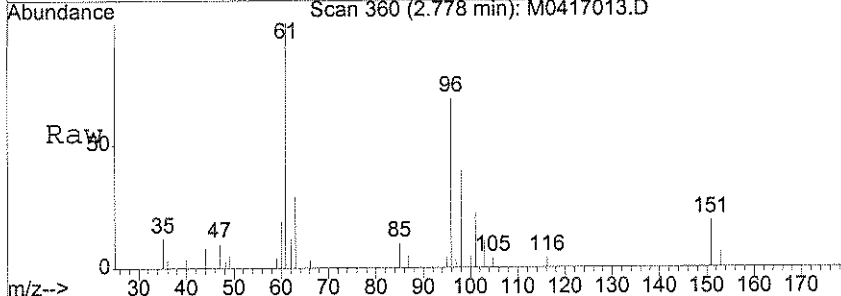
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	141		N.D.	
69) Isopropylbenzene	10.86	105	321		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	62		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	370		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	141		N.D.	
76) n-Propylbenzene	11.26	120	58		N.D.	
77) 2-Chlorotoluene	11.37	91	166		N.D.	
78) 4-Chlorotoluene	11.49	91	83		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	209		N.D.	
80) tert-Butylbenzene	11.76	119	161		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	229		N.D.	
82) sec-butylbenzene	11.98	105	324		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	71		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.12	146	71		N.D.	
86) 1,2-Dichlorobenzene	12.59	146	71		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.01	225	66		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



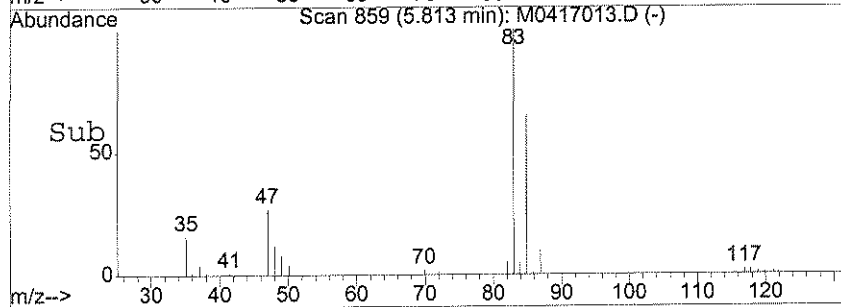
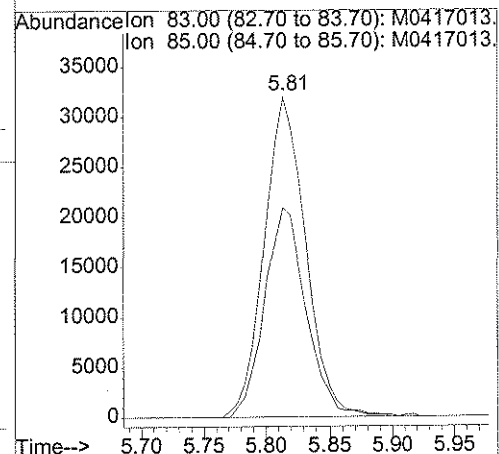
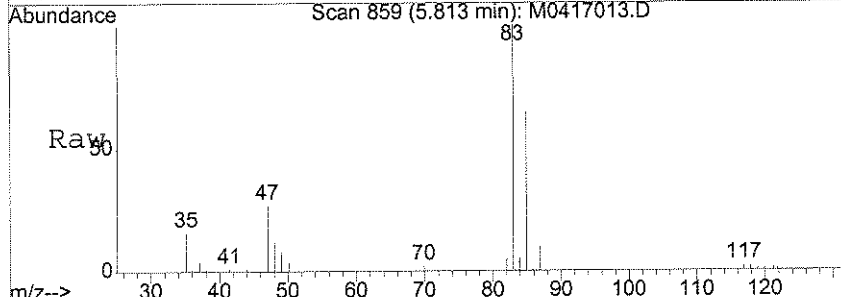
#9  
 1,1-Dichloroethene  
 Concen: 1.94 ug/l  
 RT: 2.78 min Scan# 360  
 Delta R.T. 0.00 min  
 Lab File: M0417013.D  
 Acq: 17 Apr 2007 12:25

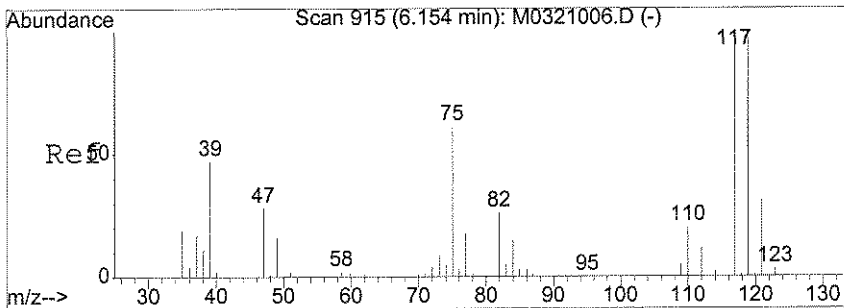
Tgt Ion	Resp	Lower	Upper
96	11299		
61	158.7	138.6	178.6
98	58.4	32.2	72.2



#31  
 Chloroform  
 Concen: 4.82 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0417013.D  
 Acq: 17 Apr 2007 12:25

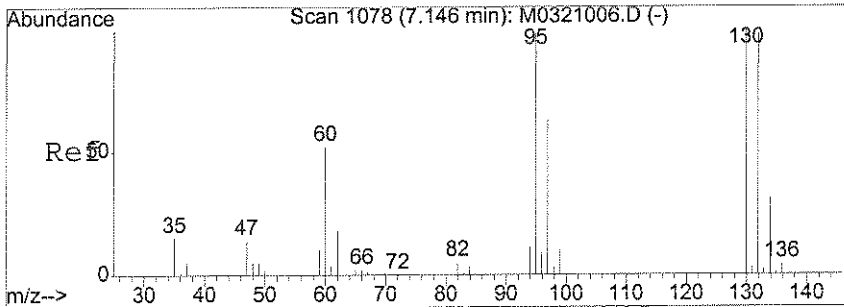
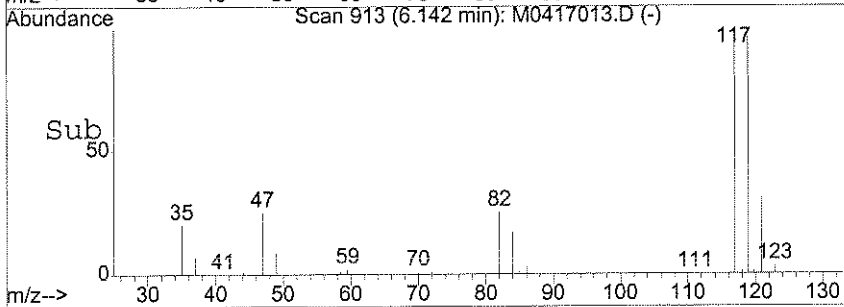
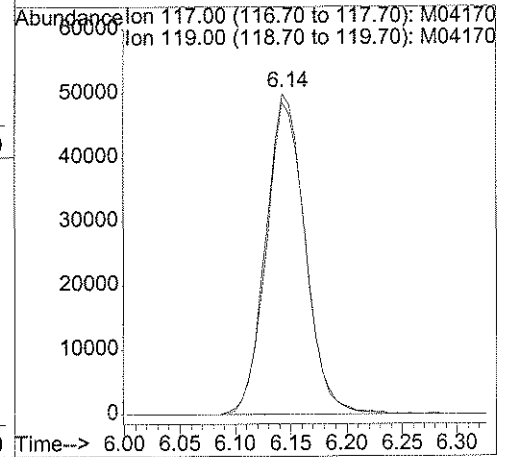
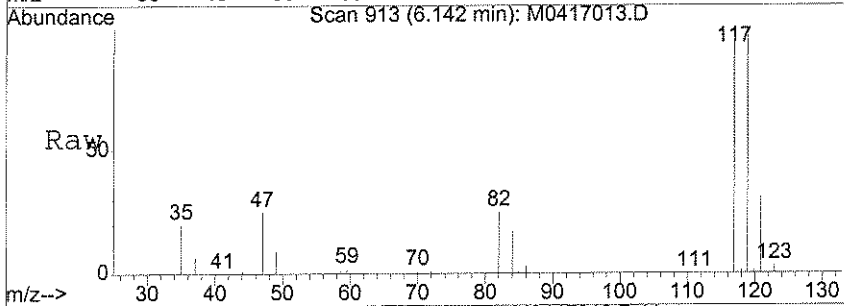
Tgt Ion	Resp	Lower	Upper
83	73779		
85	65.4	41.2	81.2





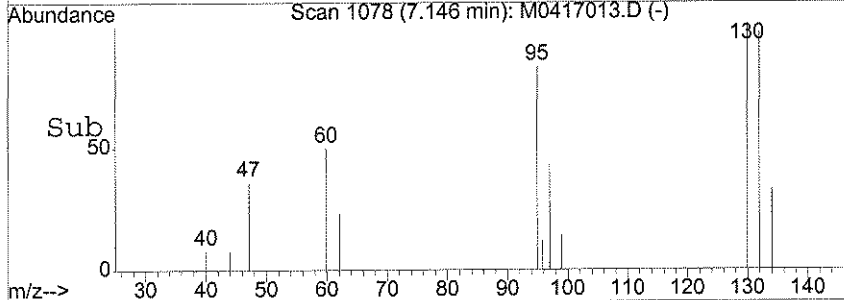
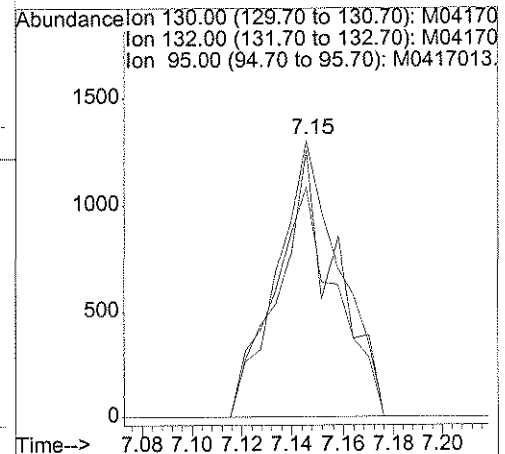
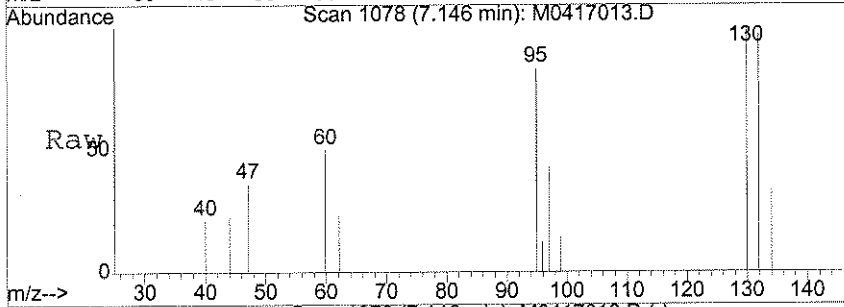
#35  
 Carbon Tetrachloride  
 Concen: 13.50 ug/l  
 RT: 6.14 min Scan# 913  
 Delta R.T. 0.00 min  
 Lab File: M0417013.D  
 Acq: 17 Apr 2007 12:25

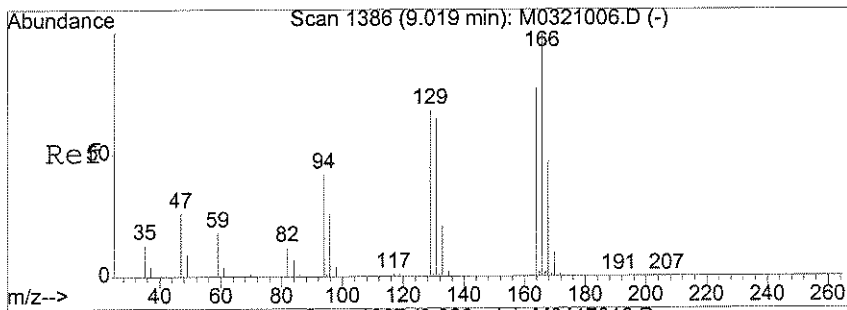
Tgt Ion	Resp	Lower	Upper
117	122850	100	
119	97.0	78.2	118.2



#41  
 Trichloroethene  
 Concen: 0.26 ug/l  
 RT: 7.15 min Scan# 1078  
 Delta R.T. 0.00 min  
 Lab File: M0417013.D  
 Acq: 17 Apr 2007 12:25

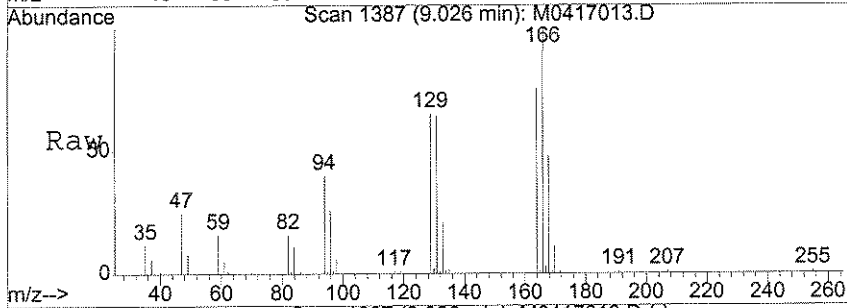
Tgt Ion	Resp	Lower	Upper
130	2226	100	
132	89.3	75.0	115.0
95	0.0	69.4	109.4#



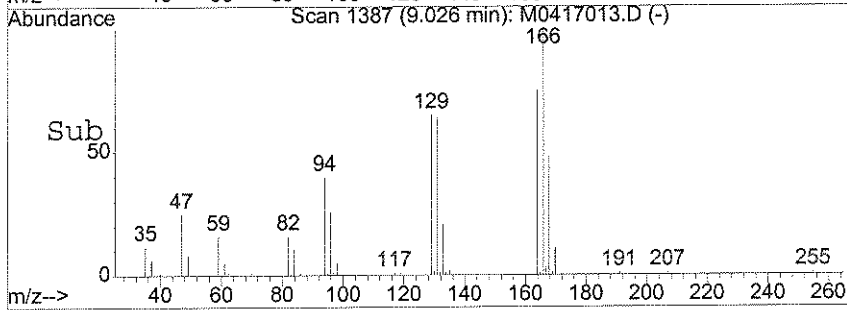
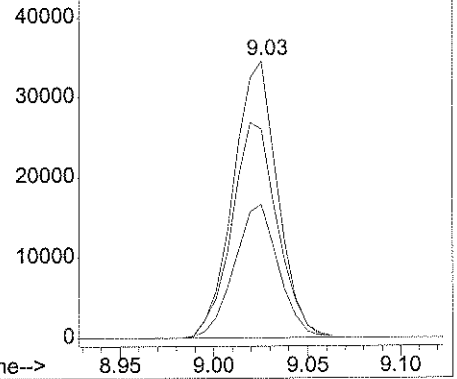


#56  
 Tetrachloroethene  
 Concen: 7.40 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0417013.D  
 Acq: 17 Apr 2007 12:25

Tgt Ion	Resp	Lower	Upper
166	56849		
164	80.1	63.3	94.9
168	48.0	39.6	59.4



Abundance Ion 165.95 (165.65 to 166.65): M04170  
 Ion 163.95 (163.65 to 164.65): M04170  
 Ion 167.95 (167.65 to 168.65): M04170





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-006  
 Lab File ID: M0417014.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 12:49  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417014.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 12:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.1	
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417014.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 12:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

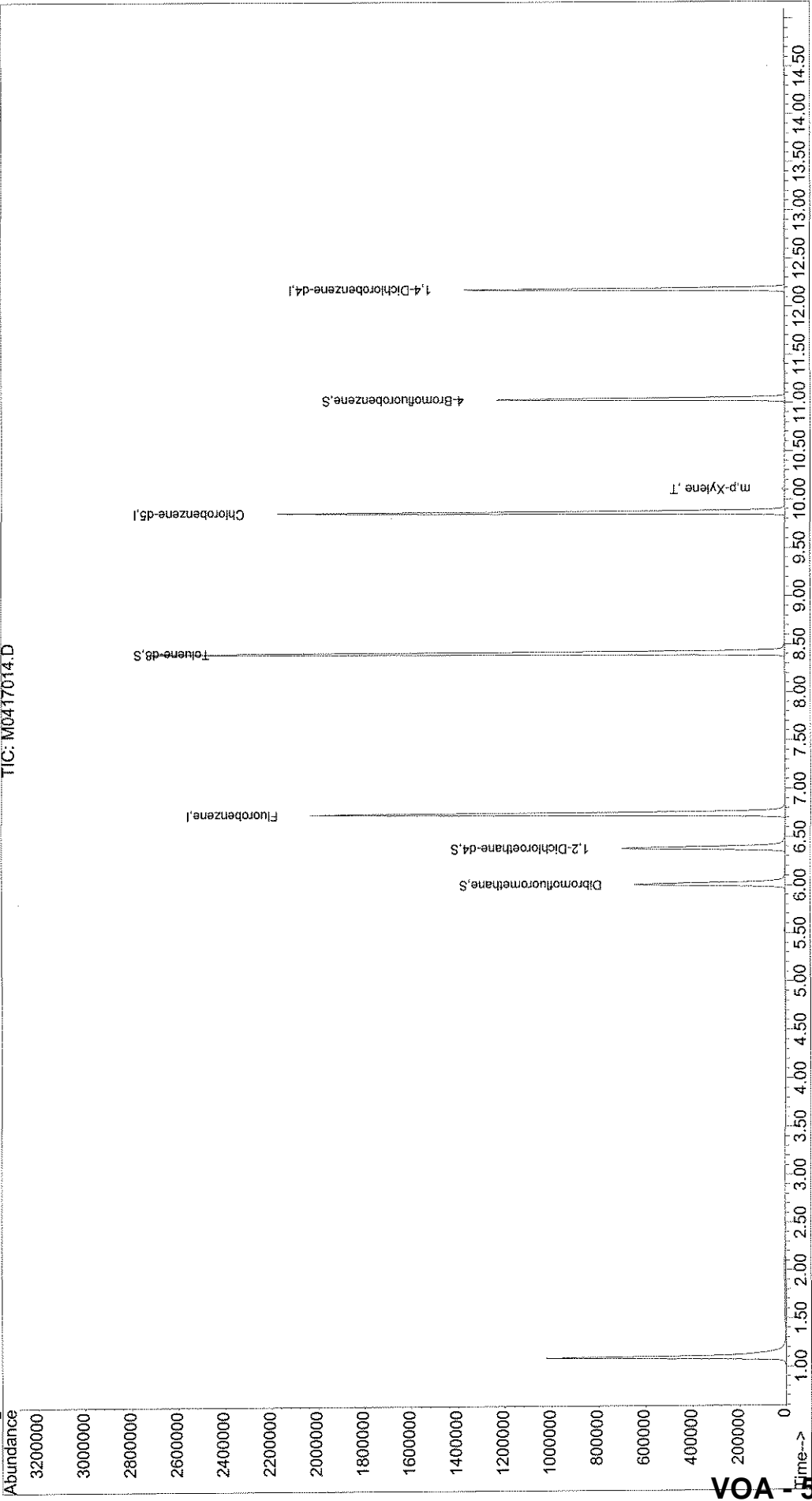
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417014.D Vial: 59  
Acq On : 17 Apr 2007 12:49 Operator: DGA  
Sample : JPL32-006 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:14 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA-51

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417014.D  
 Acq On : 17 Apr 2007 12:49  
 Sample : JPL32-006  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:14 2007

Vial: 59  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1905788	50.00	ug/l	0.00 97.64%
50) Chlorobenzene-d5	9.87	82	635982	50.00	ug/l	0.00 90.43%
70) 1,4-Dichlorobenzene-d4	12.18	152	337002	50.00	ug/l	0.00 86.58%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	449921	50.04	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery =	100.08%		
37) 1,2-Dichloroethane-d4	6.39	65	520939	50.61	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery =	101.22%		
51) Toluene-d8	8.41	98	1839331	52.36	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery =	104.72%		
72) 4-Bromofluorobenzene	11.04	95	381137	54.83	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery =	109.66%		

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.89	96	130	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.02	76	92	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417014.D  
 Acq On : 17 Apr 2007 12:49  
 Sample : JPL32-006  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:14 2007

Vial: 59  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.76	43	75		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.38	96	58		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	d
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.80	83	102		N.D.	
32) 1,1,1-Trichloroethane	5.95	97	58		N.D.	
34) Cyclohexane	5.99	56	213		N.D.	
35) Carbon Tetrachloride	6.28	117	74		N.D.	
36) 1,1-Dichloropropene	6.19	75	56		N.D.	
38) Benzene	6.41	78	861		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	7.32	83	62		N.D.	
43) 1,2-Dichloropropane	7.43	63	72		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.60	83	60		N.D.	
47) 2-Chloroethyl vinyl ether	8.05	63	58		N.D.	
48) cis-1,3-Dichloropropene	8.17	75	67		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	690		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.88	97	125		N.D.	
56) Tetrachloroethene	9.02	166	414		N.D.	
57) 1,3-Dichloropropane	9.04	76	55		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	477		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	3009		N.D.	
65) m,p-Xylene	10.11	106	3219	1.09	ug/l	99
66) o-xylene	10.51	106	1346		N.D.	
67) Styrene	10.59	104	58		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417014.D 8260B.M Wed Apr 25 09:14:34 2007

*[Handwritten Signature]*  
 Page 2  
 VOA - 53

Quantitation Report

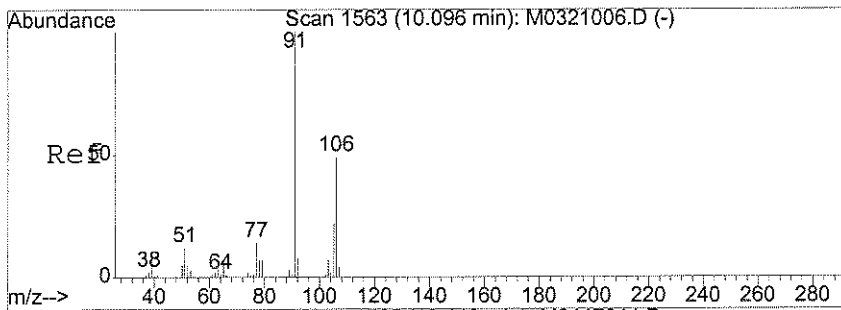
Data File : X:\MSVOA\MOBY\041707\M0417014.D  
 Acq On : 17 Apr 2007 12:49  
 Sample : JPL32-006  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:14 2007

Vial: 59  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

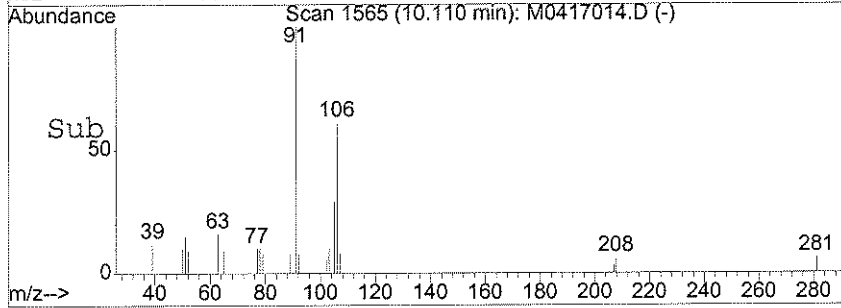
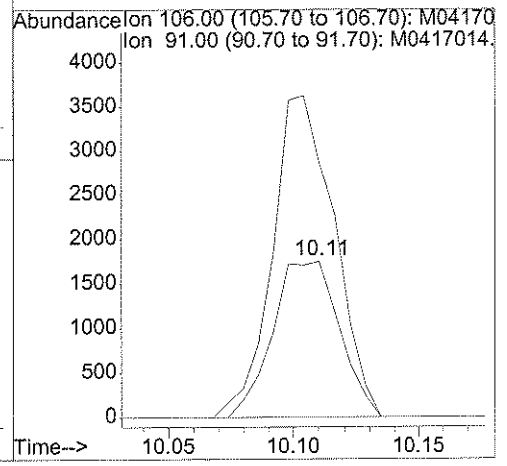
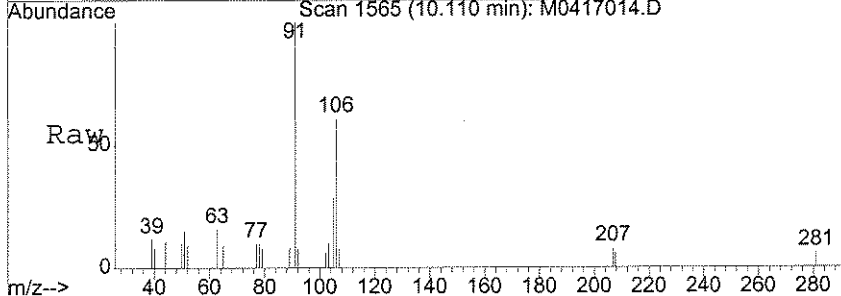
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	163		N.D.	
69) Isopropylbenzene	10.86	105	213		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.04	156	61		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	391		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	62		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.35	91	61		N.D.	
78) 4-Chlorotoluene	11.47	91	257		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	75		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	180		N.D.	
82) sec-butylbenzene	11.98	105	198		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	192		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	192		N.D.	
86) 1,2-Dichlorobenzene	12.30	146	85		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.33	225	66		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



#65  
 m,p-Xylene  
 Concen: 1.09 ug/l  
 RT: 10.11 min Scan# 1565  
 Delta R.T. 0.00 min  
 Lab File: M0417014.D  
 Acq: 17 Apr 2007 12:49

Tgt Ion	Resp	Lower	Upper
106	100		
91	191.5	169.6	209.6





1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-007  
 Lab File ID: M0417015.D  
 Date Collected: 04/05/2007  
 Date/Time Analyzed: 04/17/2007 13:13  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 13:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/05/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 13:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

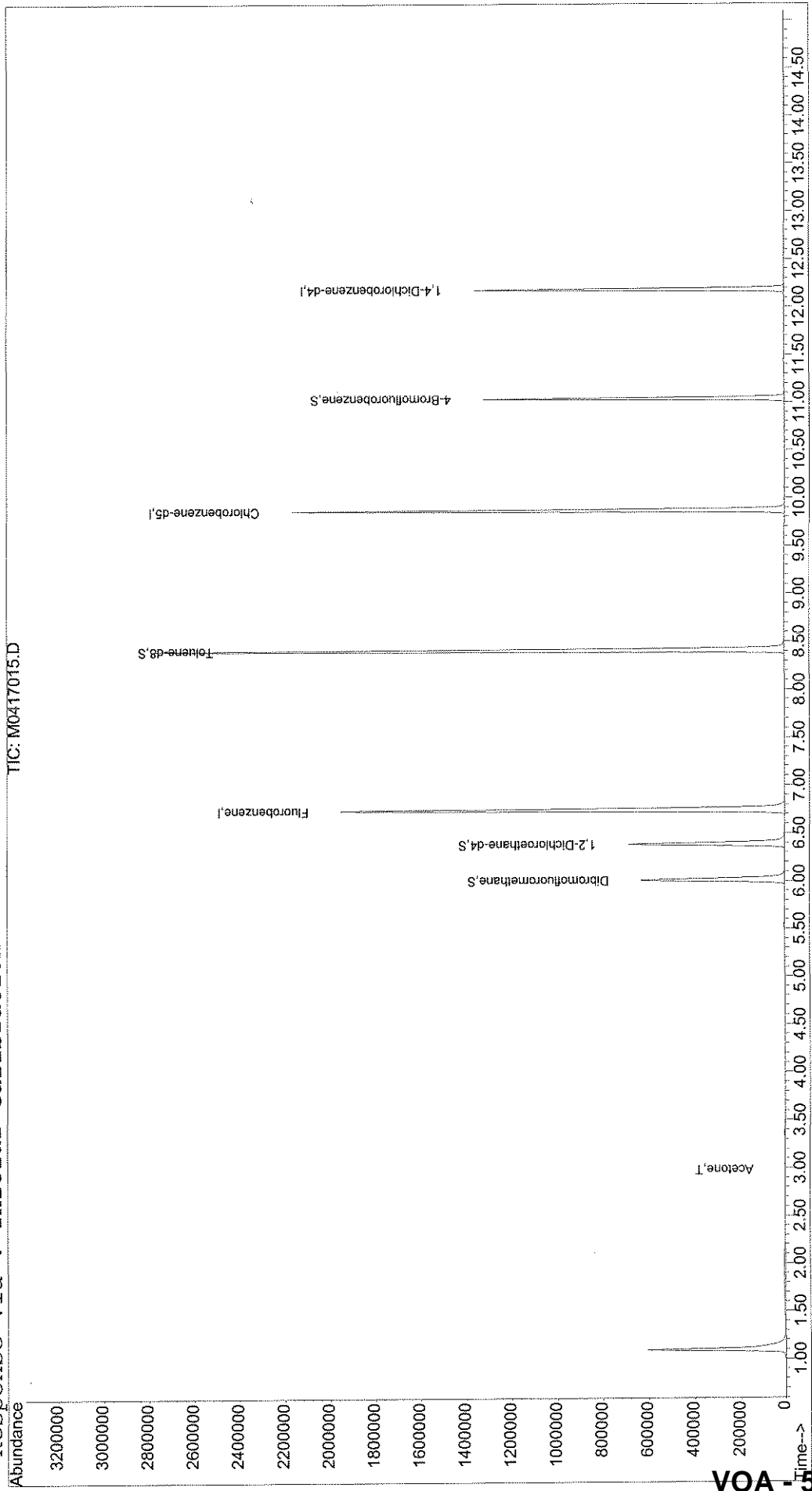
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417015.D Vial: 60  
Acq On : 17 Apr 2007 13:13 Operator: DGA  
Sample : JPL32-007 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:15 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417015.D  
 Acq On : 17 Apr 2007 13:13  
 Sample : JPL32-007  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:15 2007

Vial: 60  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1879222	50.00	ug/l	0.00 96.28%
50) Chlorobenzene-d5	9.87	82	633240	50.00	ug/l	0.00 90.04%
70) 1,4-Dichlorobenzene-d4	12.18	152	335402	50.00	ug/l	0.00 86.17%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	447059	50.43	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.86%
37) 1,2-Dichloroethane-d4	6.39	65	504319	49.69	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.38%
51) Toluene-d8	8.41	98	1825210	52.19	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.38%
72) 4-Bromofluorobenzene	11.04	95	384702	55.61	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.22%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.45	50	61	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.84	96	69	N.D.		
6) Chloroethane	1.94	64	95	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.83	96	64	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.99	43	2554	1.49	ug/l	94
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	2361	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	3.47	84	1924	N.D.		
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417015.D 8260B.M Wed Apr 25 09:15:44 2007

*J* 04/25/07  
 VOA-60 Page 1

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417015.D  
 Acq On : 17 Apr 2007 13:13  
 Sample : JPL32-007  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:15 2007

Vial: 60  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.69	43	59		N.D.	
24) Chloroprene	4.66	53	66		N.D.	
25) 2,2-Dichloropropane	5.36	77	93		N.D.	
26) cis-1,2-Dichloroethene	5.42	96	112		N.D.	
27) 2-Butanone	5.42	43	69		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.73	41	61		N.D.	
31) Chloroform	5.83	83	76		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.19	75	62		N.D.	
38) Benzene	6.41	78	472		N.D.	
39) 1,2-Dichloroethane	6.39	62	275		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	7.38	63	61		N.D.	
44) Dibromomethane	7.59	93	74		N.D.	
45) Methyl methacrylate	7.40	41	135		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.27	75	62		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	213		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.95	97	57		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	9.54	107	71		N.D.	
61) Chlorobenzene	9.89	112	69		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	331		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.27	104	65		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417015.D  
 Acq On : 17 Apr 2007 13:13  
 Sample : JPL32-007  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:15 2007

Vial: 60  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	66		N.D.	
69) Isopropylbenzene	10.87	105	155		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	466		N.D.	
75) 1,2,3-Trichloropropane	11.20	110	66		N.D.	
76) n-Propylbenzene	11.28	120	57		N.D.	
77) 2-Chlorotoluene	11.36	91	122		N.D.	
78) 4-Chlorotoluene	11.48	91	85		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	136		N.D.	
80) tert-Butylbenzene	11.76	119	119		N.D.	
81) 1,2,4-Trimethylbenzene	11.83	105	65		N.D.	
82) sec-butylbenzene	11.99	105	293		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	55		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	128		N.D.	
86) 1,2-Dichlorobenzene	12.52	146	57		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417015.D 8260B.M Wed Apr 25 09:15:45 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 13:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 13:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-008

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 13:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

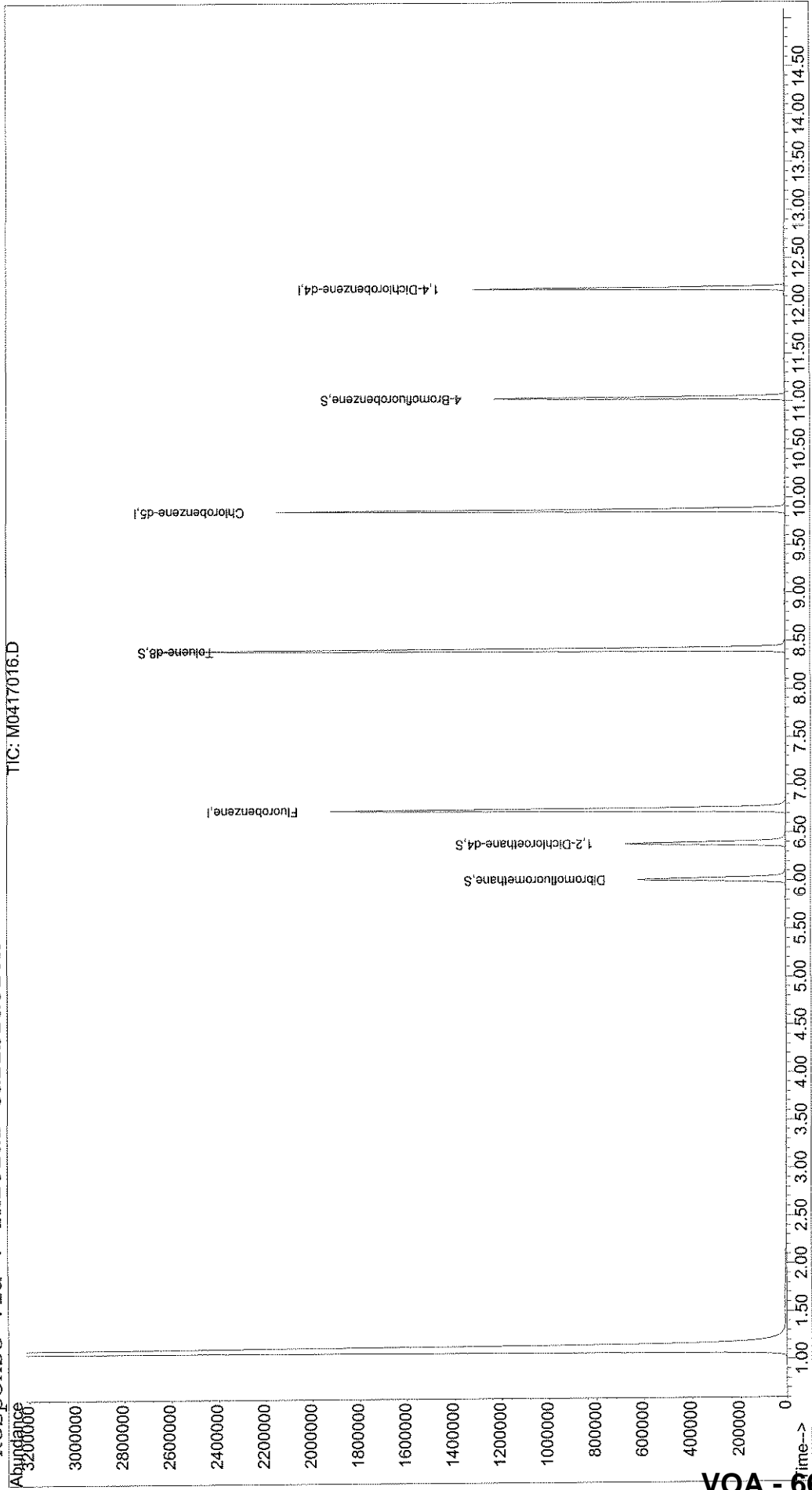
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417016.D  
Acq On : 17 Apr 2007 13:37  
Sample : JPL32-008  
Misc : #3 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:22 2007  
Vial: 61  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417016.D  
 Acq On : 17 Apr 2007 13:37  
 Sample : JPL32-008  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:22 2007

Vial: 61  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1844658	50.00	ug/l	0.00 94.50%
50) Chlorobenzene-d5	9.87	82	617447	50.00	ug/l	0.00 87.79%
70) 1,4-Dichlorobenzene-d4	12.19	152	323413	50.00	ug/l	0.00 83.09%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	437689	50.29	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.58%
37) 1,2-Dichloroethane-d4	6.39	65	503208	50.51	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.02%
51) Toluene-d8	8.41	98	1794149	52.61	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.22%
72) 4-Bromofluorobenzene	11.04	95	376193	56.39	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	112.78%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.84	96	59	N.D.	
6) Chloroethane	1.95	64	72	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	3120	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.41	43	55	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	4.54	63	1468	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417016.D  
 Acq On : 17 Apr 2007 13:37  
 Sample : JPL32-008  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:22 2007

Vial: 61  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	5.28	77	55		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.50	43	65		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.82	83	2325		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	6.01	56	66		N.D.	
35) Carbon Tetrachloride	6.19	117	57		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	641		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.14	130	653		N.D.	
42) Methylcyclohexane	7.29	83	186		N.D.	
43) 1,2-Dichloropropane	7.39	63	61		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.42	41	64		N.D.	
46) Bromodichloromethane	7.73	83	265		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	663		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.94	97	57		N.D.	
56) Tetrachloroethene	9.02	166	994		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	449		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.01	131	61		N.D.	
64) Ethylbenzene	9.99	91	452		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	10.50	106	228		N.D.	
67) Styrene	10.51	104	288		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417016.D  
 Acq On : 17 Apr 2007 13:37  
 Sample : JPL32-008  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:22 2007

Vial: 61  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.04	105	424		N.D.	
71) trans-1,4-Dichloro-2-buten	11.07	53	60		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.30	83	56		N.D.	
75) 1,2,3-Trichloropropane	11.25	110	59		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.47	91	70		N.D.	
78) 4-Chlorotoluene	11.49	91	60		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	66		N.D.	
80) tert-Butylbenzene	11.75	119	56		N.D.	
81) 1,2,4-Trimethylbenzene	11.99	105	180		N.D.	
82) sec-butylbenzene	11.99	105	180		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	1273		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.12	146	1273		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	535		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-009  
 Lab File ID: M0417017.D  
 Date Collected: 04/06/2007  
 Date/Time Analyzed: 04/17/2007 14:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-009  
 Lab File ID: M0417017.D  
 Date Collected: 04/06/2007  
 Date/Time Analyzed: 04/17/2007 14:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

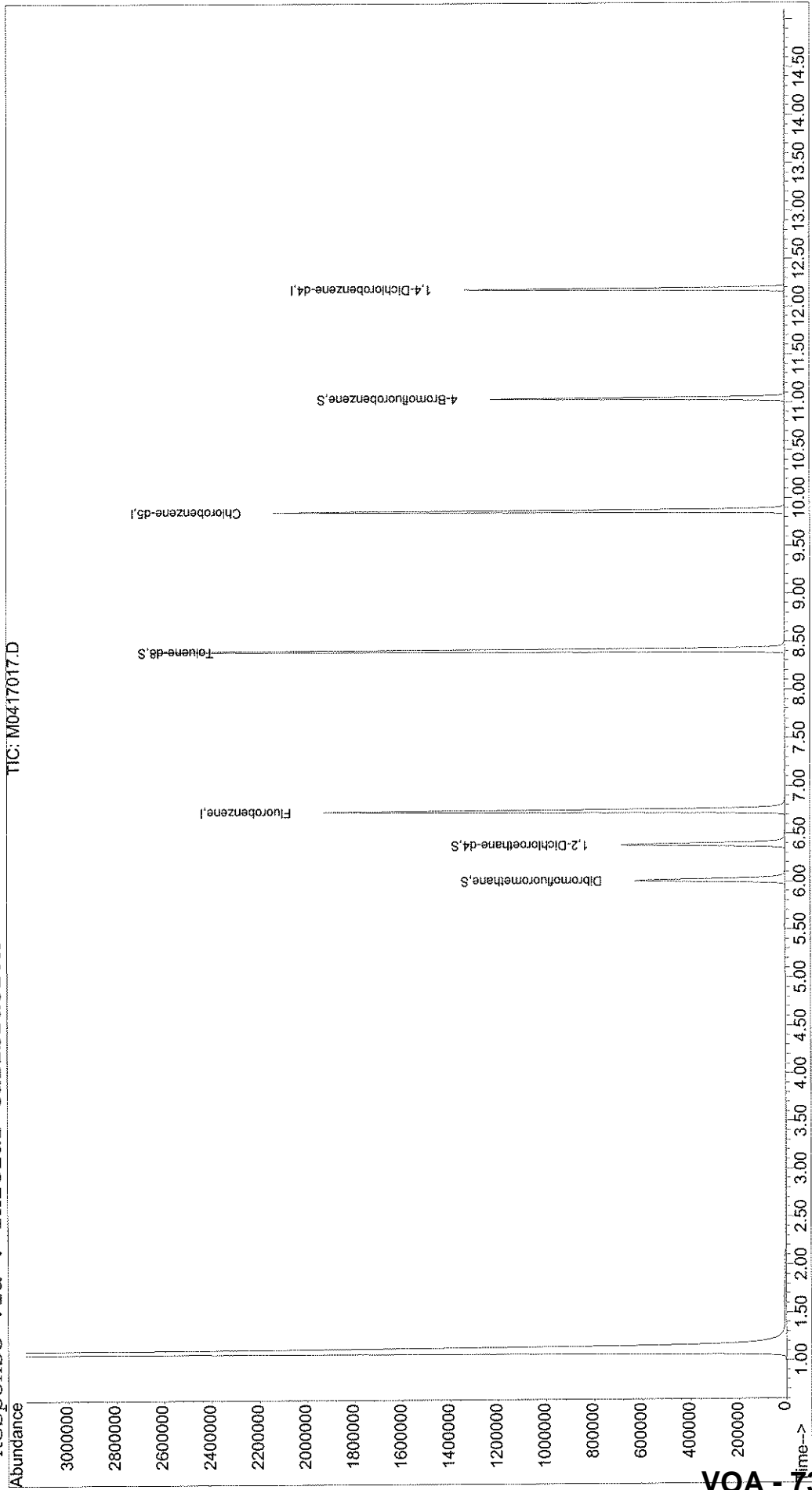
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417017.D Vial: 62  
Acq On : 17 Apr 2007 14:01 Operator: DGA  
Sample : JPL32-009 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:16 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA - 73

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417017.D  
 Acq On : 17 Apr 2007 14:01  
 Sample : JPL32-009  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:16 2007

Vial: 62  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1824404	50.00	ug/l	0.00 93.47%
50) Chlorobenzene-d5	9.87	82	609414	50.00	ug/l	0.00 86.65%
70) 1,4-Dichlorobenzene-d4	12.19	152	324094	50.00	ug/l	0.00 83.27%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	436782	50.75	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.50%
37) 1,2-Dichloroethane-d4	6.39	65	506452	51.40	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.80%
51) Toluene-d8	8.41	98	1774650	52.73	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.46%
72) 4-Bromofluorobenzene	11.04	95	371213	55.53	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.06%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	1.52	62	59	N.D.	
5) Bromomethane	1.87	96	94	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.97	43	172	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	1199	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	3.30	40	62	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	3.95	96	78	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	4.54	63	356	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417017.D 8260B.M Wed Apr 25 09:16:58 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417017.D  
 Acq On : 17 Apr 2007 14:01  
 Sample : JPL32-009  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:16 2007

Vial: 62  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.54	43	227		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.82	83	764		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.22	75	115		N.D.	
38) Benzene	6.40	78	988		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.14	130	305		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.40	93	56		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.73	83	129		N.D.	
47) 2-Chloroethyl vinyl ether	8.13	63	61		N.D.	
48) cis-1,3-Dichloropropene	8.36	75	61		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	640		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.74	97	56		N.D.	
56) Tetrachloroethene	9.02	166	562		N.D.	
57) 1,3-Dichloropropane	9.24	76	62		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	414		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.00	91	563		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	10.49	106	66		N.D.	
67) Styrene	10.52	104	68		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417017.D 8260B.M Wed Apr 25 09:16:59 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417017.D  
 Acq On : 17 Apr 2007 14:01  
 Sample : JPL32-009  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:16 2007

Vial: 62  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	282		N.D.	
69) Isopropylbenzene	10.87	105	58		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	447		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	187		N.D.	
76) n-Propylbenzene	11.45	120	85		N.D.	
77) 2-Chlorotoluene	11.27	91	208		N.D.	
78) 4-Chlorotoluene	11.48	91	168		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	61		N.D.	
80) tert-Butylbenzene	11.77	119	227		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	124		N.D.	
82) sec-butylbenzene	11.98	105	137		N.D.	
83) 1,3-Dichlorobenzene	12.13	146	246		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.22	146	180		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	230		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.36	157	55		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.53	225	60		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

(#) = qualifier out of range (m) = manual integration  
 M0417017.D 8260B.M Wed Apr 25 09:16:59 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.28	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	1.0	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-22-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-010

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

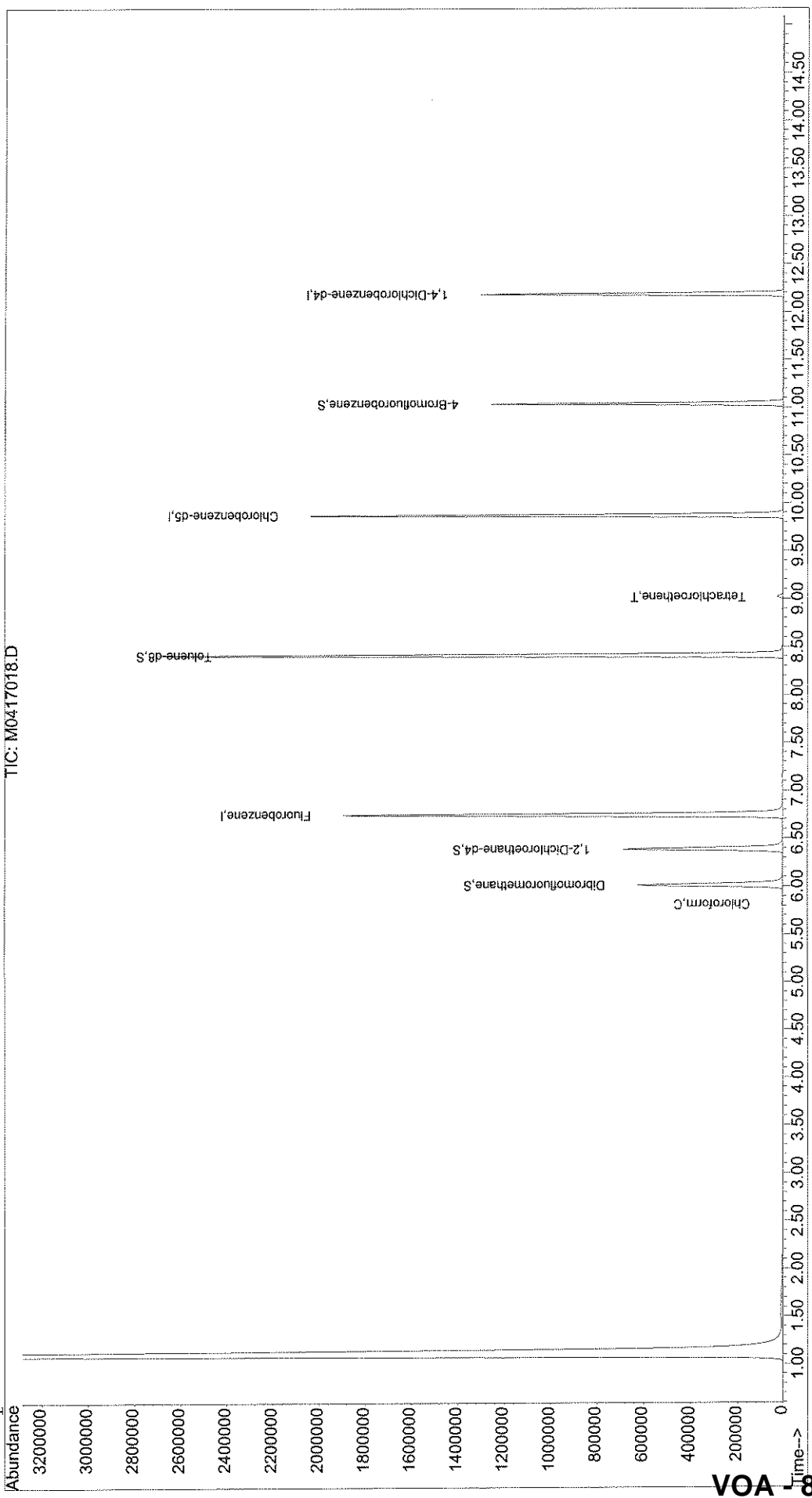
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417018.D Vial: 63  
Acq On : 17 Apr 2007 14:25 Operator: DGA  
Sample : JPL32-010 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:24 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA-80

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417018.D  
 Acq On : 17 Apr 2007 14:25  
 Sample : JPL32-010  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:24 2007

Vial: 63  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	1836817	50.00	ug/l	0.00	94.10%
50) Chlorobenzene-d5	9.86	82	606336	50.00	ug/l	0.00	86.21%
70) 1,4-Dichlorobenzene-d4	12.18	152	320583	50.00	ug/l	0.00	82.36%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	430121	49.63	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.26%	
37) 1,2-Dichloroethane-d4	6.39	65	505947	51.00	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.00%	
51) Toluene-d8	8.41	98	1779569	53.14	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.28%	
72) 4-Bromofluorobenzene	11.04	95	369483	55.87	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.74%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.83	96	57	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	2.75	96	56	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.98	43	117	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	438	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	3.93	73	55	N.D.	
22) 1,1-Dichloroethane	4.54	63	2769	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417018.D  
 Acq On : 17 Apr 2007 14:25  
 Sample : JPL32-010  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:24 2007

Vial: 63  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	5.31	77	84	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.48	43	94	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	5.65	128	58	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.81	83	4173	0.28	ug/l	97
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	0.00	56	0	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	6.09	75	63	N.D.		
38) Benzene	6.40	78	711	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.	d	
41) Trichloroethene	7.15	130	1484	N.D.		
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	0.00	41	0	N.D.		
46) Bromodichloromethane	7.73	83	420	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	8.27	75	71	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	8.48	92	575	N.D.		
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	8.87	69	61	N.D.		
55) 1,1,2-Trichloroethane	8.86	97	59	N.D.		
56) Tetrachloroethene	9.03	166	7523	1.03	ug/l	89
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.	d	
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	9.90	112	132	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	9.88	131	61	N.D.		
64) Ethylbenzene	9.99	91	662	N.D.		
65) m,p-Xylene	0.00	106	0	N.D.	d	
66) o-xylene	10.52	106	94	N.D.		
67) Styrene	0.00	104	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417018.D 8260B.M Wed Apr 25 09:25:07 2007

*J. O. 4/25/07*  
 Page 2  
**VOA - 82**

Quantitation Report

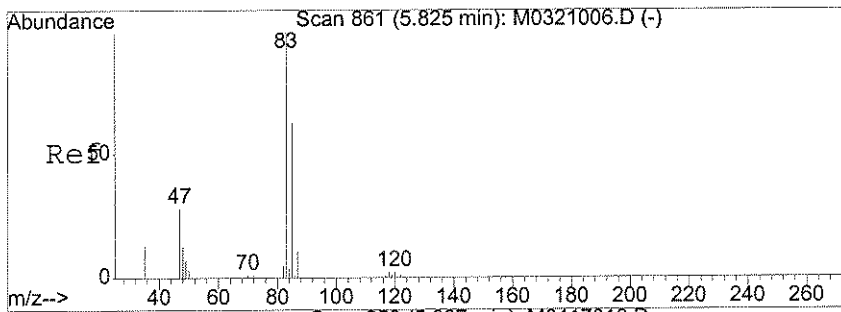
Data File : X:\MSVOA\MOBY\041707\M0417018.D  
 Acq On : 17 Apr 2007 14:25  
 Sample : JPL32-010  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:24 2007

Vial: 63  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

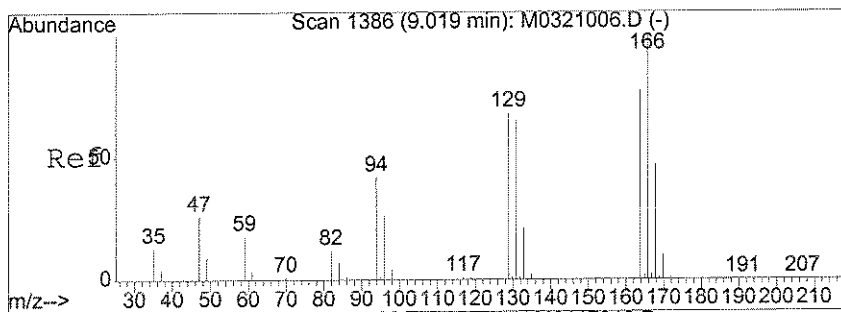
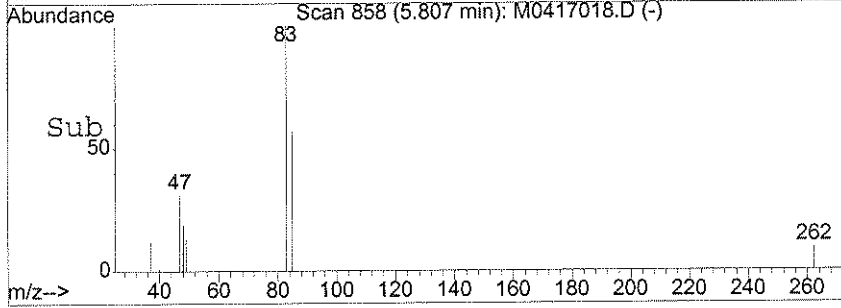
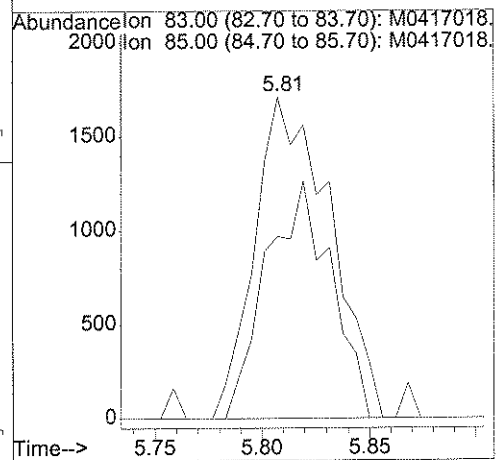
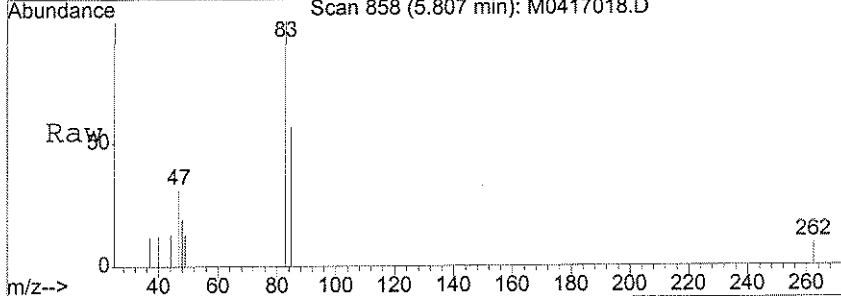
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	292		N.D.	
71) trans-1,4-Dichloro-2-buten	11.03	53	97		N.D.	
73) Bromobenzene	11.03	156	57		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.05	83	93		N.D.	
75) 1,2,3-Trichloropropane	11.20	110	55		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.35	91	66		N.D.	
78) 4-Chlorotoluene	11.48	91	72		N.D.	
79) 1,3,5-Trimethylbenzene	11.60	105	59		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.94	105	68		N.D.	
82) sec-butylbenzene	11.98	105	78		N.D.	
83) 1,3-Dichlorobenzene	12.20	146	81		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	81		N.D.	
86) 1,2-Dichlorobenzene	12.59	146	78		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



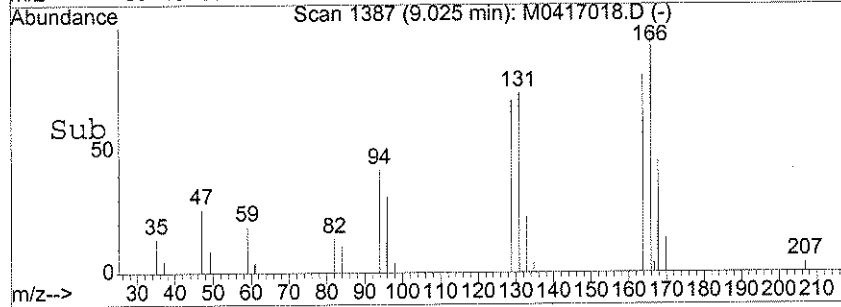
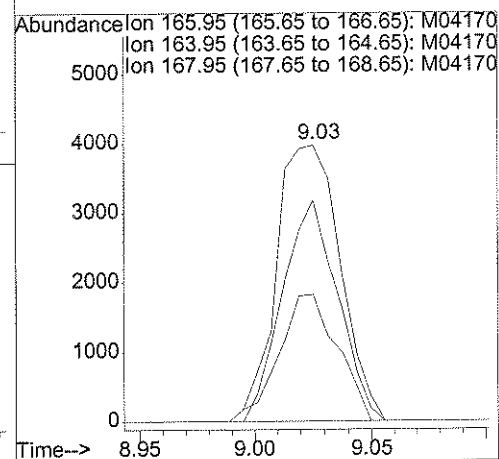
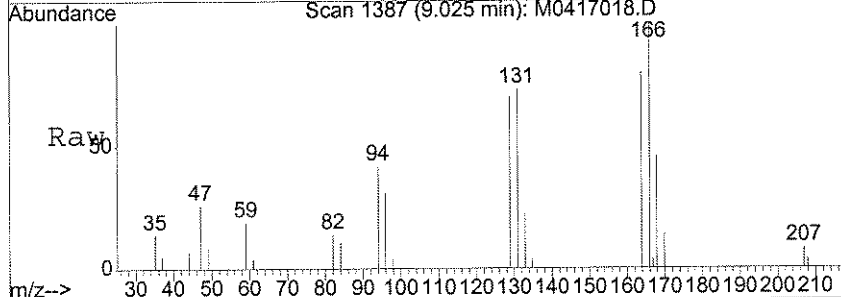
#31  
 Chloroform  
 Concen: 0.28 ug/l  
 RT: 5.81 min Scan# 858  
 Delta R.T. -0.01 min  
 Lab File: M0417018.D  
 Acq: 17 Apr 2007 14:25

Tgt Ion:	83	Resp:	4173
Ion Ratio	Lower	Upper	
83	100		
85	63.5	41.2	81.2



#56  
 Tetrachloroethene  
 Concen: 1.03 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0417018.D  
 Acq: 17 Apr 2007 14:25

Tgt Ion:	166	Resp:	7523
Ion Ratio	Lower	Upper	
166	100		
164	69.4	63.3	94.9
168	42.1	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-6-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-6-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.91	J
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-6-1Q07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 14:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

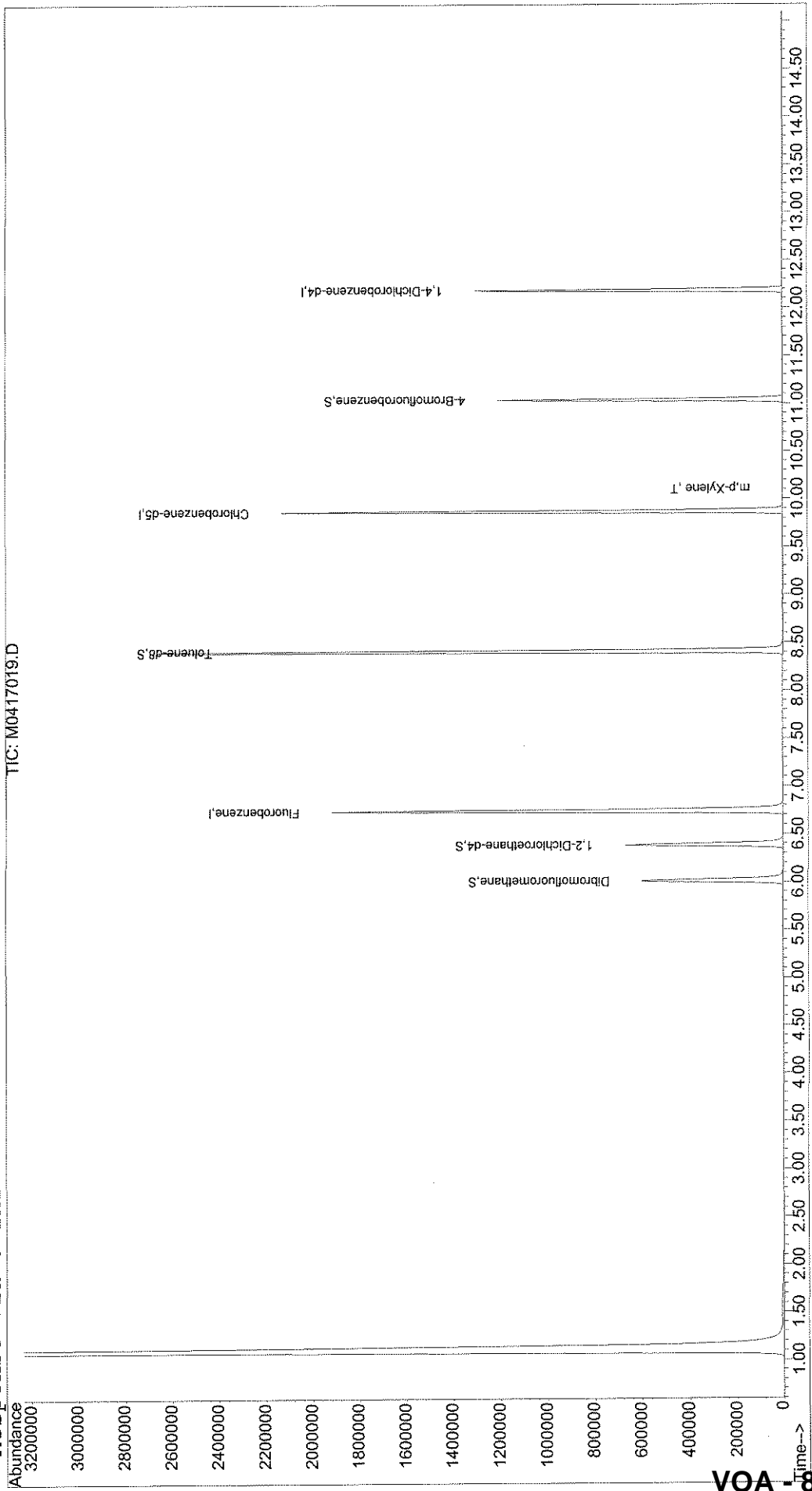
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417019.D Vial: 64  
Acq On : 17 Apr 2007 14:49 Operator: DGA  
Sample : JPL32-011 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:27 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417019.D  
 Acq On : 17 Apr 2007 14:49  
 Sample : JPL32-011  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:27 2007

Vial: 64  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1818907	50.00	ug/l	0.00 93.19%
50) Chlorobenzene-d5	9.87	82	606714	50.00	ug/l	0.00 86.27%
70) 1,4-Dichlorobenzene-d4	12.18	152	319005	50.00	ug/l	0.00 81.96%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	428340	49.92	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.84%
37) 1,2-Dichloroethane-d4	6.39	65	501205	51.02	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.04%
51) Toluene-d8	8.41	98	1768088	52.76	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.52%
72) 4-Bromofluorobenzene	11.04	95	357675	54.36	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	108.72%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.42	50	56	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.83	96	84	N.D.	
6) Chloroethane	1.89	64	76	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.04	76	2090	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.38	43	65	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	3.77	96	60	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	4.54	63	1637	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417019.D  
 Acq On : 17 Apr 2007 14:49  
 Sample : JPL32-011  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:27 2007

Vial: 64  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.33	96	67		N.D.	
27) 2-Butanone	5.52	43	58		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.82	83	2308		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	556		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.15	130	659		N.D.	
42) Methylcyclohexane	7.41	83	62		N.D.	
43) 1,2-Dichloropropane	7.41	63	58		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.48	41	62		N.D.	
46) Bromodichloromethane	7.72	83	422		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.07	75	60		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	809		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.98	97	66		N.D.	
56) Tetrachloroethene	9.02	166	800		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	9.36	129	65		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	68		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.18	131	56		N.D.	
64) Ethylbenzene	9.99	91	517		N.D.	
65) m,p-Xylene	10.10	106	1007	0.91	ug/l	99
66) o-xylene	10.50	106	65		N.D.	
67) Styrene	10.52	104	62		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417019.D 8260B.M Wed Apr 25 09:27:21 2007

*J* 04/25/07  
 Page 2  
**VOA-90**

Quantitation Report

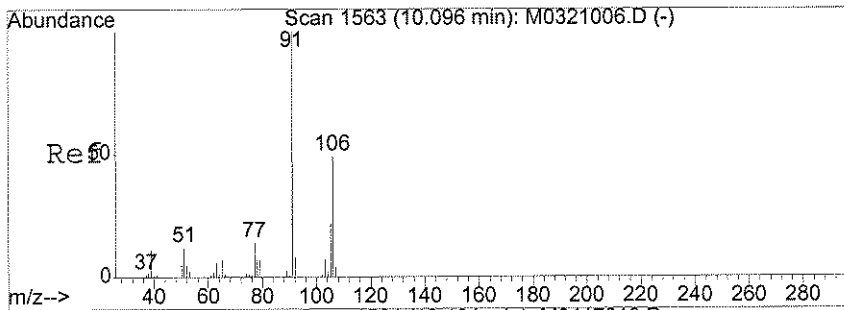
Data File : X:\MSVOA\MOBY\041707\M0417019.D  
 Acq On : 17 Apr 2007 14:49  
 Sample : JPL32-011  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:27 2007

Vial: 64  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

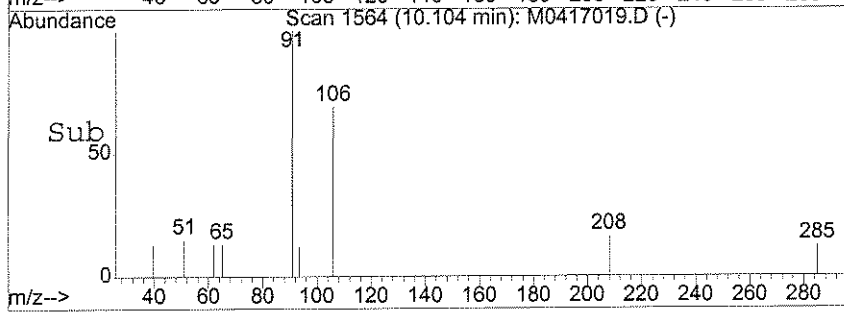
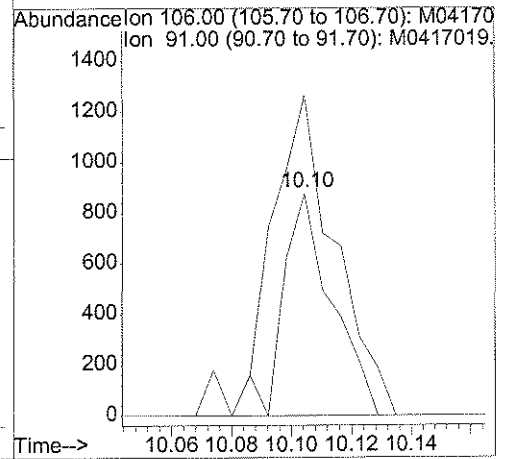
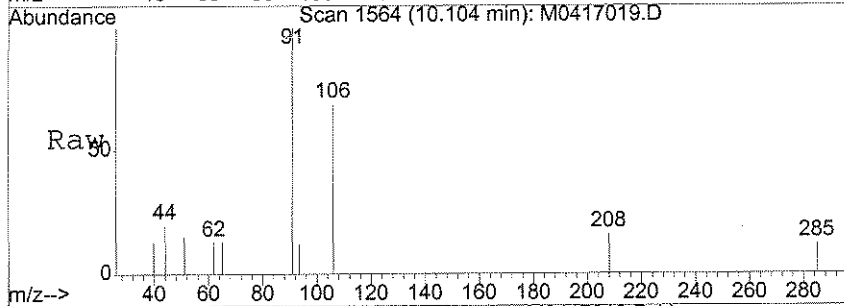
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	301		N.D.	
69) Isopropylbenzene	10.86	105	74		N.D.	
71) trans-1,4-Dichloro-2-buten	11.25	53	75		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.30	83	73		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	59		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.39	91	59		N.D.	
78) 4-Chlorotoluene	11.39	91	59		N.D.	
79) 1,3,5-Trimethylbenzene	11.53	105	72		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.95	105	56		N.D.	
82) sec-butylbenzene	11.95	105	56		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	1234		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	
85) 1,4-Dichlorobenzene	12.12	146	1234		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	633		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.33	225	69		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



#65  
 m,p-Xylene  
 Concen: 0.91 ug/l  
 RT: 10.10 min Scan# 1564  
 Delta R.T. -0.00 min  
 Lab File: M0417019.D  
 Acq: 17 Apr 2007 14:49

Tgt Ion: 106 Resp: 1007  
 Ion Ratio Lower Upper  
 106 100  
 91 188.5 169.6 209.6



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 15:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 15:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.8	
95-47-6	o-Xylene	0.36	J
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-012

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 15:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

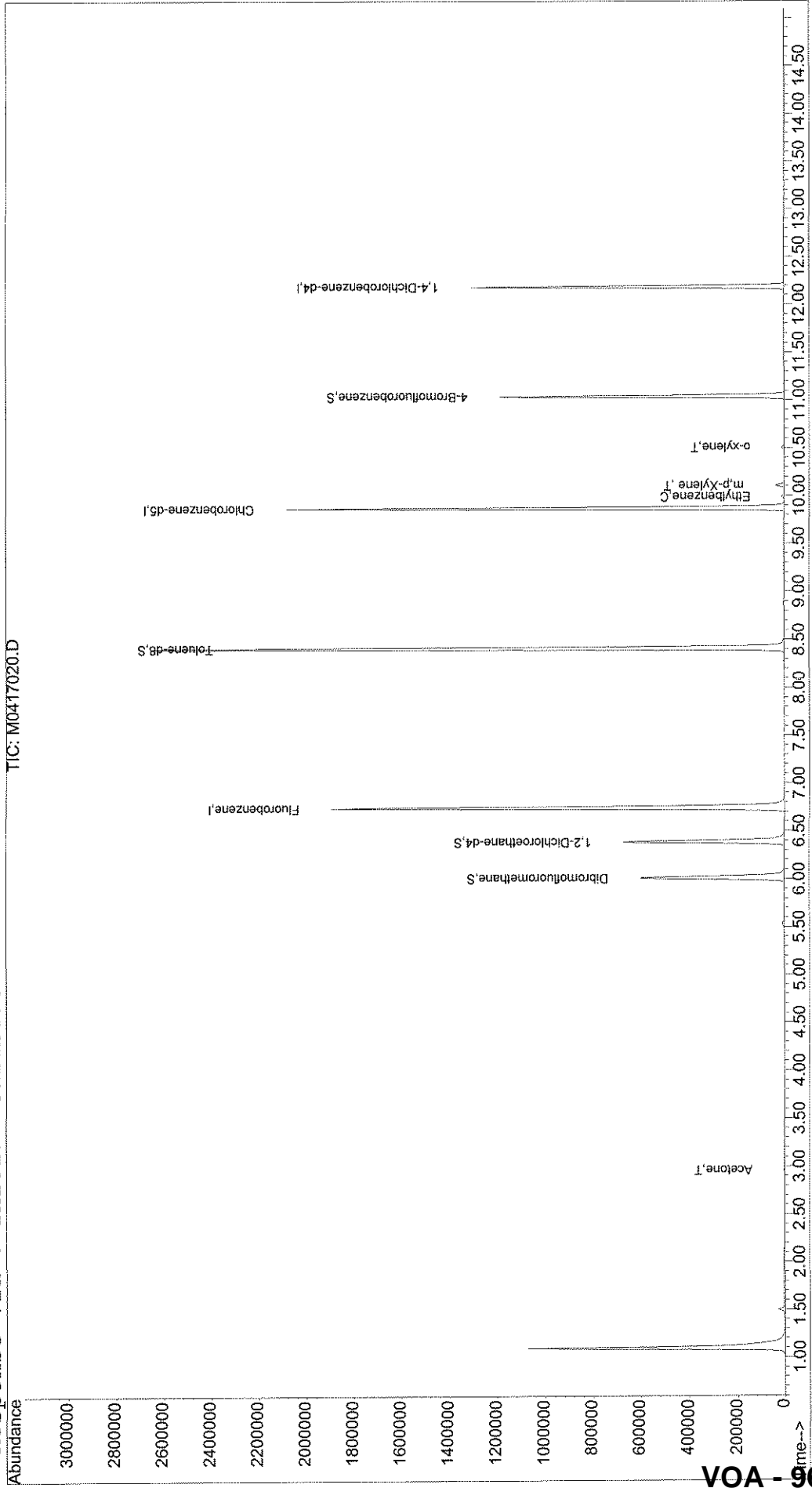
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417020.D  
Acq On : 17 Apr 2007 15:13  
Sample : JPL32-012  
Misc : #1 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:28 2007  
Vial: 65  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417020.D  
 Acq On : 17 Apr 2007 15:13  
 Sample : JPL32-012  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:28 2007

Vial: 65  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	1808933	50.00	ug/l	0.00 92.67%
50) Chlorobenzene-d5	9.87	82	594858	50.00	ug/l	0.00 84.58%
70) 1,4-Dichlorobenzene-d4	12.19	152	322528	50.00	ug/l	0.00 82.86%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	431638	50.58	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.16%
37) 1,2-Dichloroethane-d4	6.39	65	496979	50.87	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.74%
51) Toluene-d8	8.41	98	1748775	53.23	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.46%
72) 4-Bromofluorobenzene	11.04	95	358985	53.96	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	107.92%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.45	50	75	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.83	96	80	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.96	43	5996	3.64 ug/l	#	89
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	198	N.D.		
15) Allyl chloride	3.34	76	56	N.D.		
16) Acetonitrile	3.32	40	62	N.D.		
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417020.D 8260B.M Wed Apr 25 09:28:52 2007

8 04/25/07  
 VOA-97 Page 1

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417020.D  
 Acq On : 17 Apr 2007 15:13  
 Sample : JPL32-012  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:28 2007

Vial: 65  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	d
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.60	128	59		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.82	83	555		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	5.98	56	170		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.15	75	82		N.D.	
38) Benzene	6.41	78	1060		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.60	41	75		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.18	75	68		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	846		N.D.	
53) trans-1,3-Dichloropropene	8.94	75	61		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.13	166	78		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	9.23	129	59		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	232		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.00	91	7405	0.23	ug/l	98
65) m,p-Xylene	10.10	106	10505	1.77	ug/l	88
66) o-xylene	10.50	106	3645	0.36	ug/l	92
67) Styrene	10.51	104	200		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417020.D 8260B.M Wed Apr 25 09:28:52 2007

*[Handwritten Signature]*  
 VOA-98 2

Quantitation Report

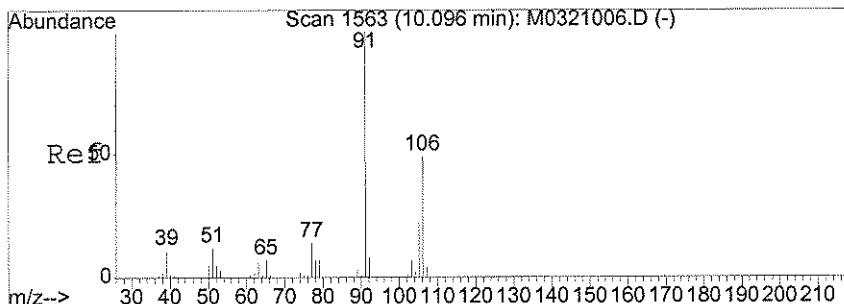
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 Acq On : 17 Apr 2007 15:13  
 Sample : JPL32-012  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:28 2007

Vial: 65  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

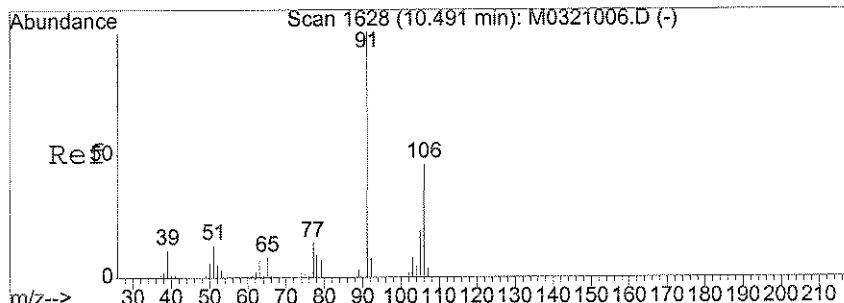
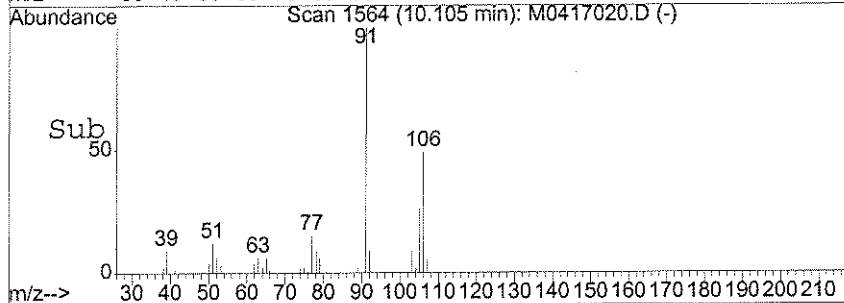
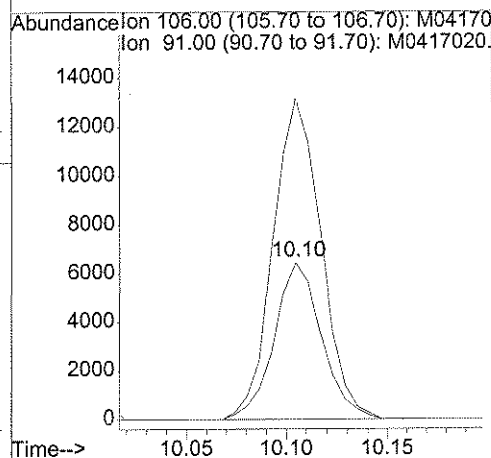
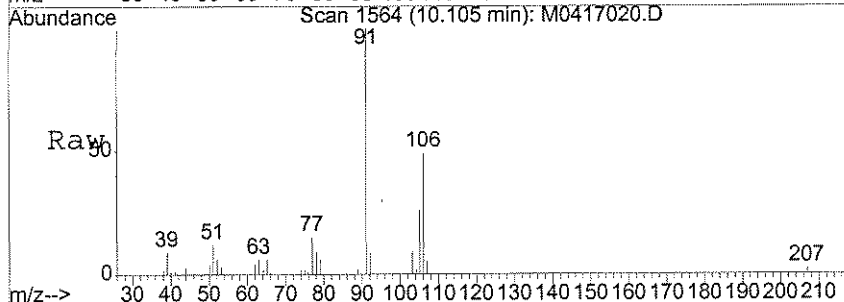
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	90		N.D.	
69) Isopropylbenzene	10.85	105	133		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	323		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	248		N.D.	
76) n-Propylbenzene	11.28	120	63		N.D.	
77) 2-Chlorotoluene	11.36	91	140		N.D.	
78) 4-Chlorotoluene	11.48	91	130		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	207		N.D.	
80) tert-Butylbenzene	11.66	119	59		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	436		N.D.	
82) sec-butylbenzene	11.98	105	57		N.D.	
83) 1,3-Dichlorobenzene	12.20	146	59		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	59		N.D.	
86) 1,2-Dichlorobenzene	12.85	146	62		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.42	225	60		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



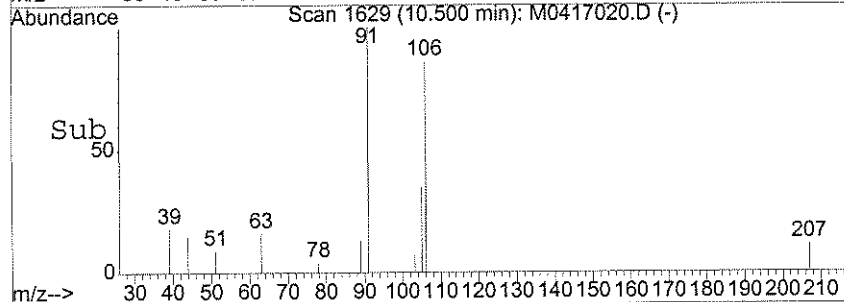
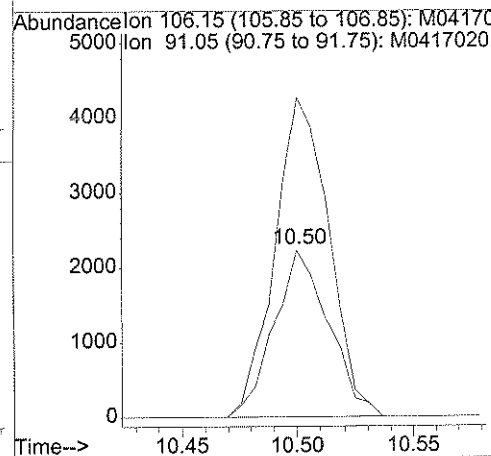
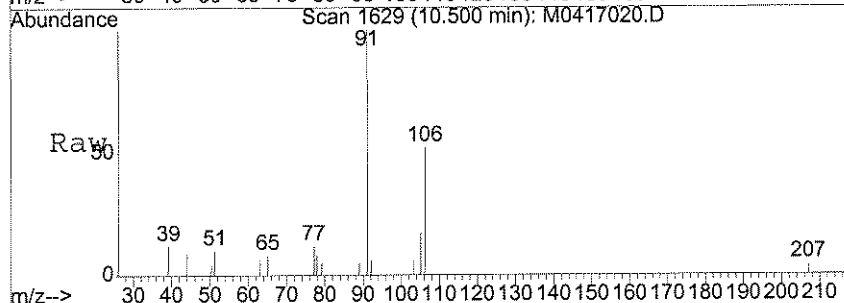
#65  
 m,p-Xylene  
 Concen: 1.77 ug/l  
 RT: 10.10 min Scan# 1564  
 Delta R.T. -0.00 min  
 Lab File: M0417020.D  
 Acq: 17 Apr 2007 15:13

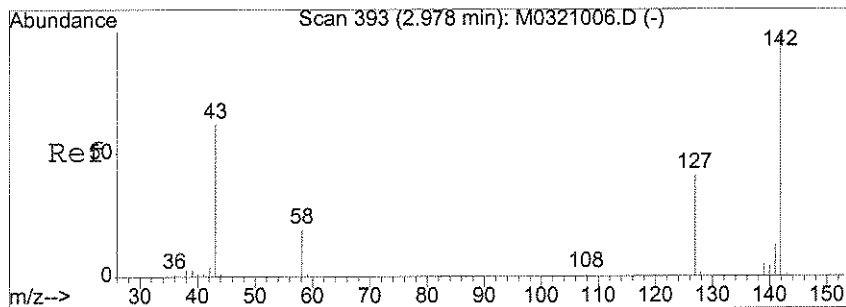
Tgt Ion:106 Resp: 10505  
 Ion Ratio Lower Upper  
 106 100  
 91 207.5 169.6 209.6



#66  
 o-xylene  
 Concen: 0.36 ug/l  
 RT: 10.50 min Scan# 1629  
 Delta R.T. -0.00 min  
 Lab File: M0417020.D  
 Acq: 17 Apr 2007 15:13

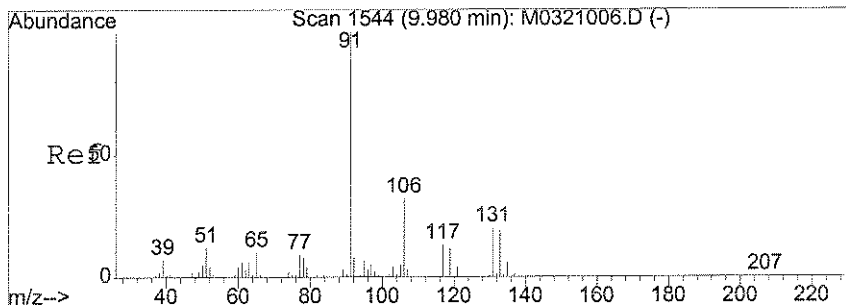
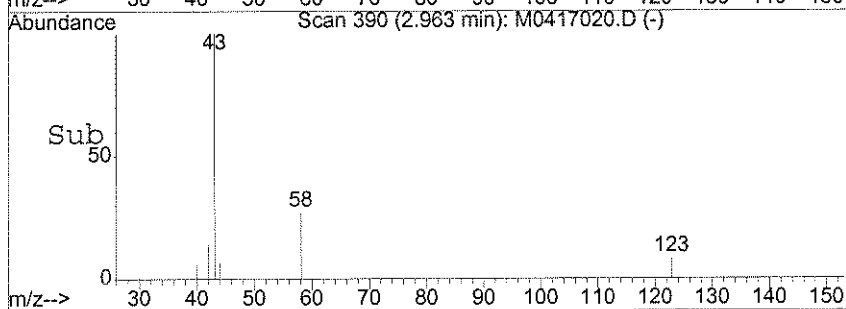
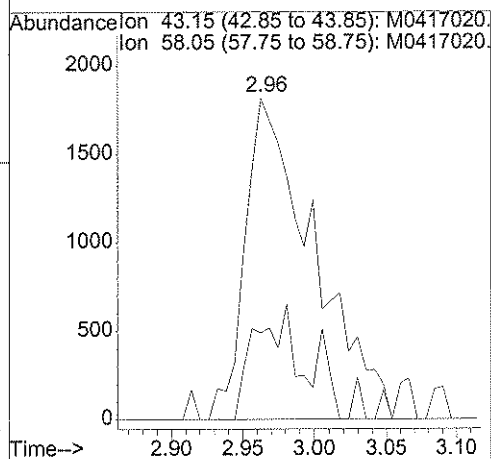
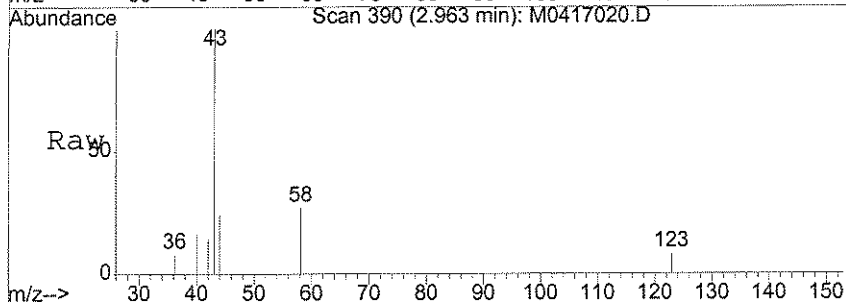
Tgt Ion:106 Resp: 3645  
 Ion Ratio Lower Upper  
 106 100  
 91 188.3 160.6 240.8





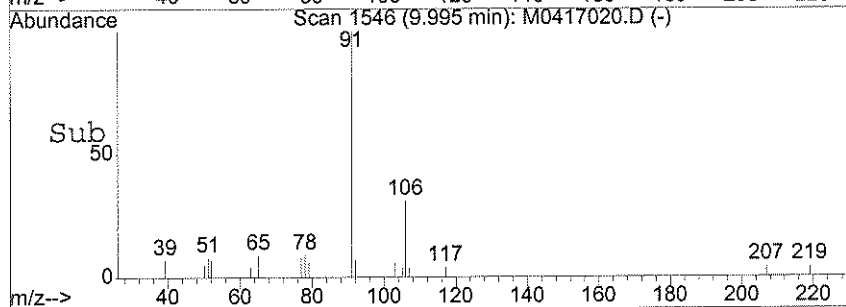
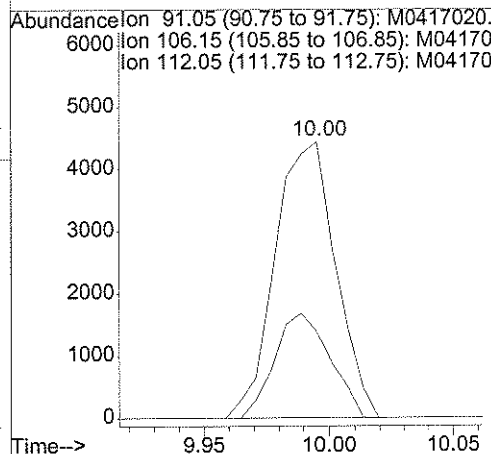
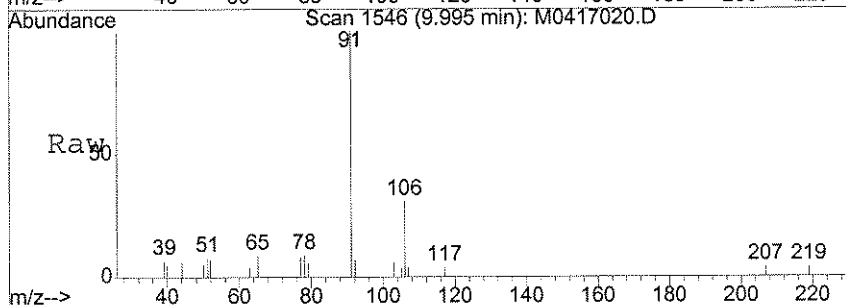
#11  
 Acetone  
 Concen: 3.64 ug/l  
 RT: 2.96 min Scan# 390  
 Delta R.T. 0.04 min  
 Lab File: M0417020.D  
 Acq: 17 Apr 2007 15:13

Tgt Ion: 43 Resp: 5996  
 Ion Ratio Lower Upper  
 43 100  
 58 21.5 21.9 32.9#



#64  
 Ethylbenzene  
 Concen: 0.23 ug/l  
 RT: 10.00 min Scan# 1546  
 Delta R.T. 0.00 min  
 Lab File: M0417020.D  
 Acq: 17 Apr 2007 15:13

Tgt Ion: 91 Resp: 7405  
 Ion Ratio Lower Upper  
 91 100  
 106 31.2 26.1 39.1  
 112 0.0 0.0 0.0



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-013  
 Lab File ID: M0417021.D  
 Date Collected: 04/06/2007  
 Date/Time Analyzed: 04/17/2007 15:37  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 15:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-013

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417021.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/06/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 15:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

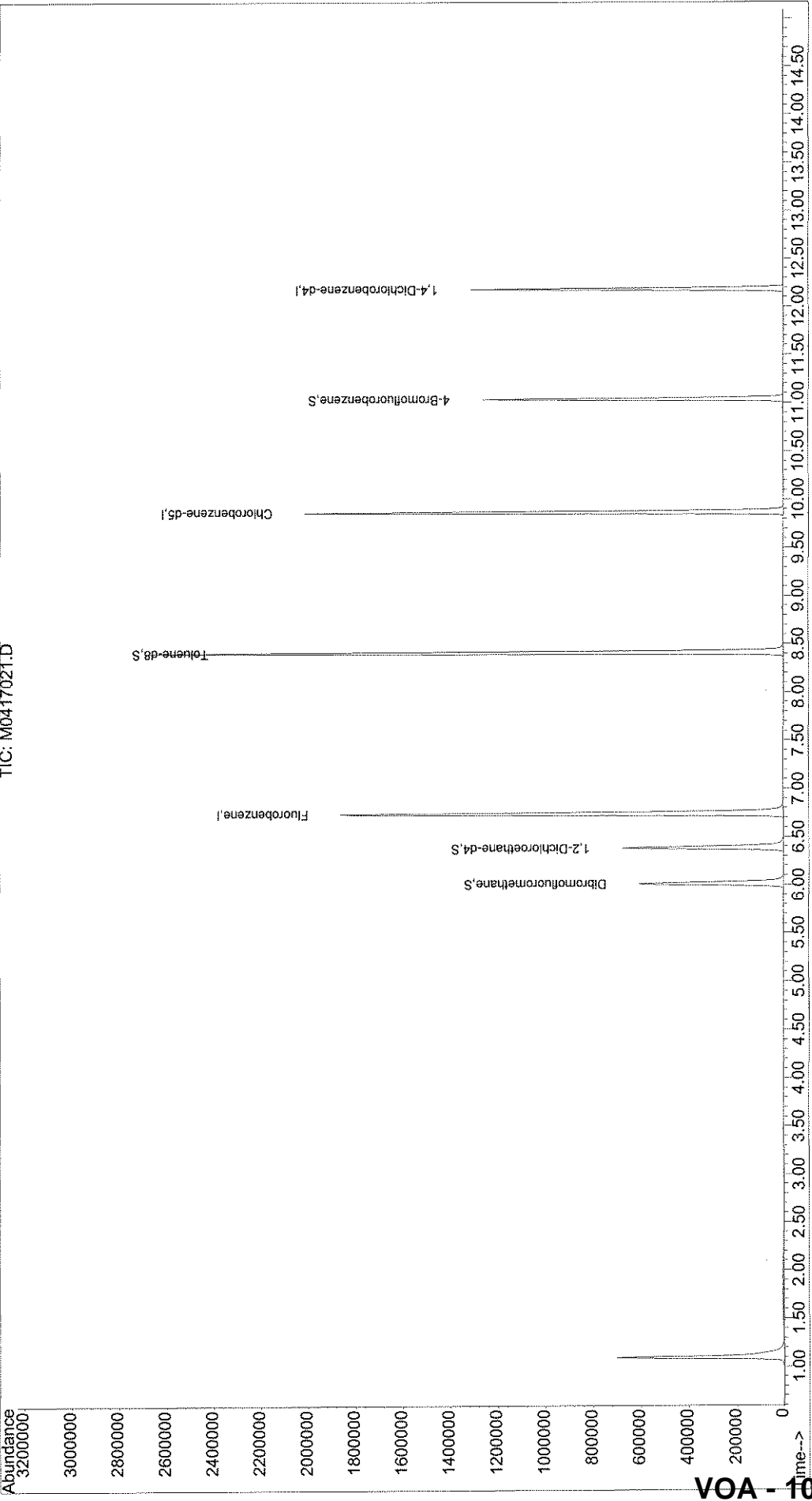
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417021.D  
Acq On : 17 Apr 2007 15:37  
Sample : JPL32-013  
Misc : #1 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:29 2007  
Vial: 66  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417021.D  
 Acq On : 17 Apr 2007 15:37  
 Sample : JPL32-013  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:29 2007

Vial: 66  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1795219	50.00	ug/l	0.00 91.97%
50) Chlorobenzene-d5	9.86	82	602841	50.00	ug/l	0.00 85.72%
70) 1,4-Dichlorobenzene-d4	12.18	152	317669	50.00	ug/l	0.00 81.61%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	421619	49.78	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.56%
37) 1,2-Dichloroethane-d4	6.38	65	495128	51.07	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.14%
51) Toluene-d8	8.41	98	1743034	52.35	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.70%
72) 4-Bromofluorobenzene	11.04	95	364861	55.68	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.36%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.82	96	77	N.D.	
6) Chloroethane	1.93	64	82	N.D.	
7) Trichlorofluoromethane	2.22	101	59	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.99	43	142	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.04	76	74	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	3.32	40	59	N.D.	
17) Methyl Acetate	3.30	43	62	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	3.93	73	58	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417021.D 8260B.M Wed Apr 25 09:29:52 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417021.D  
 Acq On : 17 Apr 2007 15:37  
 Sample : JPL32-013  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:29 2007

Vial: 66  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	5.31	77	68		N.D.	
26) cis-1,2-Dichloroethene	5.41	96	58		N.D.	
27) 2-Butanone	5.54	43	58		N.D.	
28) Propionitrile	5.59	54	61		N.D.	
29) Bromochloromethane	5.65	128	58		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.81	83	61		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.27	117	65		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	563		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.16	130	63		N.D.	
42) Methylcyclohexane	7.32	83	56		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.42	41	58		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.49	92	644		N.D.	
53) trans-1,3-Dichloropropene	8.64	75	64		N.D.	
54) Ethyl methacrylate	8.75	69	77		N.D.	
55) 1,1,2-Trichloroethane	8.87	97	57		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	8.90	76	55		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	229		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	408		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.52	104	94		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417021.D  
 Acq On : 17 Apr 2007 15:37  
 Sample : JPL32-013  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:29 2007

Vial: 66  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.73	173	143		N.D.	
69) Isopropylbenzene	11.04	105	282		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.19	156	71		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.30	83	56		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	263		N.D.	
76) n-Propylbenzene	11.04	120	57		N.D.	
77) 2-Chlorotoluene	11.36	91	72		N.D.	
78) 4-Chlorotoluene	11.48	91	124		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	182		N.D.	
80) tert-Butylbenzene	11.90	119	61		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	61		N.D.	
82) sec-butylbenzene	11.99	105	56		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	111		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	111		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.48	225	58		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-5

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

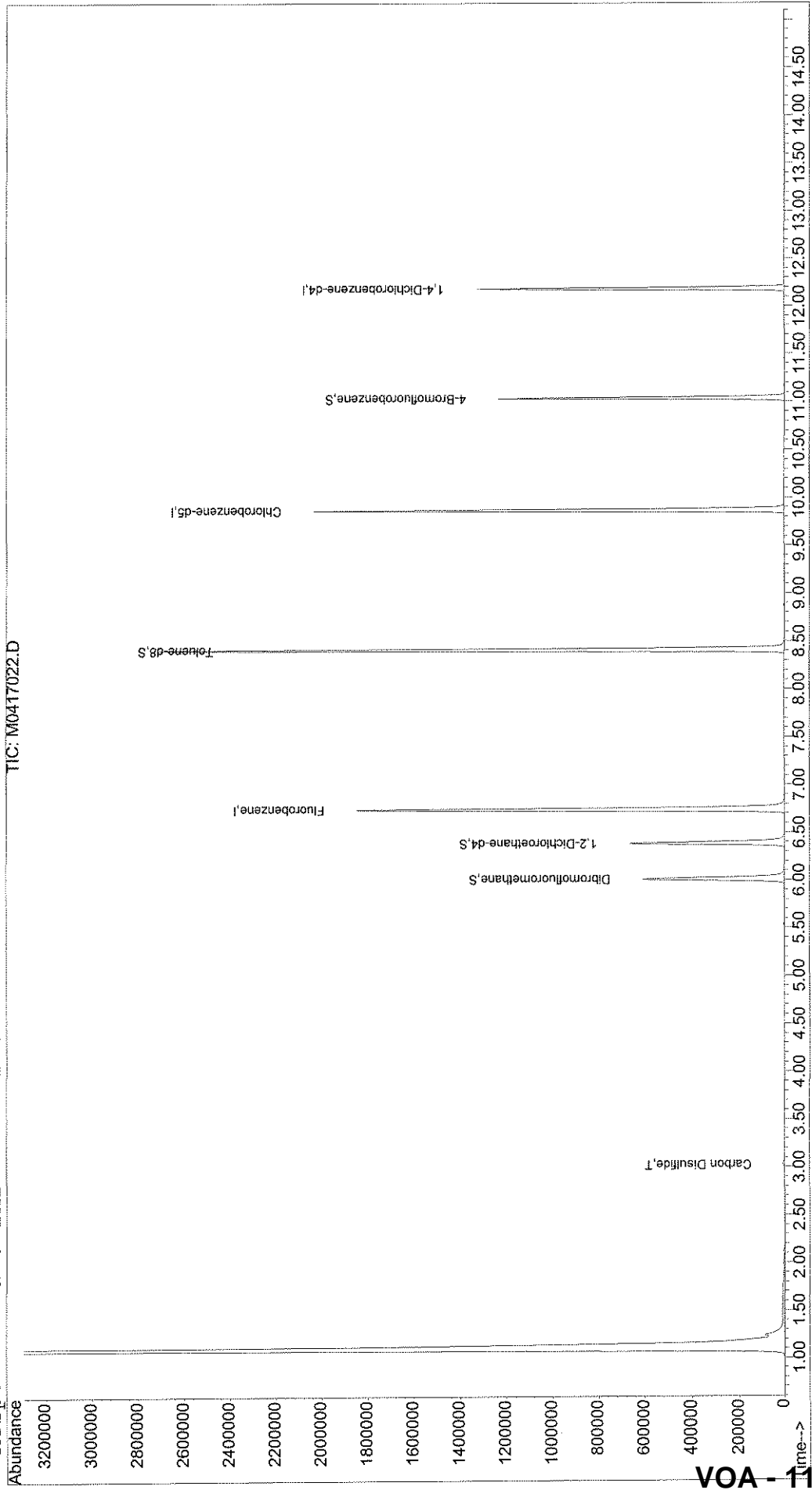
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417022.D Vial: 67  
Acq On : 17 Apr 2007 16:01 Operator: DGA  
Sample : JPL32-014 Inst : MOBY  
Misc : #1 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:31 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA-112

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417022.D  
 Acq On : 17 Apr 2007 16:01  
 Sample : JPL32-014  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:31 2007

Vial: 67  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	1802294	50.00	ug/l	0.00 92.33%
50) Chlorobenzene-d5	9.87	82	609316	50.00	ug/l	0.00 86.64%
70) 1,4-Dichlorobenzene-d4	12.18	152	319441	50.00	ug/l	0.00 82.07%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	428974	50.45	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.90%
37) 1,2-Dichloroethane-d4	6.38	65	497624	51.13	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.26%
51) Toluene-d8	8.41	98	1774778	52.74	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.48%
72) 4-Bromofluorobenzene	11.04	95	360166	54.66	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.32%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.45	50	60	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.86	96	61	N.D.	
6) Chloroethane	1.96	64	363	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	2.96	108	73	N.D.	
14) Carbon Disulfide	3.03	76	11979	0.67 ug/l	100
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417022.D 8260B.M Wed Apr 25 09:31:24 2007

*[Handwritten Signature]*  
 Page 1  
**VOA-113**

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417022.D  
 Acq On : 17 Apr 2007 16:01  
 Sample : JPL32-014  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:31 2007

Vial: 67  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.35	96	98		N.D.	
27) 2-Butanone	5.37	43	75		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	700		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	7.30	83	65		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.72	93	82		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.12	75	74		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	1008		N.D.	
53) trans-1,3-Dichloropropene	8.89	75	60		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.85	97	68		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	9.12	129	63		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.91	112	77		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	9.88	131	58		N.D.	
64) Ethylbenzene	9.99	91	693		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	10.52	106	60		N.D.	
67) Styrene	10.53	104	81		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417022.D 8260B.M Wed Apr 25 09:31:25 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417022.D  
 Acq On : 17 Apr 2007 16:01  
 Sample : JPL32-014  
 Misc : #1 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:31 2007

Vial: 67  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	64		N.D.	
69) Isopropylbenzene	11.02	105	300		N.D.	
71) trans-1,4-Dichloro-2-buten	10.75	53	67		N.D.	
73) Bromobenzene	11.04	156	118		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	404		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	60		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.28	91	70		N.D.	
78) 4-Chlorotoluene	11.28	91	70		N.D.	
79) 1,3,5-Trimethylbenzene	11.32	105	74		N.D.	
80) tert-Butylbenzene	12.01	119	64		N.D.	
81) 1,2,4-Trimethylbenzene	11.98	105	235		N.D.	
82) sec-butylbenzene	11.98	105	235		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	78		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	78		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-015  
 Lab File ID: M0417023.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 16:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) ug/L	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-015  
 Lab File ID: M0417023.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 16:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-4

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-015  
 Lab File ID: M0417023.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 16:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

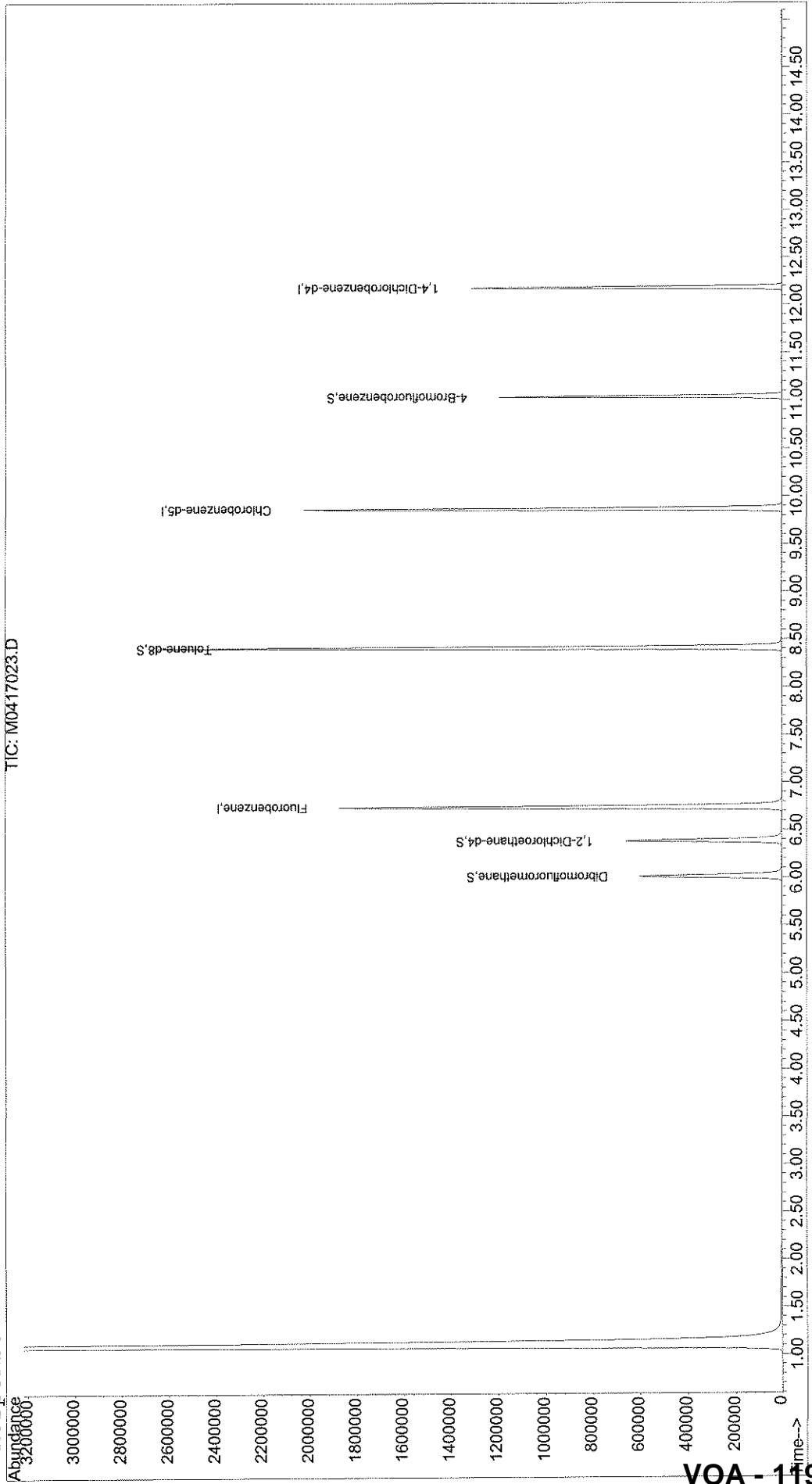
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417023.D Vial: 68  
Acq On : 17 Apr 2007 16:25 Operator: DGA  
Sample : JPL32-015 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:32 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA-119



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417023.D  
 Acq On : 17 Apr 2007 16:25  
 Sample : JPL32-015  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:32 2007

Vial: 68  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)	Rcv (Ar )
1) Fluorobenzene	6.74	96	1774536	50.00	ug/l	0.00	90.91%
50) Chlorobenzene-d5	9.87	82	595127	50.00	ug/l	0.00	84.62%
70) 1,4-Dichlorobenzene-d4	12.18	152	317668	50.00	ug/l	0.00	81.61%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	422068	50.41	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.82%	
37) 1,2-Dichloroethane-d4	6.38	65	486967	50.81	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.62%	
51) Toluene-d8	8.41	98	1733994	52.75	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.50%	
72) 4-Bromofluorobenzene	11.04	95	349921	53.40	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	106.80%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.41	50	194	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.79	96	165	N.D.		
6) Chloroethane	1.97	64	72	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.84	96	55	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.97	43	226	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.03	76	750	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.87	96	120	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417023.D 8260B.M Wed Apr 25 09:32:38 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417023.D  
 Acq On : 17 Apr 2007 16:25  
 Sample : JPL32-015  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:32 2007

Vial: 68  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.69	43	57		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.28	96	56		N.D.	
27) 2-Butanone	5.44	43	57		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.81	83	545		N.D.	
32) 1,1,1-Trichloroethane	6.06	97	126		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.14	117	58		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	823		N.D.	
39) 1,2-Dichloroethane	6.40	62	315		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.66	41	68		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.32	75	58		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	580		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	8.96	166	67		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.91	112	55		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.14	131	78		N.D.	
64) Ethylbenzene	9.98	91	191		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	10.50	106	56		N.D.	
67) Styrene	0.00	104	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417023.D 8260B.M Wed Apr 25 09:32:39 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417023.D  
 Acq On : 17 Apr 2007 16:25  
 Sample : JPL32-015  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:32 2007

Vial: 68  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	115		N.D.	
69) Isopropylbenzene	11.04	105	393		N.D.	
71) trans-1,4-Dichloro-2-buten	10.81	53	55		N.D.	
73) Bromobenzene	11.03	156	140		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	157		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	222		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.37	91	80		N.D.	
78) 4-Chlorotoluene	11.49	91	201		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.97	105	63		N.D.	
82) sec-butylbenzene	11.97	105	63		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	68		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	153		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.25	157	75		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.33	225	70		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-016

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.95	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-3

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-016  
 Lab File ID: M0417024.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 16:49  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.87	J
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-3

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-016

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417024.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 16:49

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

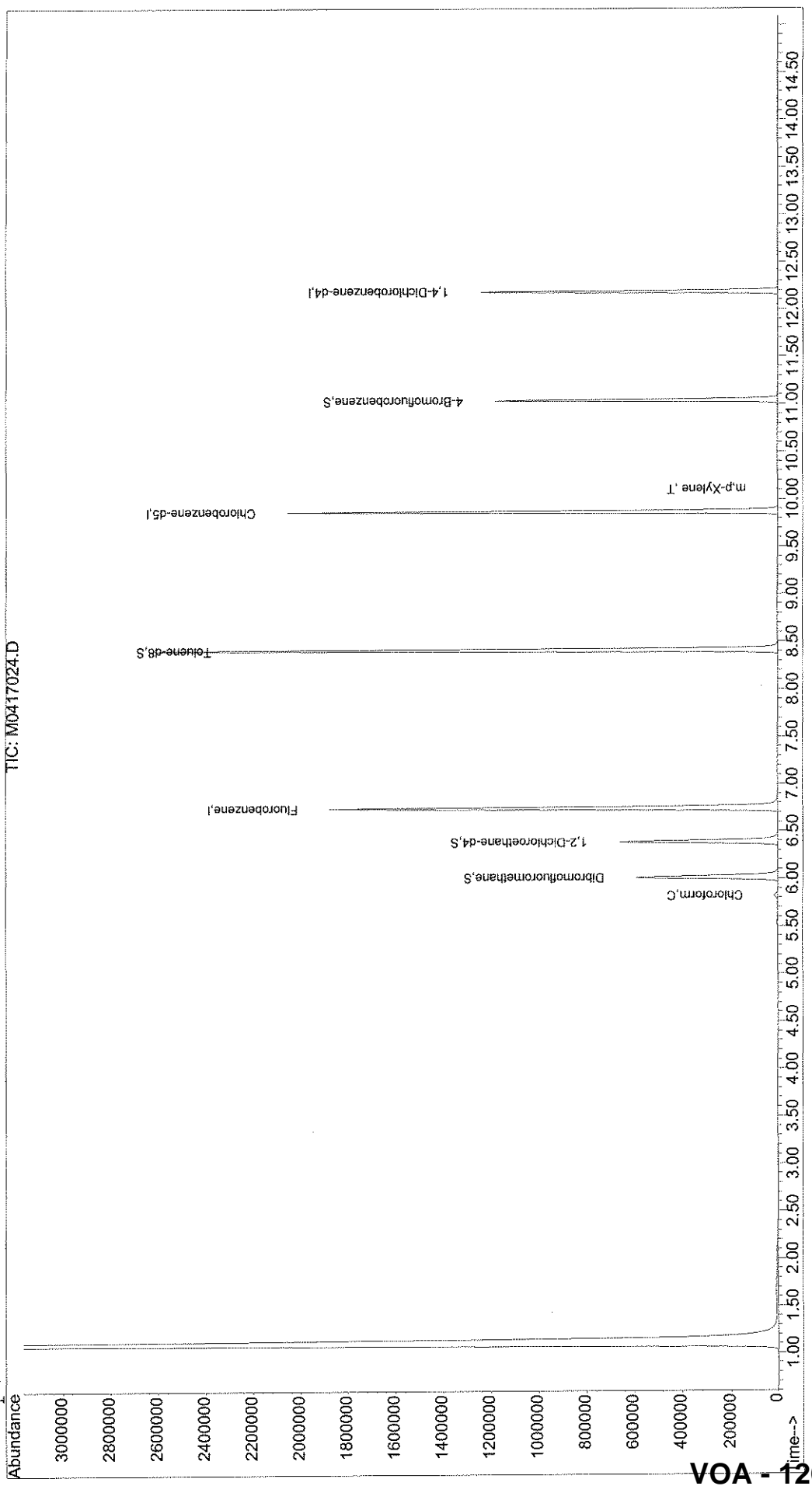
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417024.D  
Acq On : 17 Apr 2007 16:49  
Sample : JPL32-016  
Misc : #2 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:33 2007  
Vial: 69  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417024.D  
 Acq On : 17 Apr 2007 16:49  
 Sample : JPL32-016  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:33 2007

Vial: 69  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1770473	50.00	ug/l	0.00 90.70%
50) Chlorobenzene-d5	9.87	82	586454	50.00	ug/l	0.00 83.39%
70) 1,4-Dichlorobenzene-d4	12.19	152	318684	50.00	ug/l	0.00 81.88%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	419008	50.16	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.32%
37) 1,2-Dichloroethane-d4	6.39	65	494577	51.73	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	103.46%
51) Toluene-d8	8.41	98	1724326	53.24	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.48%
72) 4-Bromofluorobenzene	11.04	95	351084	53.41	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	106.82%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.42	50	61	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	1.95	64	61	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	2.73	96	57	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.98	43	199	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.02	76	406	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.28	43	60	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417024.D 8260B.M Wed Apr 25 09:33:49 2007



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417024.D  
 Acq On : 17 Apr 2007 16:49  
 Sample : JPL32-016  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:33 2007

Vial: 69  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	4.67	53	55		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.50	96	69		N.D.	
27) 2-Butanone	5.56	43	141		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.76	41	62		N.D.	
31) Chloroform	5.82	83	13522	0.95	ug/l	96
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	6.00	56	71		N.D.	
35) Carbon Tetrachloride	6.13	117	56		N.D.	
36) 1,1-Dichloropropene	6.16	75	55		N.D.	
38) Benzene	6.41	78	767		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	510		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.52	93	68		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.71	83	389		N.D.	
47) 2-Chloroethyl vinyl ether	8.12	63	56		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.49	92	322		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.03	166	157		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	9.32	129	61		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	61		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	518		N.D.	
65) m,p-Xylene	10.11	106	589	0.87	ug/l	93
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.66	104	59		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417024.D 8260B.M Wed Apr 25 09:33:50 2007

*[Handwritten Signature]*  
 VOA-128 Page 2

Quantitation Report

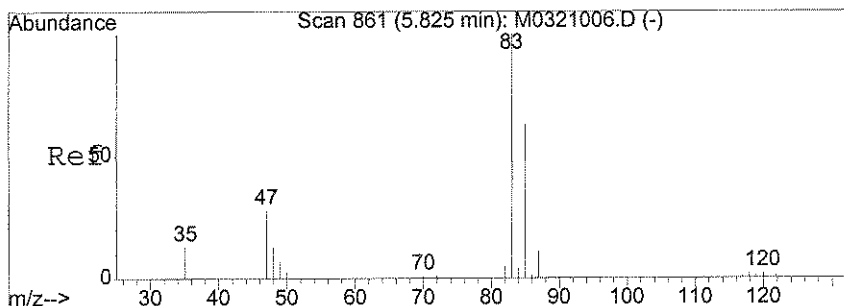
Data File : X:\MSVOA\MOBY\041707\M0417024.D  
 Acq On : 17 Apr 2007 16:49  
 Sample : JPL32-016  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:33 2007

Vial: 69  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

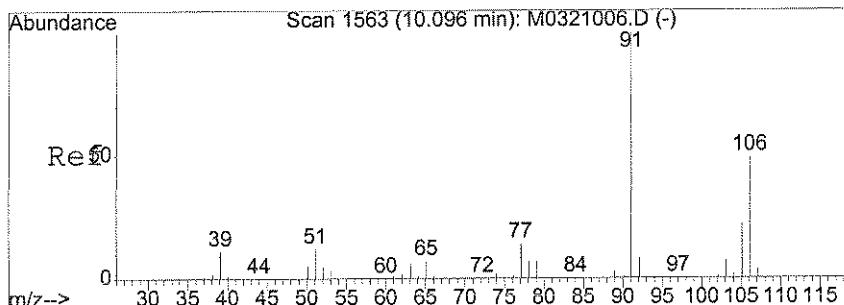
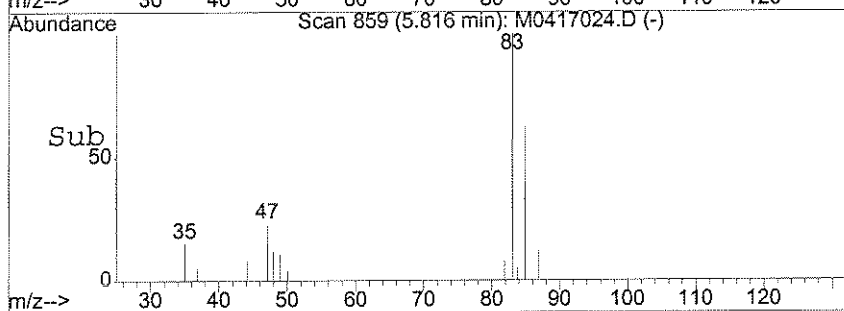
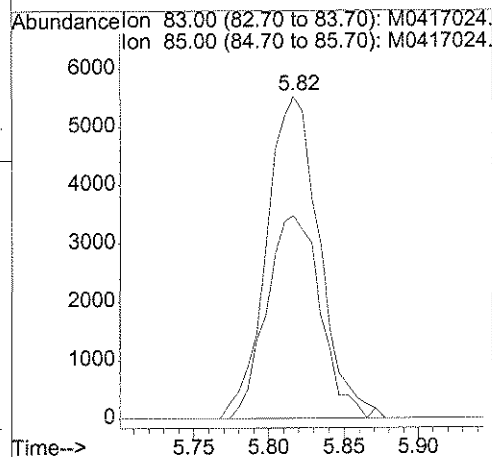
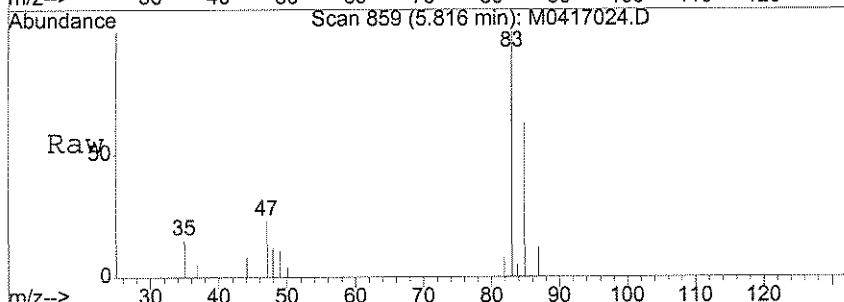
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.73	173	66		N.D.	
69) Isopropylbenzene	11.03	105	735		N.D.	
71) trans-1,4-Dichloro-2-buten	10.98	53	64		N.D.	
73) Bromobenzene	11.05	156	56		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.38	83	77		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	224		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.26	91	139		N.D.	
78) 4-Chlorotoluene	11.67	91	65		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.75	119	60		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	78		N.D.	
82) sec-butylbenzene	11.99	105	57		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	63		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



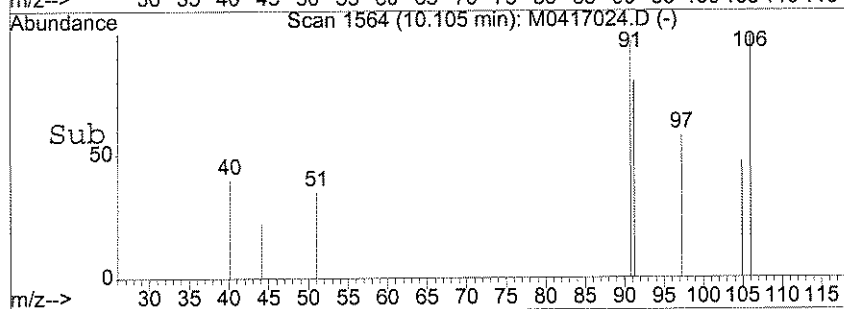
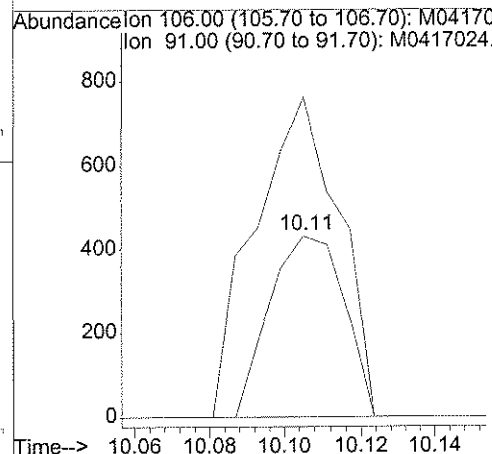
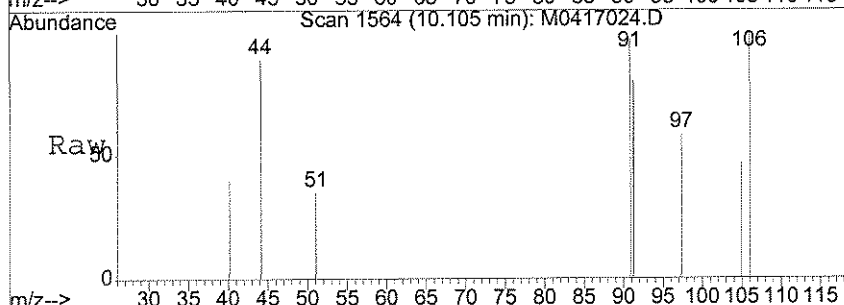
#31  
 Chloroform  
 Concen: 0.95 ug/l  
 RT: 5.82 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0417024.D  
 Acq: 17 Apr 2007 16:49

Tgt Ion: 83 Resp: 13522  
 Ion Ratio Lower Upper  
 83 100  
 85 64.4 41.2 81.2



#65  
 m,p-Xylene  
 Concen: 0.87 ug/l  
 RT: 10.11 min Scan# 1564  
 Delta R.T. -0.00 min  
 Lab File: M0417024.D  
 Acq: 17 Apr 2007 16:49

Tgt Ion: 106 Resp: 589  
 Ion Ratio Lower Upper  
 106 100  
 91 199.8 169.6 209.6



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:13

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

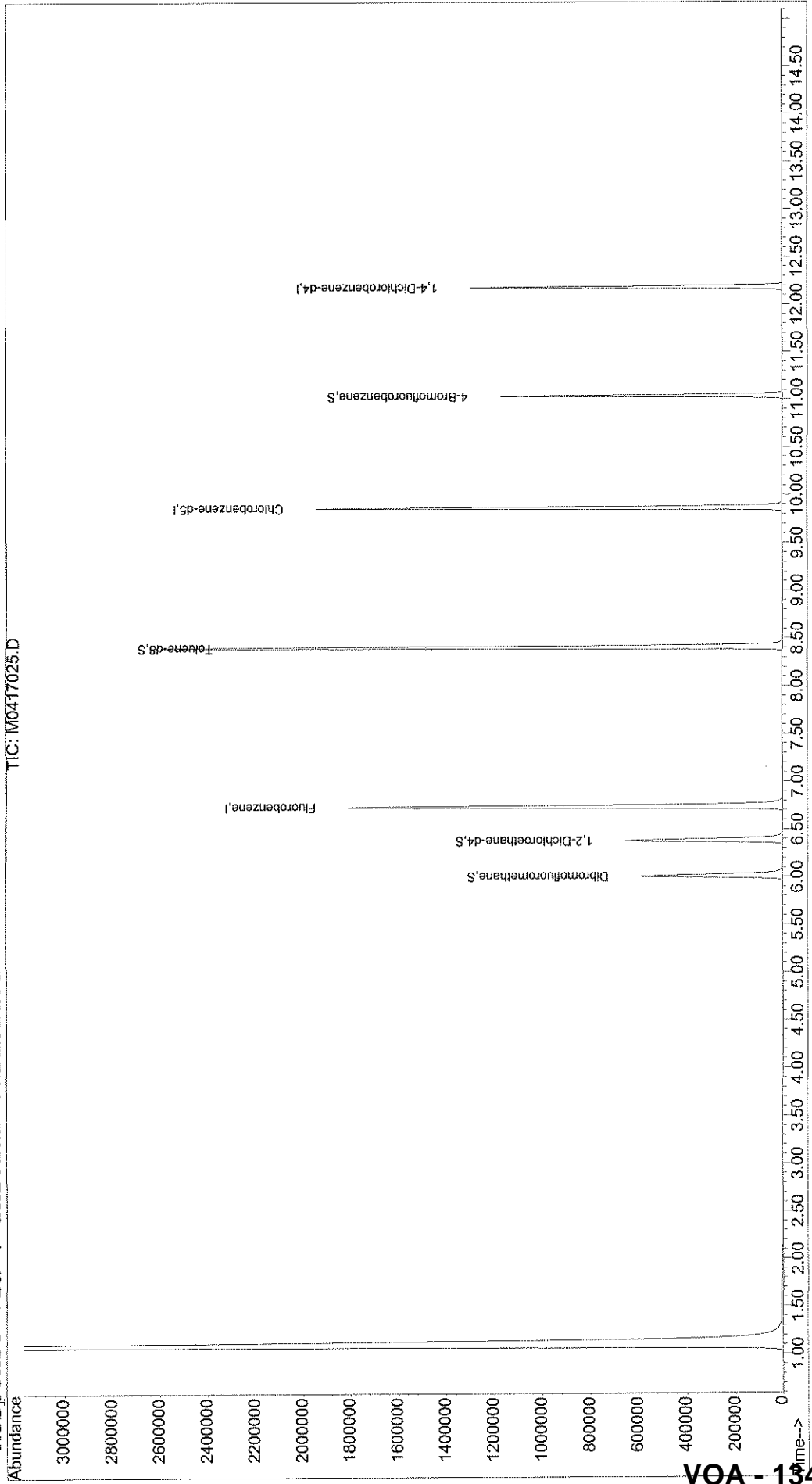
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417025.D  
Acq On : 17 Apr 2007 17:13  
Sample : JPL32-017  
Misc : #2 5ml +IS/SS(524)  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:34 2007  
Vial: 70  
Operator: DGA  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA - 134

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417025.D  
 Acq On : 17 Apr 2007 17:13  
 Sample : JPL32-017  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:34 2007

Vial: 70  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973M

Last Update : Tue Apr 24 12:05:10 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)	Rcv (Ar )
1) Fluorobenzene	6.74	96	1761186	50.00	ug/l	0.00	90.23%
50) Chlorobenzene-d5	9.87	82	582899	50.00	ug/l	0.00	82.88%
70) 1,4-Dichlorobenzene-d4	12.18	152	312692	50.00	ug/l	0.00	80.34%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	418684	50.39	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.78%	
37) 1,2-Dichloroethane-d4	6.38	65	487153	51.22	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	102.44%	
51) Toluene-d8	8.41	98	1717168	53.34	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.68%	
72) 4-Bromofluorobenzene	11.04	95	349327	54.16	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	108.32%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	1.53	62	57	N.D.		
5) Bromomethane	1.86	96	58	N.D.		
6) Chloroethane	1.93	64	82	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.96	43	58	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	905	N.D.		
15) Allyl chloride	3.20	76	55	N.D.		
16) Acetonitrile	3.32	40	58	N.D.		
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	3.93	53	59	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration

M0417025.D 8260B.M Wed Apr 25 09:35:06 2007



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417025.D  
 Acq On : 17 Apr 2007 17:13  
 Sample : JPL32-017  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:34 2007

Vial: 70  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.57	43	62		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	5.30	77	66		N.D.	
26) cis-1,2-Dichloroethene	5.26	96	58		N.D.	
27) 2-Butanone	5.52	43	55		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.57	128	56		N.D.	
30) Methacrylonitrile	5.76	41	56		N.D.	
31) Chloroform	5.82	83	2004		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	705		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	1293		N.D.	
42) Methylcyclohexane	7.19	83	55		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.72	83	58		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	185		N.D.	
53) trans-1,3-Dichloropropene	8.88	75	55		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.88	97	57		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	9.18	129	55		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	202		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	134		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417025.D 8260B.M Wed Apr 25 09:35:07 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417025.D  
 Acq On : 17 Apr 2007 17:13  
 Sample : JPL32-017  
 Misc : #2 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:34 2007

Vial: 70  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	0.00	173	0		N.D.	
69) Isopropylbenzene	11.03	105	368		N.D.	
71) trans-1,4-Dichloro-2-buten	11.04	53	61		N.D.	
73) Bromobenzene	11.03	156	119		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	244		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	57		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.27	91	69		N.D.	
78) 4-Chlorotoluene	11.27	91	69		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	60		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.59	105	69		N.D.	
82) sec-butylbenzene	12.10	105	61		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	61		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	<u>ug/L</u>	
75-71-8	Dichlorodifluoromethane	0.50		U
74-87-3	Chloromethane	0.50		U
75-01-4	Vinyl chloride	0.50		U
74-83-9	Bromomethane	0.50		U
75-00-3	Chloroethane	0.50		U
75-69-4	Trichlorofluoromethane	0.50		U
75-35-4	1,1-Dichloroethene	0.50		U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50		U
75-09-2	Methylene chloride	0.50		U
1634-04-4	Methyl tert-butyl ether	0.50		U
156-60-5	trans-1,2-Dichloroethene	0.50		U
75-34-3	1,1-Dichloroethane	0.50		U
594-20-7	2,2-Dichloropropane	0.50		U
156-59-2	cis-1,2-Dichloroethene	0.50		U
78-93-3	2-Butanone	5.0		U
74-97-5	Bromochloromethane	0.50		U
67-66-3	Chloroform	0.50		U
71-55-6	1,1,1-Trichloroethane	0.50		U
56-23-5	Carbon tetrachloride	0.50		U
563-58-6	1,1-Dichloropropene	0.50		U
71-43-2	Benzene	0.50		U
107-06-2	1,2-Dichloroethane	0.50		U
79-01-6	Trichloroethene	0.50		U
78-87-5	1,2-Dichloropropane	0.50		U
74-95-3	Dibromomethane	0.50		U
75-27-4	Bromodichloromethane	0.50		U
10061-01-	cis-1,3-Dichloropropene	0.50		U
108-10-1	4-Methyl-2-pentanone	5.0		U
108-88-3	Toluene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.85	J
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-25-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-018

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417026.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 17:37

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

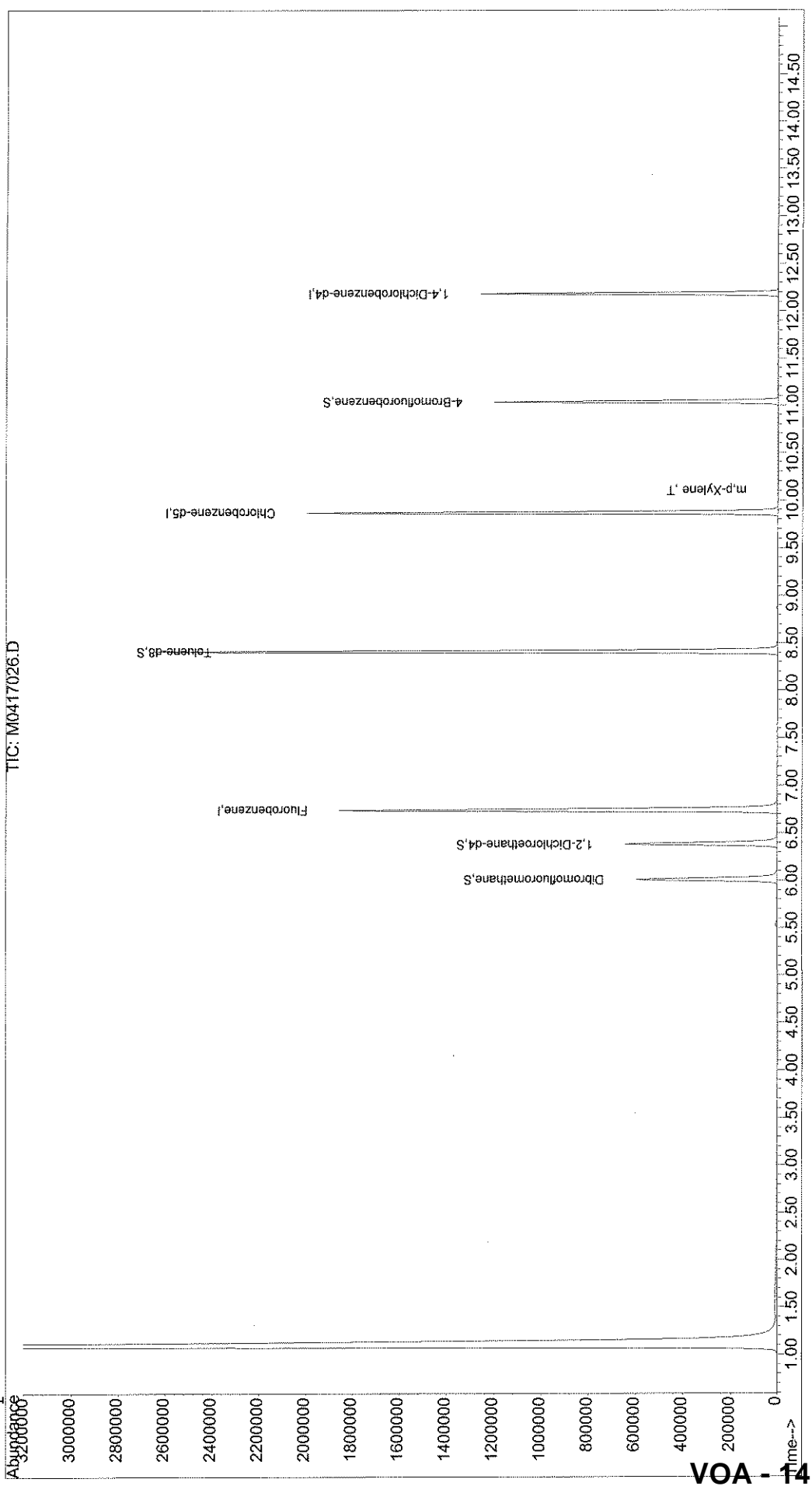
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417026.D Vial: 71  
Acq On : 17 Apr 2007 17:37 Operator: DGA  
Sample : JPL32-018 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:36 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417026.D  
 Acq On : 17 Apr 2007 17:37  
 Sample : JPL32-018  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:36 2007

Vial: 71  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	1758115	50.00	ug/l	0.00	90.07%
50) Chlorobenzene-d5	9.87	82	591839	50.00	ug/l	0.00	84.15%
70) 1,4-Dichlorobenzene-d4	12.18	152	307452	50.00	ug/l	0.00	78.99%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	415325	50.07	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.14%	
37) 1,2-Dichloroethane-d4	6.38	65	482017	50.77	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	101.54%	
51) Toluene-d8	8.41	98	1717181	52.53	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.06%	
72) 4-Bromofluorobenzene	11.04	95	349109	55.05	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.10%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.81	96	55	N.D.	
6) Chloroethane	0.00	64	0	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.04	76	195	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	3.91	73	66	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration

M0417026.D 8260B.M Wed Apr 25 09:36:25 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417026.D  
 Acq On : 17 Apr 2007 17:37  
 Sample : JPL32-018  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:36 2007

Vial: 71  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.52	43	131		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.59	41	62		N.D.	
31) Chloroform	5.82	83	1891		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.26	75	57		N.D.	
38) Benzene	6.41	78	687		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.62	93	61		N.D.	
45) Methyl methacrylate	7.55	41	58		N.D.	
46) Bromodichloromethane	7.73	83	68		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	299		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	59		N.D.	
57) 1,3-Dichloropropane	9.25	76	56		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	87		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	292		N.D.	
65) m,p-Xylene	10.10	106	354	0.85	ug/l	94
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	



Quantitation Report

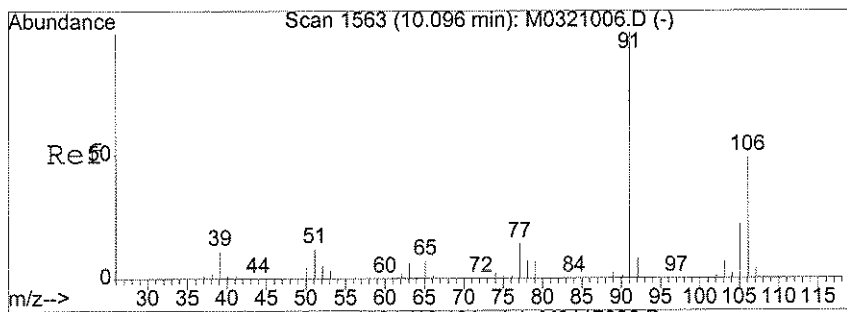
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 Acq On : 17 Apr 2007 17:37  
 Sample : JPL32-018  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:36 2007

Vial: 71  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

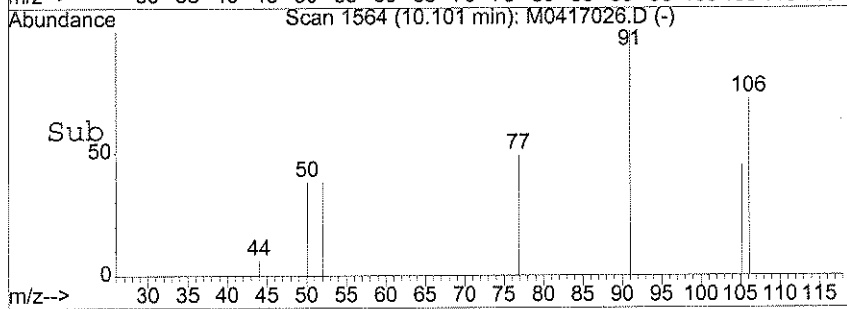
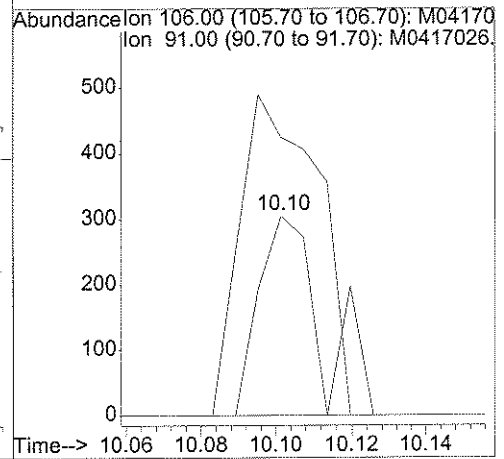
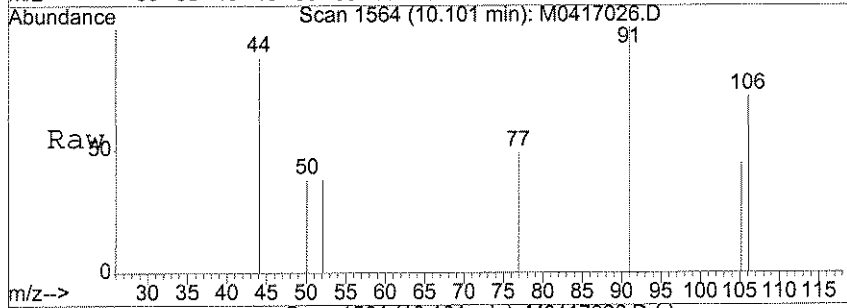
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	123		N.D.	
69) Isopropylbenzene	11.03	105	608		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	380		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	136		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.29	91	55		N.D.	
78) 4-Chlorotoluene	11.49	91	82		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	112		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.98	105	173		N.D.	
82) sec-butylbenzene	11.98	105	173		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.38	157	59		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.32	225	71		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



#65  
 m,p-Xylene  
 Concen: 0.85 ug/l  
 RT: 10.10 min Scan# 1564  
 Delta R.T. -0.01 min  
 Lab File: M0417026.D  
 Acq: 17 Apr 2007 17:37

Tgt Ion: 106	Resp: 354
Ion Ratio Lower	Upper
106 100	
91 198.9	169.6 209.6



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-019  
 Lab File ID: M0417027.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 18:01  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-019

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417027.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	0.94	J
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-019

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417027.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:01

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

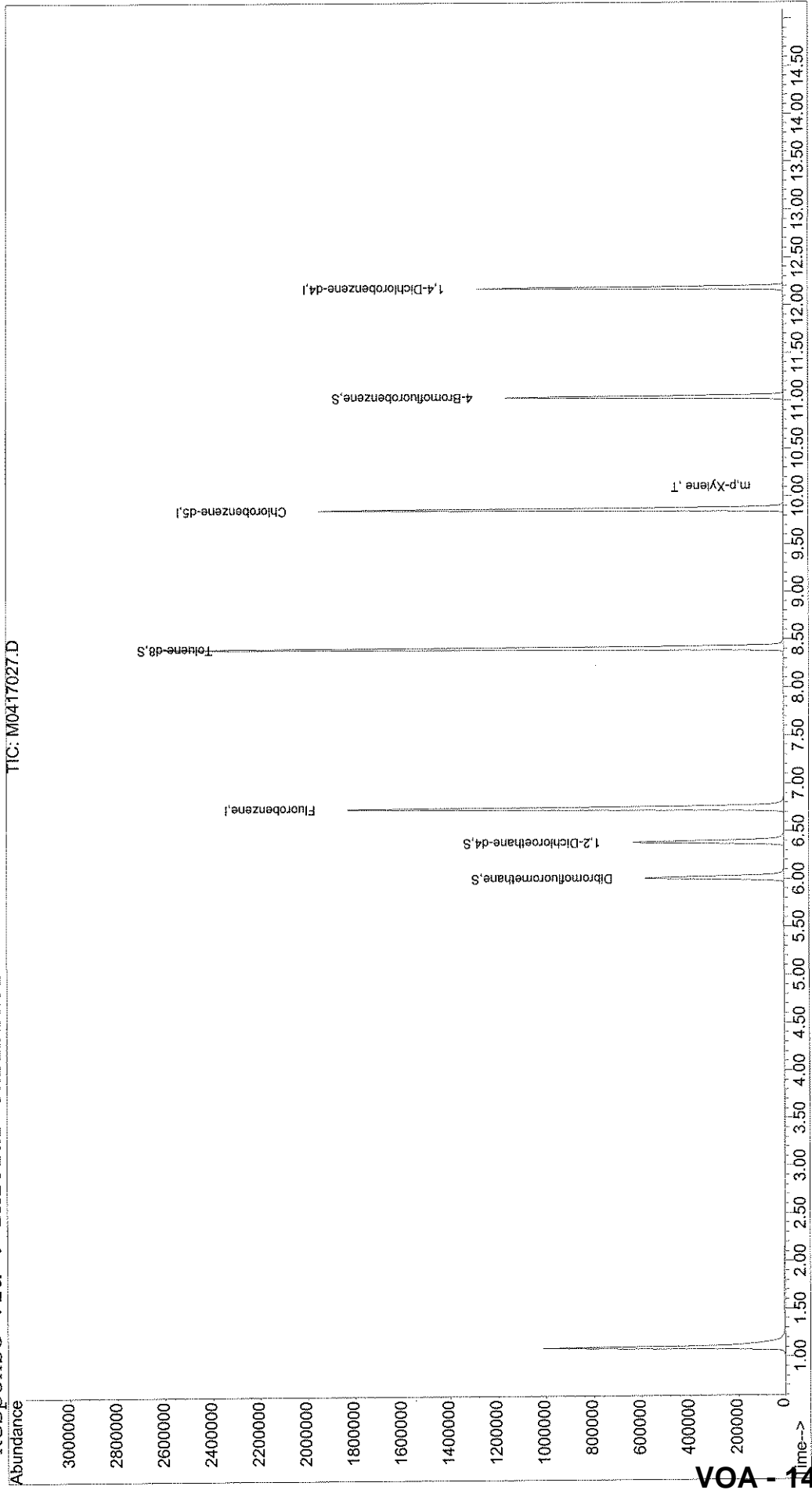
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417027.D Vial: 72  
Acq On : 17 Apr 2007 18:01 Operator: DGA  
Sample : JPL32-019 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:37 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417027.D  
 Acq On : 17 Apr 2007 18:01  
 Sample : JPL32-019  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:37 2007

Vial: 72  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	1751884	50.00	ug/l	0.00	89.75%
50) Chlorobenzene-d5	9.87	82	579908	50.00	ug/l	0.00	82.46%
70) 1,4-Dichlorobenzene-d4	12.18	152	308749	50.00	ug/l	0.00	79.32%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	410038	49.61	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.22%	
37) 1,2-Dichloroethane-d4	6.39	65	476302	50.34	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.68%	
51) Toluene-d8	8.41	98	1706498	53.28	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.56%	
72) 4-Bromofluorobenzene	11.04	95	348747	54.76	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.52%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.42	50	59	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.02	76	332	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration

M0417027.D 8260B.M Wed Apr 25 09:37:34 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417027.D  
 Acq On : 17 Apr 2007 18:01  
 Sample : JPL32-019  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:37 2007

Vial: 72  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.68	43	64	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.		
27) 2-Butanone	5.39	43	87	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.81	83	158	N.D.		
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	6.01	56	55	N.D.		
35) Carbon Tetrachloride	0.00	117	0	N.D.		
36) 1,1-Dichloropropene	6.28	75	63	N.D.		
38) Benzene	6.41	78	895	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.	d	
41) Trichloroethene	0.00	130	0	N.D.		
42) Methylcyclohexane	0.00	83	0	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	7.59	93	56	N.D.		
45) Methyl methacrylate	7.40	41	56	N.D.		
46) Bromodichloromethane	0.00	83	0	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	8.47	92	511	N.D.		
53) trans-1,3-Dichloropropene	8.79	75	70	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	8.89	97	64	N.D.		
56) Tetrachloroethene	0.00	166	0	N.D.		
57) 1,3-Dichloropropane	9.21	76	60	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.	d	
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	9.90	112	160	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	10.22	131	62	N.D.		
64) Ethylbenzene	9.99	91	1982	N.D.		
65) m,p-Xylene	10.10	106	1317	0.94	ug/l	90
66) o-xylene	10.50	106	530	N.D.		
67) Styrene	0.00	104	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0417027.D 8260B.M Wed Apr 25 09:37:35 2007

*[Handwritten Signature]*  
 Page 2  
 VOA - 151



Quantitation Report

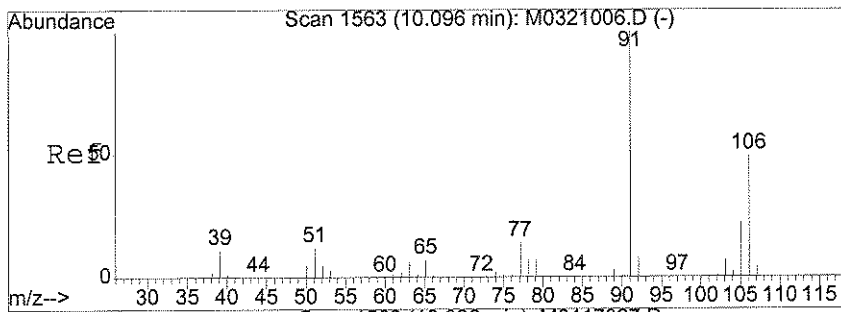
Data File : X:\MSVOA\MOBY\041707\M0417027.D  
 Acq On : 17 Apr 2007 18:01  
 Sample : JPL32-019  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:37 2007

Vial: 72  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

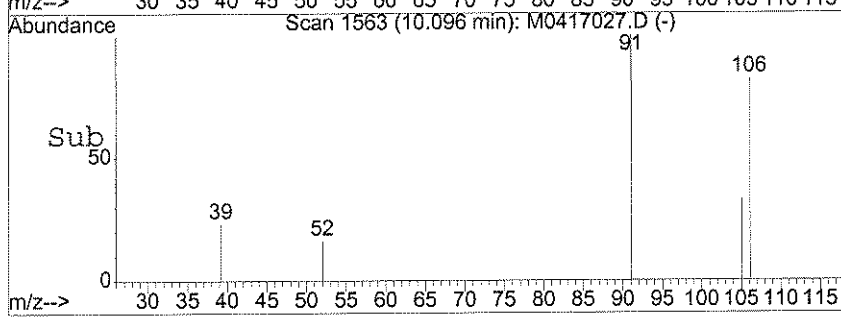
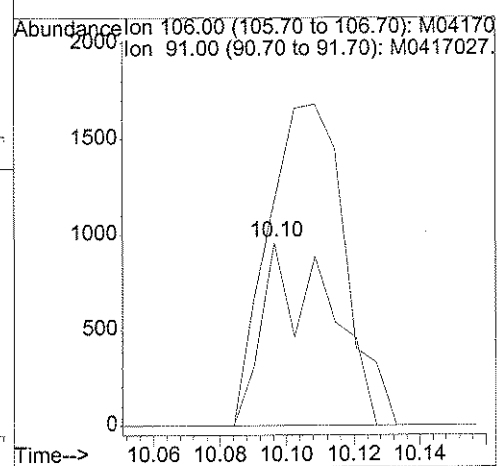
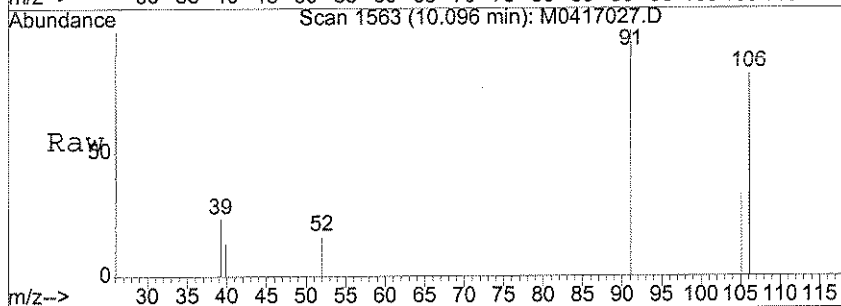
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	170		N.D.	
69) Isopropylbenzene	10.86	105	69		N.D.	
71) trans-1,4-Dichloro-2-buten	11.01	53	63		N.D.	
73) Bromobenzene	11.05	156	63		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	241		N.D.	
75) 1,2,3-Trichloropropane	0.00	110	0		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.27	91	75		N.D.	
78) 4-Chlorotoluene	11.27	91	75		N.D.	
79) 1,3,5-Trimethylbenzene	11.57	105	70		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	64		N.D.	
82) sec-butylbenzene	11.81	105	64		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



#65  
 m,p-Xylene  
 Concen: 0.94 ug/l  
 RT: 10.10 min Scan# 1563  
 Delta R.T. -0.01 min  
 Lab File: M0417027.D  
 Acq: 17 Apr 2007 18:01

Tgt Ion	Resp	Lower	Upper
106	1317		
106	100		
91	203.7	169.6	209.6



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL32  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-020  
 Lab File ID: M0417028.D  
 Date Collected: 04/09/2007  
 Date/Time Analyzed: 04/17/2007 18:25  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-020

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_(uL)

Soil Aliquot Volume: \_\_\_\_\_(uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL32-020

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417028.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/09/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/17/2007 18:25

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

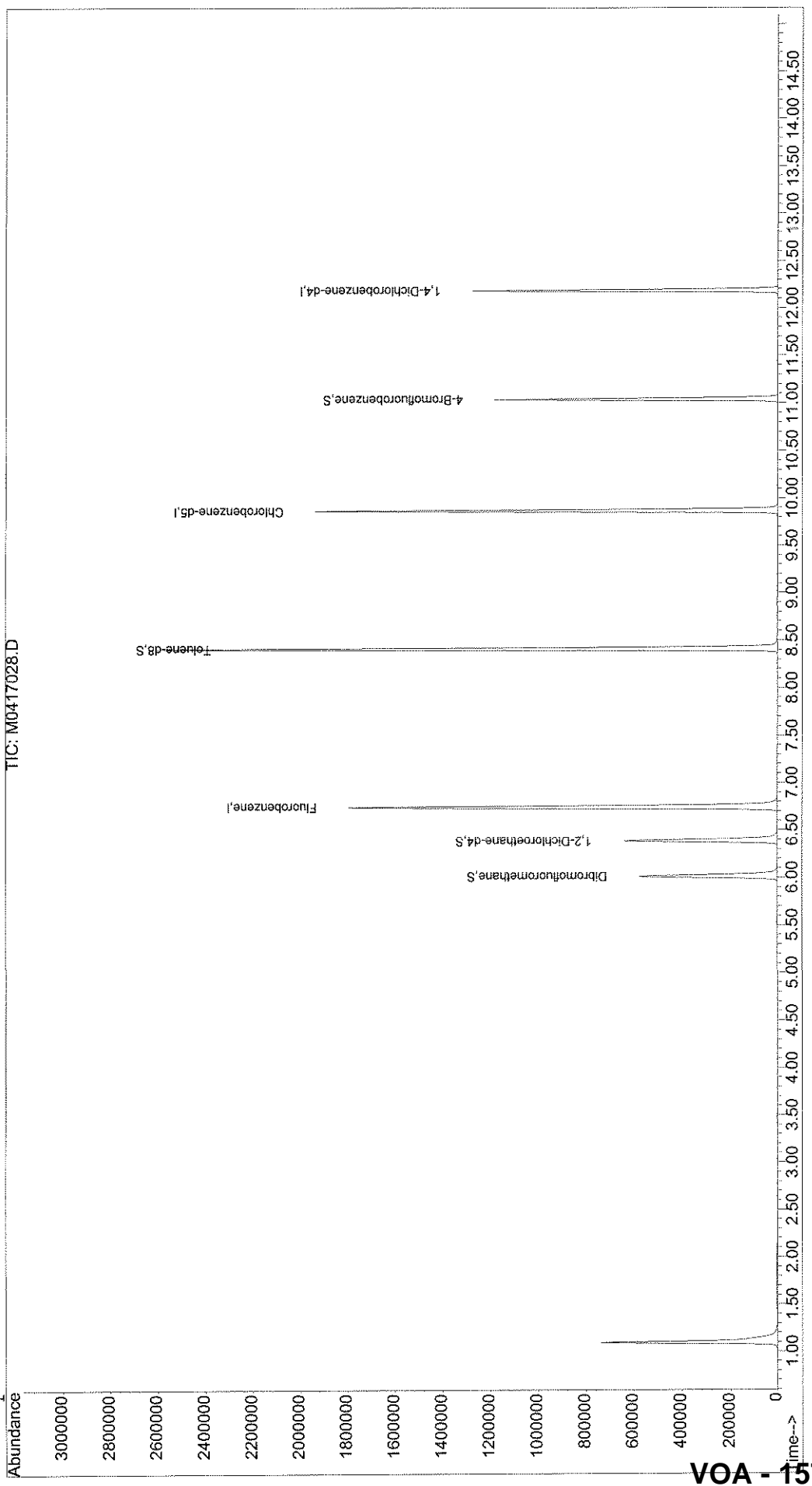
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417028.D Vial: 73  
Acq On : 17 Apr 2007 18:25 Operator: DGA  
Sample : JPL32-020 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 25 9:38 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 24 12:05:10 2007  
Response via : Initial Calibration



VOA - 157

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417028.D  
 Acq On : 17 Apr 2007 18:25  
 Sample : JPL32-020  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:38 2007

Vial: 73  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	1738103	50.00	ug/l	0.00	89.05%
50) Chlorobenzene-d5	9.87	82	571368	50.00	ug/l	0.00	81.24%
70) 1,4-Dichlorobenzene-d4	12.18	152	302362	50.00	ug/l	0.00	77.68%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	407530	49.70	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.40%	
37) 1,2-Dichloroethane-d4	6.39	65	471626	50.24	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.48%	
51) Toluene-d8	8.41	98	1691090	53.59	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	107.18%	
72) 4-Bromofluorobenzene	11.04	95	346176	55.50	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	111.00%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.41	50	62	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.86	96	80	N.D.	
6) Chloroethane	1.94	64	59	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	0.00	43	0	N.D.	d
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	634	N.D.	
15) Allyl chloride	3.22	76	59	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	3.48	84	1658	N.D.	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	3.90	53	57	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417028.D  
 Acq On : 17 Apr 2007 18:25  
 Sample : JPL32-020  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:38 2007

Vial: 73  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.53	43	55		N.D.	
28) Propionitrile	5.45	54	65		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	5.87	97	62		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	873		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	7.46	63	86		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.85	83	87		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.25	75	59		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	518		N.D.	
53) trans-1,3-Dichloropropene	8.92	75	93		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.92	97	59		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	9.04	76	80		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.85	112	78		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.88	91	2341		N.D.	
65) m,p-Xylene	0.00	106	0		N.D.	d
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	0.00	104	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0417028.D 8260B.M Wed Apr 25 09:38:38 2007



Quantitation Report

Data File : X:\MSVOA\MOBY\041707\M0417028.D  
 Acq On : 17 Apr 2007 18:25  
 Sample : JPL32-020  
 Misc : #3 5ml +IS/SS(524)  
 MS Integration Params: rteint.p  
 Quant Time: Apr 25 9:38 2007

Vial: 73  
 Operator: DGA  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 24 12:05:10 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	123		N.D.	
69) Isopropylbenzene	10.95	105	59		N.D.	
71) trans-1,4-Dichloro-2-buten	11.03	53	76		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	342		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	130		N.D.	
76) n-Propylbenzene	11.27	120	55		N.D.	
77) 2-Chlorotoluene	11.27	91	66		N.D.	
78) 4-Chlorotoluene	11.27	91	66		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.77	119	135		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	65		N.D.	
82) sec-butylbenzene	11.81	105	65		N.D.	
83) 1,3-Dichlorobenzene	0.00	146	0		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	0.00	146	0		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

**TIC FORM**

SDG JPL32

VOLATILES ANALYSIS

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-24-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-002  
 Lab File ID: M0417010.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417010.D Vial: 55  
Acq On : 17 Apr 2007 11:12 Operator: DGA  
Sample : JPL32-002 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417010.D 8260B.M Wed Apr 18 09:41:36 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-24-2

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-003  
 Lab File ID: M0417011.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417011.D Vial: 56  
Acq On : 17 Apr 2007 11:37 Operator: DGA  
Sample : JPL32-003 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417011.D 8260B.M Wed Apr 18 09:43:05 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-24-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-004  
 Lab File ID: M0417012.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417012.D Vial: 57  
Acq On : 17 Apr 2007 12:01 Operator: DGA  
Sample : JPL32-004 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417012.D 8260B.M Wed Apr 18 09:45:24 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-5-1Q07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-005  
 Lab File ID: M0417013.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417013.D Vial: 58  
Acq On : 17 Apr 2007 12:25 Operator: DGA  
Sample : JPL32-005 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417013.D 8260B.M Wed Apr 18 09:47:02 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-006  
 Lab File ID: M0417014.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417014.D Vial: 59  
Acq On : 17 Apr 2007 12:49 Operator: DGA  
Sample : JPL32-006 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417014.D 8260B.M Wed Apr 18 09:52:00 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-8-4/5/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-007  
 Lab File ID: M0417015.D  
 Date Collected: 04/06/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417015.D Vial: 60  
Acq On : 17 Apr 2007 13:13 Operator: DGA  
Sample : JPL32-007 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417015.D 8260B.M Wed Apr 18 11:55:06 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-22-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-008  
 Lab File ID: M0417016.D  
 Date Collected: 04/07/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417016.D Vial: 61  
Acq On : 17 Apr 2007 13:37 Operator: DGA  
Sample : JPL32-008 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417016.D 8260B.M Wed Apr 18 10:56:08 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-22-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL32-009

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/07/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/17/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417017.D Vial: 62  
Acq On : 17 Apr 2007 14:01 Operator: DGA  
Sample : JPL32-009 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417017.D 8260B.M Wed Apr 18 10:57:48 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-22-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-010  
 Lab File ID: M0417018.D  
 Date Collected: 04/07/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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04				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417018.D Vial: 63  
Acq On : 17 Apr 2007 14:25 Operator: DGA  
Sample : JPL32-010 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417018.D 8260B.M Wed Apr 18 10:59:21 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-6-1Q07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-011  
 Lab File ID: M0417019.D  
 Date Collected: 04/07/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417019.D Vial: 64  
Acq On : 17 Apr 2007 14:49 Operator: DGA  
Sample : JPL32-011 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417019.D 8260B.M Wed Apr 18 11:00:49 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-012  
 Lab File ID: M0417020.D  
 Date Collected: 04/07/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417020.D Vial: 65  
Acq On : 17 Apr 2007 15:13 Operator: DGA  
Sample : JPL32-012 Inst : MOBY  
Misc : #1 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417020.D 8260B.M Wed Apr 18 11:51:01 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-9-4/6/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-013  
 Lab File ID: M0417021.D  
 Date Collected: 04/07/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417021.D Vial: 66  
Acq On : 17 Apr 2007 15:37 Operator: DGA  
Sample : JPL32-013 Inst : MOBY  
Misc : #1 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417021.D 8260B.M Tue Apr 24 15:57:19 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-25-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-014  
 Lab File ID: M0417022.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417022.D Vial: 67  
Acq On : 17 Apr 2007 16:01 Operator: DGA  
Sample : JPL32-014 Inst : MOBY  
Misc : #1 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417022.D 8260B.M Tue Apr 24 15:57:37 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-25-4

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-015  
 Lab File ID: M0417023.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417023.D Vial: 68  
Acq On : 17 Apr 2007 16:25 Operator: DGA  
Sample : JPL32-015 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417023.D 8260B.M Tue Apr 24 15:57:47 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-25-3

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-016  
 Lab File ID: M0417024.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417024.D Vial: 69  
Acq On : 17 Apr 2007 16:49 Operator: DGA  
Sample : JPL32-016 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417024.D 8260B.M Tue Apr 24 15:57:57 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-25-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL32

Run Sequence: R016952

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL32-017

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0417025.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/10/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/17/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417025.D Vial: 70  
Acq On : 17 Apr 2007 17:13 Operator: DGA  
Sample : JPL32-017 Inst : MOBY  
Misc : #2 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417025.D 8260B.M Tue Apr 24 15:58:10 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-25-1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-018  
 Lab File ID: M0417026.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417026.D Vial: 71  
Acq On : 17 Apr 2007 17:37 Operator: DGA  
Sample : JPL32-018 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0417026.D 8260B.M Tue Apr 24 15:58:21 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-019  
 Lab File ID: M0417027.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417027.D Vial: 72  
Acq On : 17 Apr 2007 18:01 Operator: DGA  
Sample : JPL32-019 Inst : MOBY  
Misc : #3 5ml +IS/SS(524) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417027.D 8260B.M Tue Apr 24 15:58:34 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-10-4/9/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL)      mL  
 Level: (LOW/MED)       
 % Moisture: not dec.       
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume:      (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: JPL32-020  
 Lab File ID: M0417028.D  
 Date Collected: 04/10/2007  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume:      (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417028.D                   Vial: 73  
Acq On    : 17 Apr 2007  18:25                           Operator: DGA  
Sample    : JPL32-020                                   Inst     : MOBY  
Misc      : #3 5ml +IS/SS(524)                        Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title      : VOA 8260- 5ML Calibration 5973M  
Library    : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417028.D  8260B.M     Tue Apr 24 15:58:51 2007



1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041707MVOWM1

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL32  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R016952  
 Lab Sample ID: B041707MVOWM1  
 Lab File ID: M0417009.D  
 Date Collected: \_\_\_\_\_  
 Date Analyzed: 04/17/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041707\M0417009.D Vial: 55  
Acq On : 17 Apr 2007 10:38 Operator: DGA  
Sample : B041707MVOWM1 Inst : MOBY  
Misc : 5ml PFW+IS/SS(MV8-38-11) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

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M0417009.D 8260B.M Wed Apr 25 12:08:03 2007

**Metals Data**

**JPL32**

## COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories Contract: JPL Groundwater MonitorinLab Code: LAUCKS SDG No.: JPL32

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-24-4	JPL32-001
MW-24-3	JPL32-002
MW-24-2	JPL32-003
MW-24-1	JPL32-004
DUPE-5-1Q07	JPL32-005
DUPE-5-1Q07MS	JPL32-005MS
DUPE-5-1Q07MSD	JPL32-005MSD
EB-8-4/5/07	JPL32-006
EB-8-4/5/07MS	JPL32-006MS
EB-8-4/5/07MSD	JPL32-006MSD
MW-22-3	JPL32-008
MW-22-2	JPL32-009
MW-22-1	JPL32-010
DUPE-6-1Q07	JPL32-011
DUPE-6-1Q07MS	JPL32-011MS
DUPE-6-1Q07MSD	JPL32-011MSD
EB-9-4/6/07	JPL32-012
EB-9-4/6/07MS	JPL32-012MS
EB-9-4/6/07MSD	JPL32-012MSD
MW-25-5	JPL32-014
MW-25-4	JPL32-015

Were ICP interelement corrections applied? Yes/No YESWere ICP background corrections applied? Yes/No NOIf yes-was raw data generated before application of background corrections? Yes/No NOComments:  
\_\_\_\_\_  
\_\_\_\_\_

I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher Name: Bill AmbacherDate: 5/1/07 Title: Inorganics/Metals Manager

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
<u>MW-25-3</u>	<u>JPL32-016</u>
<u>MW-25-2</u>	<u>JPL32-017</u>
<u>MW-25-1</u>	<u>JPL32-018</u>
<u>MW-25-1D</u>	<u>JPL32-018D</u>
<u>MW-25-1MS</u>	<u>JPL32-018MS</u>
<u>EB-10-4/9/07</u>	<u>JPL32-019</u>

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

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I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: *Bill Ambacher*

Name: Bill Ambacher

Date: 5/1/07

Title: Inorganics/Metals Manager

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-24-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-001

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	4.85		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-24-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-002

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	6.95		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-24-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-003

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	7.95		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-24-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-004

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.46		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-5-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-005

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	7.79		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-8-4/5/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-006

Level (low/med): LOW

Date Received: 04/06/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.59			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-22-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-008

Level (low/med): LOW

Date Received: 04/07/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	9.59		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-22-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-009

Level (low/med): LOW

Date Received: 04/07/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.50		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-22-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-010

Level (low/med): LOW

Date Received: 04/07/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	7.96		E	M	R017204

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-6-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-011

Level (low/med): LOW

Date Received: 04/07/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	7.99		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-9-4/6/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-012

Level (low/med): LOW

Date Received: 04/07/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.59		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-25-5

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-014

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.63		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-25-4

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-015

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	9.49		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-25-3

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-016

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	9.59		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-25-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-017

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.67		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-25-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-018

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	1.75			M	R016781

Color Before: Colorless Clarity Before: Clear Texture: \_\_\_\_\_

Color After: Colorless Clarity After: Clear Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-10-4/9/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL32

Matrix (soil/water): Water

Lab Sample ID: JPL32-019

Level (low/med): LOW

Date Received: 04/10/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.60		E	M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL32**



## COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL32

Client Identification	Lab Sample Work Order Number
MW-24-3	JPL32-002DL
MW-24-2	JPL32-003DL
MW-24-1	JPL32-004
DUPE-5-1Q07	JPL32-005
DUPE-5-1Q07MS	JPL32-005MS
DUPE-5-1Q07MSD	JPL32-005MSD
EB-8-4/5/07	JPL32-006
EB-8-4/5/07MS	JPL32-006MS
EB-8-4/5/07MSD	JPL32-006MSD
MW-22-3	JPL32-008DL
MW-22-2	JPL32-009DL
MW-22-1	JPL32-010DL
DUPE-6-1Q07	JPL32-011DL
DUPE-6-1Q07MS	JPL32-011MS
DUPE-6-1Q07MSD	JPL32-011MSD
EB-9-4/6/07	JPL32-012
EB-9-4/6/07MS	JPL32-012MS
EB-9-4/6/07MSD	JPL32-012MSD
MW-25-5	JPL32-014DL
MW-25-4	JPL32-015DL
MW-25-3	JPL32-016DL

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher

Name: Bill Ambacher

Date: 4/30/07

Title: Inorganics/Metals Manager

**COVER PAGE - INORGANIC ANALYSES DATA PACKAGE**

Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL32

Client Identification	Lab Sample Work Order Number
MW-25-2	JPL32-017DL
MW-25-1	JPL32-018DL
EB-10-4/9/07	JPL32-019

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher

Name: Bill Ambacher

Date: 4/30/07

Title: Inorganics/Metals Manager

## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-24-3 **Date/Time Collected:** 04/05/2007 08:07  
**Lab Sample ID:** JPL32-002 **Date/Time Received:** 04/06/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-24-2 **Date/Time Collected:** 04/05/2007 08:41  
**Lab Sample ID:** JPL32-003 **Date/Time Received:** 04/06/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	51		2.0	0.28	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL32  
Sample Number: MW-24-1 Date/Time Collected: 04/05/2007 09:16  
Lab Sample ID: JPL32-004 Date/Time Received: 04/06/2007 08:30  
Method: E300.0 Unit: mg/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Nitrate - N	14797-55-8	1	2.7		0.040	0.055	04/06/2007	04/06/2007	R016613
Nitrite - N	14797-65-0	1	0.050	U	0.050	0.017	04/06/2007	04/06/2007	R016613
Sulfate as SO4	14808-79-8	10	30		10	1.7	04/06/2007	04/06/2007	R016613
Chloride	16887-00-6	10	28		2.0	0.76	04/06/2007	04/06/2007	R016613
Orthophosphate	7723-14-0	1	0.10	U	0.10	0.33	04/06/2007	04/06/2007	R016613

Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	50	1900		50	7.0	04/23/2007	04/24/2007	R017072

**Laucks Testing Laboratories, Inc.**

**Final Results**

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** DUPE-5-1Q07 **Date/Time Collected:** 04/05/2007 00:00  
**Lab Sample ID:** JPL32-005 **Date/Time Received:** 04/06/2007 08:30  
**Method:** E300.0 **Unit:** mg/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Nitrate - N	14797-55-8	1	2.7		0.040	0.055	04/06/2007	04/06/2007	R016613
Nitrite - N	14797-65-0	1	0.050	U	0.050	0.017	04/06/2007	04/06/2007	R016613
Sulfate as SO4	14808-79-8	10	34		10	1.7	04/06/2007	04/06/2007	R016613
Chloride	16887-00-6	10	32		2.0	0.76	04/06/2007	04/06/2007	R016613
Orthophosphate	7723-14-0	1	0.10	U	0.10	0.33	04/06/2007	04/06/2007	R016613

**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	50	2000		50	7.0	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** EB-8-4/5/07 **Date/Time Collected:** 04/05/2007 09:02  
**Lab Sample ID:** JPL32-006 **Date/Time Received:** 04/06/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/23/2007	04/24/2007	R017072



Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL32  
Sample Number: MW-22-3 Date/Time Collected: 04/06/2007 07:42  
Lab Sample ID: JPL32-008 Date/Time Received: 04/07/2007 09:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL32  
Sample Number: MW-22-2 Date/Time Collected: 04/06/2007 08:38  
Lab Sample ID: JPL32-009 Date/Time Received: 04/07/2007 09:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-22-1 **Date/Time Collected:** 04/06/2007 09:23  
**Lab Sample ID:** JPL32-010 **Date/Time Received:** 04/07/2007 09:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** DUPE-6-1Q07 **Date/Time Collected:** 04/06/2007 00:00  
**Lab Sample ID:** JPL32-011 **Date/Time Received:** 04/07/2007 09:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	2.5	U	2.5	0.35	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** EB-9-4/6/07 **Date/Time Collected:** 04/06/2007 08:51  
**Lab Sample ID:** JPL32-012 **Date/Time Received:** 04/07/2007 09:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/23/2007	04/24/2007	R017072

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-25-5 **Date/Time Collected:** 04/09/2007 08:07  
**Lab Sample ID:** JPL32-014 **Date/Time Received:** 04/10/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	04/20/2007	04/21/2007	R016992

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-25-4 **Date/Time Collected:** 04/09/2007 08:36  
**Lab Sample ID:** JPL32-015 **Date/Time Received:** 04/10/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2.5	7.5		2.5	0.35	04/20/2007	04/21/2007	R016992

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL32  
Sample Number: MW-25-3 Date/Time Collected: 04/09/2007 09:04  
Lab Sample ID: JPL32-016 Date/Time Received: 04/10/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	9.3		4.0	0.56	04/20/2007	04/21/2007	R016992



Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL32  
Sample Number: MW-25-2 Date/Time Collected: 04/09/2007 09:31  
Lab Sample ID: JPL32-017 Date/Time Received: 04/10/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	14		4.0	0.56	04/20/2007	04/21/2007	R016992

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** MW-25-1 **Date/Time Collected:** 04/09/2007 10:17  
**Lab Sample ID:** JPL32-018 **Date/Time Received:** 04/10/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	04/20/2007	04/21/2007	R016992

**Laucks Testing Laboratories, Inc.**

**Final Results**

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL32  
**Sample Number:** EB-10-4/9/07 **Date/Time Collected:** 04/09/2007 10:04  
**Lab Sample ID:** JPL32-019 **Date/Time Received:** 04/10/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	04/20/2007	04/21/2007	R016992

**SAMPLE DATA**

SDG JPL33

VOLATILES ANALYSIS

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-2

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-001

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/10/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 19:12

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-001  
 Lab File ID: M0419018.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 19:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-2

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-001  
 Lab File ID: M0419018.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 19:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

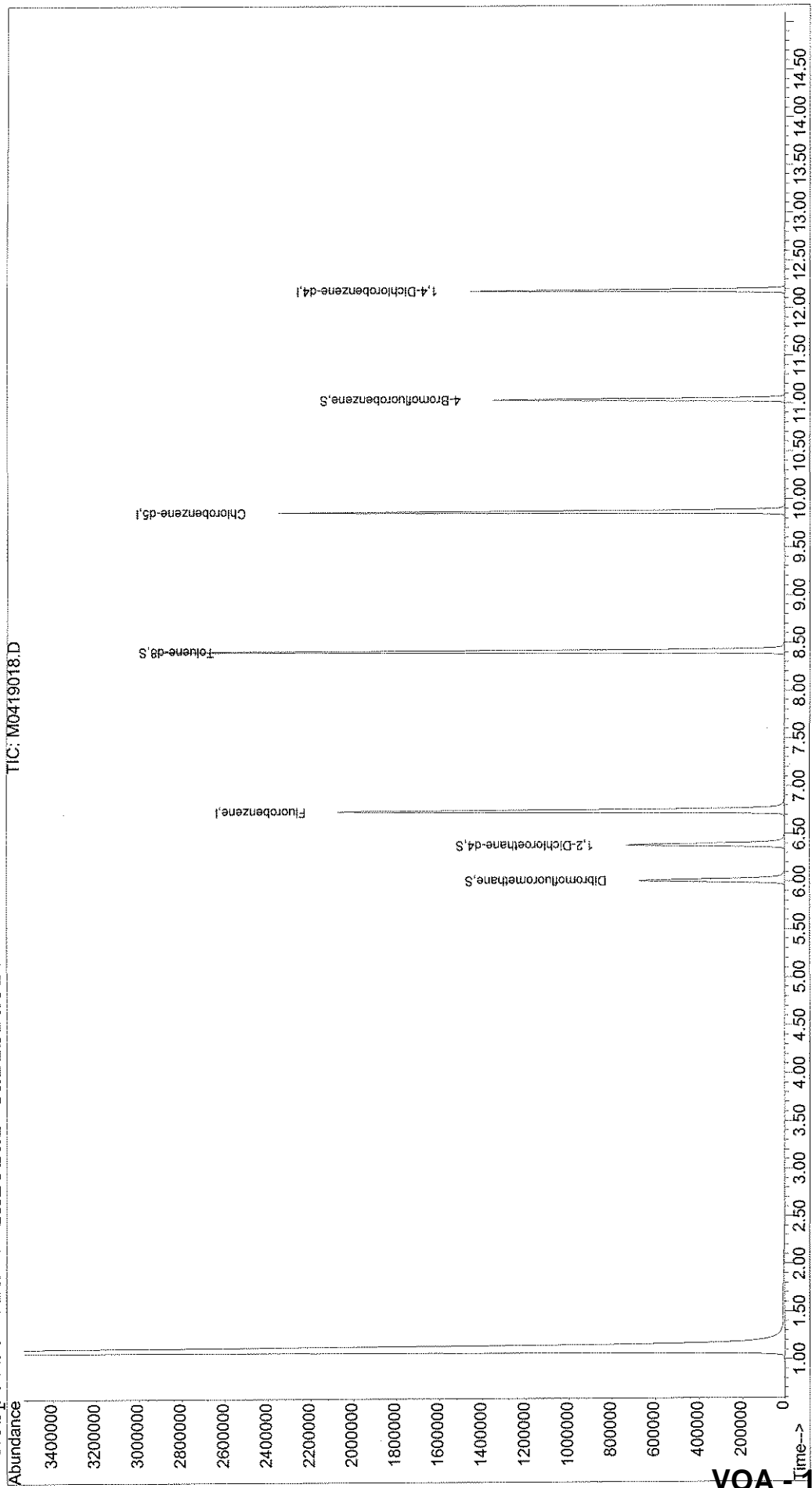
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419018.D Vial: 61  
Acq On : 19 Apr 2007 19:12 Operator: LH  
Sample : JPL33-001 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:09 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA-18



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419018.D  
 Acq On : 19 Apr 2007 19:12  
 Sample : JPL33-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:09 2007

Vial: 61  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1996280	50.00	ug/l	0.00 102.27%
50) Chlorobenzene-d5	9.87	82	671472	50.00	ug/l	0.00 95.47%
70) 1,4-Dichlorobenzene-d4	12.18	152	363372	50.00	ug/l	0.00 93.36%

System Monitoring Compounds

33) Dibromofluoromethane	6.02	111	475145	50.45	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.90%
37) 1,2-Dichloroethane-d4	6.39	65	538532	49.95	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.90%
51) Toluene-d8	8.41	98	1959843	52.85	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.70%
72) 4-Bromofluorobenzene	11.04	95	411700	54.93	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.86%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.85	96	58	N.D.	
6) Chloroethane	1.98	64	66	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	2.82	96	66	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	2.86	43	65	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	2951	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	3.77	96	64	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419018.D  
 Acq On : 19 Apr 2007 19:12  
 Sample : JPL33-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:09 2007

Vial: 61  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.67	43	58		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.50	43	66		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.42	78	793		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.36	93	57		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.85	83	69		N.D.	
47) 2-Chloroethyl vinyl ether	8.15	63	61		N.D.	
48) cis-1,3-Dichloropropene	8.08	75	84		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.47	92	427		N.D.	
53) trans-1,3-Dichloropropene	8.60	75	57		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.88	97	57		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	9.14	76	62		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	378		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.10	91	1193		N.D.	
65) m,p-Xylene	10.11	106	631		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.51	104	58		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419018.D  
 Acq On : 19 Apr 2007 19:12  
 Sample : JPL33-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:09 2007

Vial: 61  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	121		N.D.	
69) Isopropylbenzene	10.88	105	64		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.39	83	55		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	190		N.D.	
76) n-Propylbenzene	11.28	120	61		N.D.	
77) 2-Chlorotoluene	11.27	91	249		N.D.	
78) 4-Chlorotoluene	11.47	91	107		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.77	119	126		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	120		N.D.	
82) sec-butylbenzene	11.98	105	138		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	69		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	191		N.D.	
86) 1,2-Dichlorobenzene	12.65	146	55		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.48	157	58		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-1

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/10/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 19:36

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-002  
 Lab File ID: M0419019.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 19:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-26-1

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-002  
 Lab File ID: M0419019.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 19:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

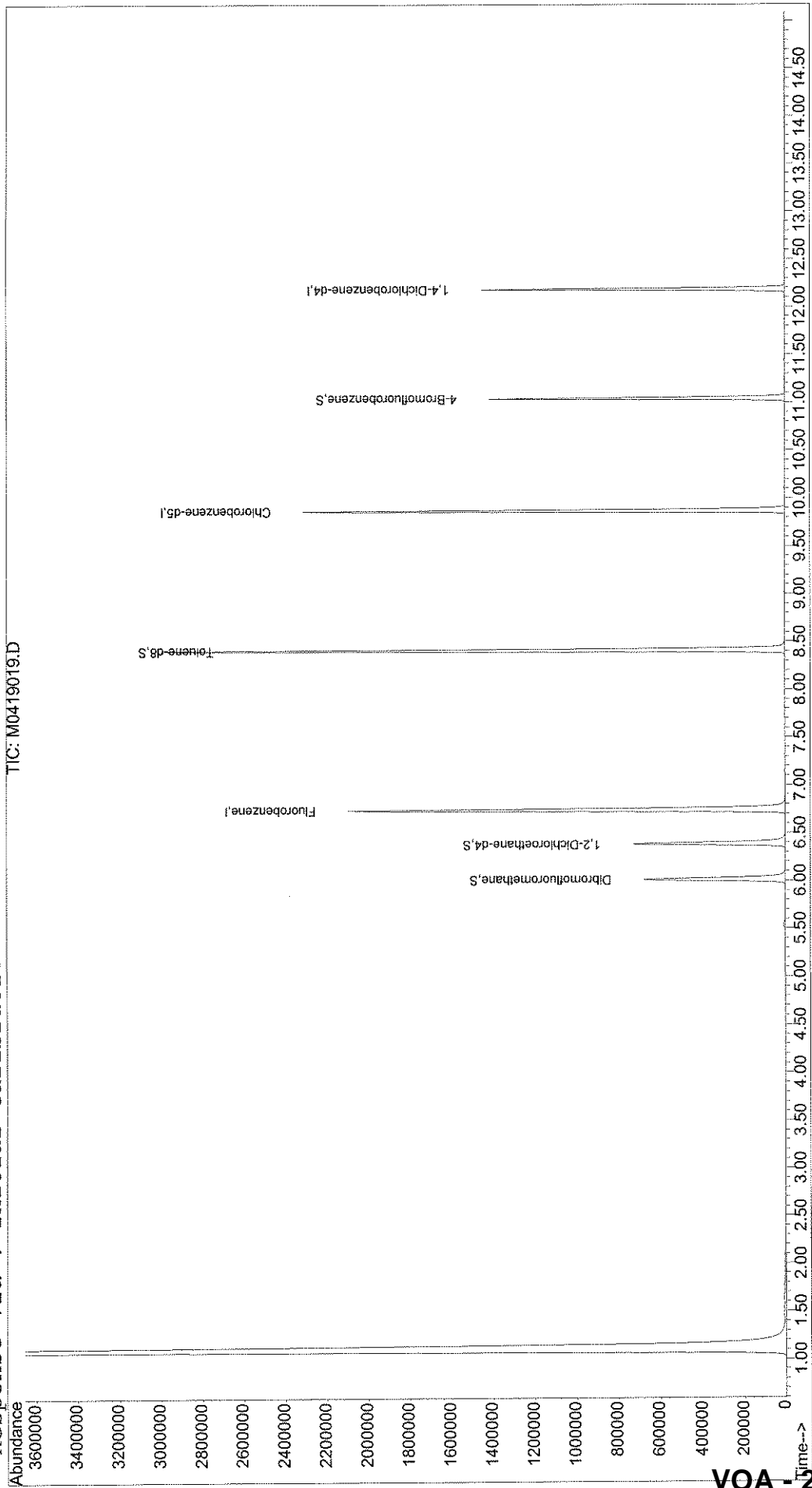
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419019.D Vial: 62  
Acq On : 19 Apr 2007 19:36 Operator: LH  
Sample : JPL33-002 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:10 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419019.D  
 Acq On : 19 Apr 2007 19:36  
 Sample : JPL33-002  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:10 2007

Vial: 62  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1982586	50.00	ug/l	0.00 101.57%
50) Chlorobenzene-d5	9.87	82	683033	50.00	ug/l	0.00 97.12%
70) 1,4-Dichlorobenzene-d4	12.18	152	364178	50.00	ug/l	0.00 93.56%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	476248	50.92	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	101.84%	
37) 1,2-Dichloroethane-d4	6.38	65	541795	50.60	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	101.20%	
51) Toluene-d8	8.41	98	1968697	52.19	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	104.38%	
72) 4-Bromofluorobenzene	11.04	95	413064	54.99	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery	=	109.98%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.83	96	90	N.D.	
6) Chloroethane	1.97	64	61	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	2.90	43	68	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.05	76	1078	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	3.33	40	55	N.D.	
17) Methyl Acetate	0.00	43	0	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	3.92	73	342	N.D.	
22) 1,1-Dichloroethane	4.54	63	272	N.D.	

(#) = qualifier out of range (m) = manual integration



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419019.D  
 Acq On : 19 Apr 2007 19:36  
 Sample : JPL33-002  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:10 2007

Vial: 62  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	4.75	53	65		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.39	96	57		N.D.	
27) 2-Butanone	5.44	43	56		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.56	128	63		N.D.	
30) Methacrylonitrile	5.83	41	55		N.D.	
31) Chloroform	5.81	83	990		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.09	117	63		N.D.	
36) 1,1-Dichloropropene	6.15	75	73		N.D.	
38) Benzene	6.41	78	635		N.D.	
39) 1,2-Dichloroethane	6.41	62	269		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	830		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.73	83	122		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	768		N.D.	
53) trans-1,3-Dichloropropene	8.85	75	62		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.87	97	64		N.D.	
56) Tetrachloroethene	9.02	166	1163		N.D.	
57) 1,3-Dichloropropane	8.92	76	61		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	155		N.D.	
62) 1-Chlorohexane	9.86	91	2376		N.D.	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	989		N.D.	
65) m,p-Xylene	10.11	106	1124		N.D.	
66) o-xylene	10.50	106	69		N.D.	
67) Styrene	10.55	104	55		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419019.D 8260B.M Fri Apr 20 10:11:03 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419019.D  
 Acq On : 19 Apr 2007 19:36  
 Sample : JPL33-002  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:10 2007

Vial: 62  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	241		N.D.	
69) Isopropylbenzene	10.86	105	58		N.D.	
71) trans-1,4-Dichloro-2-buten	10.74	53	58		N.D.	
73) Bromobenzene	11.09	156	65		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	321		N.D.	
75) 1,2,3-Trichloropropane	11.03	110	159		N.D.	
76) n-Propylbenzene	11.28	120	86		N.D.	
77) 2-Chlorotoluene	11.37	91	57		N.D.	
78) 4-Chlorotoluene	11.37	91	57		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	63		N.D.	
82) sec-butylbenzene	11.99	105	167		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	151		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	151		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	68		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.33	225	67		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/10/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 20:00

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-003  
 Lab File ID: M0419020.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 20:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

EB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-003  
 Lab File ID: M0419020.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 20:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

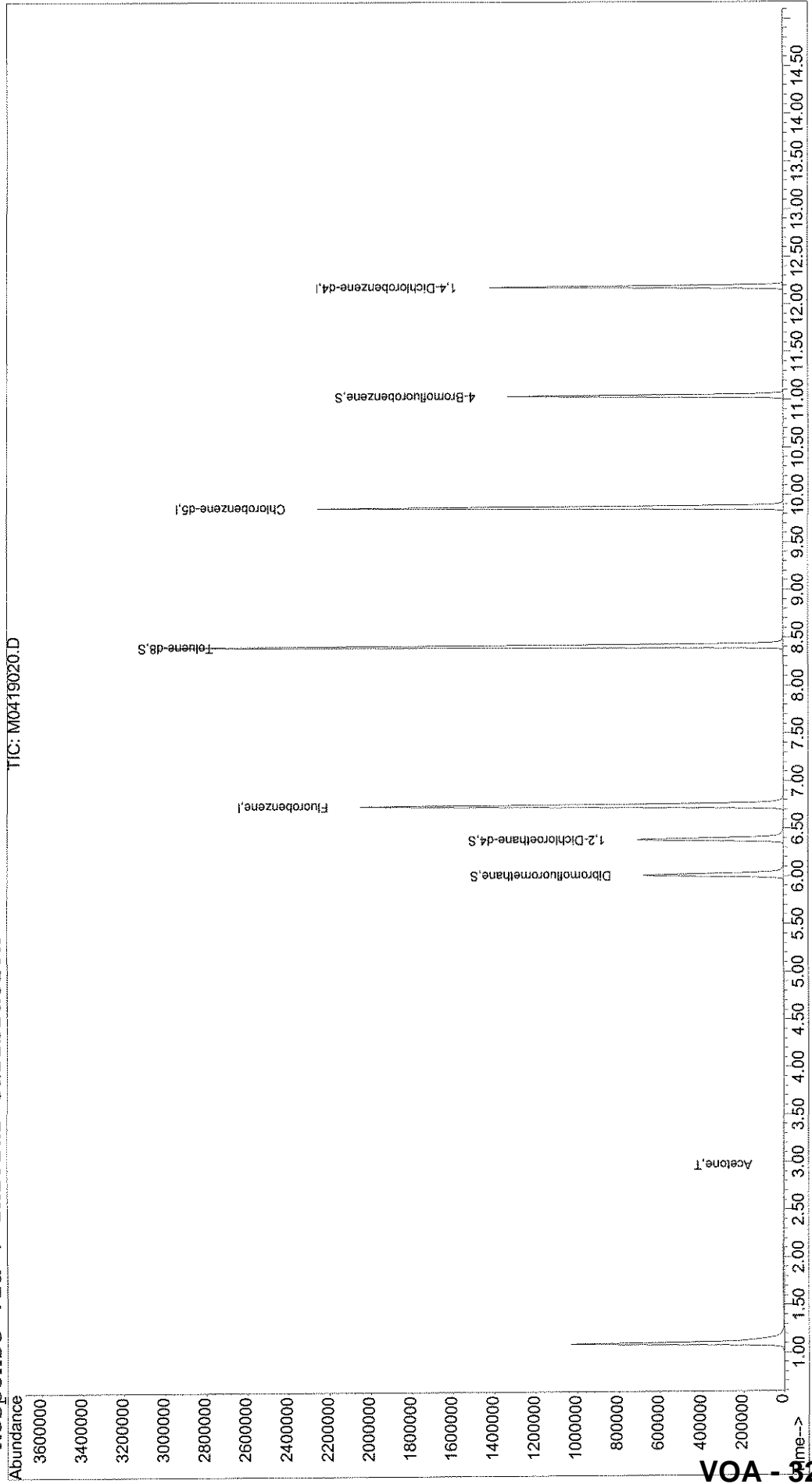
CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
96-12-8	1,2-Dibromo-3-chloropropane	0.50		U
120-82-1	1,2,4-Trichlorobenzene	0.50		U
87-68-3	Hexachlorobutadiene	0.50		U
91-20-3	Naphthalene	0.50		U
87-61-6	1,2,3-Trichlorobenzene	0.50		U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419020.D Vial: 63  
Acq On : 19 Apr 2007 20:00 Operator: LH  
Sample : JPL33-003 Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:13 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 32

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419020.D  
 Acq On : 19 Apr 2007 20:00  
 Sample : JPL33-003  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:13 2007

Vial: 63  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Dev(Max)
						Rcv(Ar )	
1) Fluorobenzene	6.74	96	1969897	50.00	ug/l	0.00	100.92%
50) Chlorobenzene-d5	9.87	82	666978	50.00	ug/l	0.00	94.84%
70) 1,4-Dichlorobenzene-d4	12.18	152	354133	50.00	ug/l	0.00	90.98%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	462622	49.78	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.56%	
37) 1,2-Dichloroethane-d4	6.38	65	532184	50.02	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.04%	
51) Toluene-d8	8.41	98	1930060	52.39	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.78%	
72) 4-Bromofluorobenzene	11.04	95	403836	55.28	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.56%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	1.41	50	238		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	0.00	96	0		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D. d	
11) Acetone	2.97	43	3491	1.95	ug/l	# 78
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.04	76	513		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D. d	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D. d	
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	4.64	63	72		N.D.	

Handwritten note: *W 4/20/07*

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419020.D  
 Acq On : 19 Apr 2007 20:00  
 Sample : JPL33-003  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:13 2007

Vial: 63  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	4.73	53	59		N.D.	
25) 2,2-Dichloropropane	5.22	77	75		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.52	43	430		N.D.	
28) Propionitrile	5.47	54	62		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.69	41	64		N.D.	
31) Chloroform	5.82	83	626		N.D.	
32) 1,1,1-Trichloroethane	6.07	97	71		N.D.	
34) Cyclohexane	5.97	56	56		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.41	78	1036		N.D.	
39) 1,2-Dichloroethane	6.40	62	357		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.42	41	55		N.D.	
46) Bromodichloromethane	7.72	83	67		N.D.	
47) 2-Chloroethyl vinyl ether	8.12	63	63		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	610		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	8.75	69	77		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.08	166	55		N.D.	
57) 1,3-Dichloropropane	8.99	76	68		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	9.55	129	67		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	302		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	1914		N.D.	
65) m,p-Xylene	10.10	106	971		N.D.	
66) o-xylene	10.50	106	554		N.D.	
67) Styrene	0.00	104	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419020.D 8260B.M Fri Apr 20 10:13:33 2007



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419020.D  
 Acq On : 19 Apr 2007 20:00  
 Sample : JPL33-003  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:13 2007

Vial: 63  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	224		N.D.	
69) Isopropylbenzene	10.86	105	161		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.04	156	67		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	418		N.D.	
75) 1,2,3-Trichloropropane	11.08	110	64		N.D.	
76) n-Propylbenzene	11.49	120	58		N.D.	
77) 2-Chlorotoluene	11.36	91	125		N.D.	
78) 4-Chlorotoluene	11.48	91	122		N.D.	
79) 1,3,5-Trimethylbenzene	11.56	105	65		N.D.	
80) tert-Butylbenzene	11.77	119	57		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	59		N.D.	
82) sec-butylbenzene	11.98	105	127		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	68		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	68		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/10/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 17:11

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-004  
 Lab File ID: M0419013.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 17:11  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-004  
 Lab File ID: M0419013.D  
 Date Collected: 04/10/2007  
 Date/Time Analyzed: 04/19/2007 17:11  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

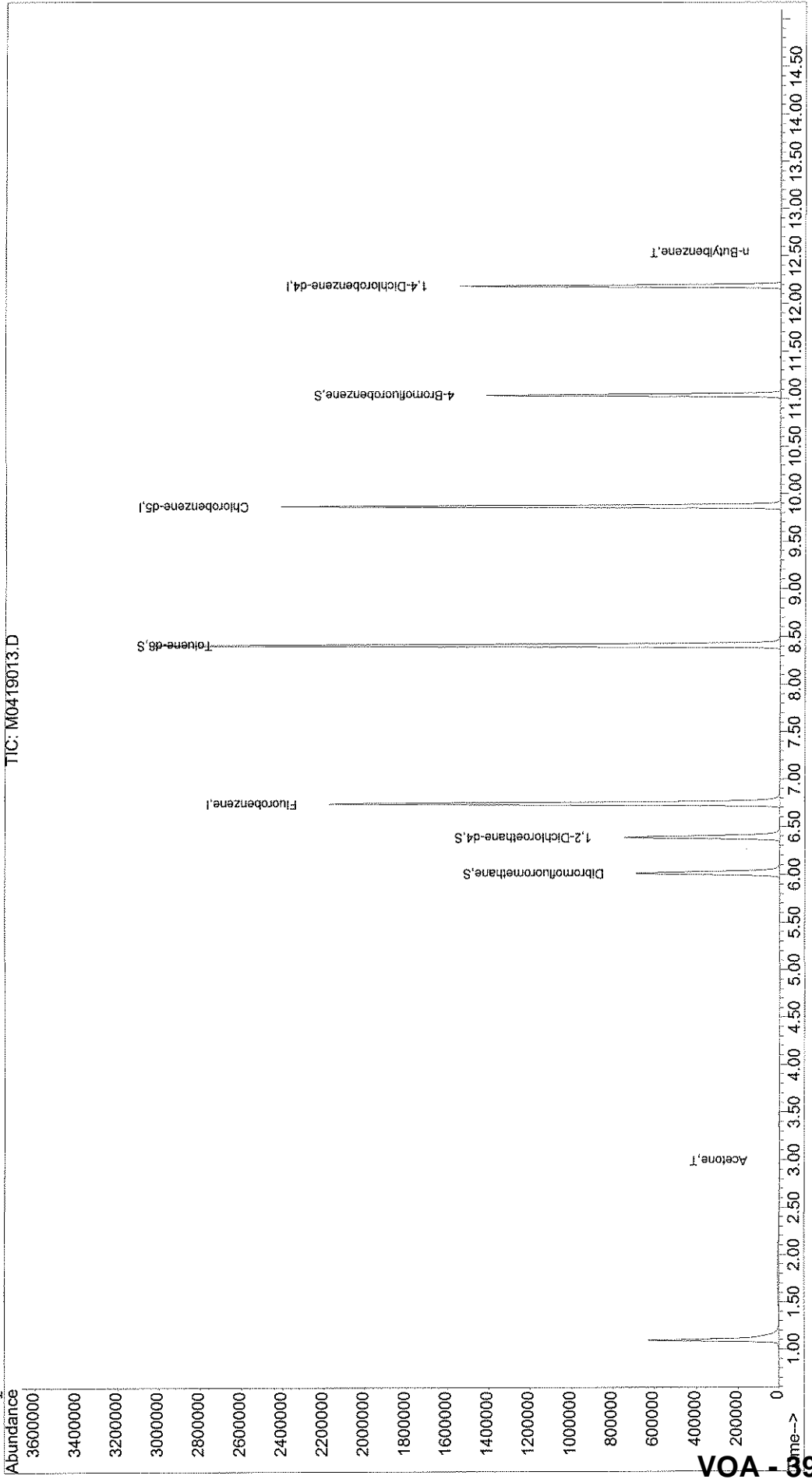
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419013.D  
Acq On : 19 Apr 2007 17:11  
Sample : JPL33-004 TB  
Misc : 5ml +IS/SS #1  
MS Integration Params: rteint.p  
Quant Time: Apr 20 9:58 2007  
Vial: 56  
Operator: LH  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 39

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419013.D  
 Acq On : 19 Apr 2007 17:11  
 Sample : JPL33-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 9:58 2007

Vial: 56  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	2071245	50.00	ug/l	0.00 106.11%
50) Chlorobenzene-d5	9.87	82	710542	50.00	ug/l	0.00 101.03%
70) 1,4-Dichlorobenzene-d4	12.19	152	375769	50.00	ug/l	0.00 96.54%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	490436	50.19	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.38%
37) 1,2-Dichloroethane-d4	6.39	65	558620	49.94	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.88%
51) Toluene-d8	8.41	98	2040518	52.00	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.00%
72) 4-Bromofluorobenzene	11.04	95	438157	56.53	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	113.06%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.45	50	126	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.84	96	62	N.D.		
6) Chloroethane	1.95	64	56	N.D.		
7) Trichlorofluoromethane	2.21	101	74	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.79	96	87	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.99	43	3096m	1.64	ug/l #	96
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.03	76	1026	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	3.32	43	58	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.86	96	127	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

NOT TARGET Analyte  
GR 5/16/07

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419013.D  
 Acq On : 19 Apr 2007 17:11  
 Sample : JPL33-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 9:58 2007

Vial: 56  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.48	43	165		N.D.	
28) Propionitrile	5.65	54	65		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.81	83	65		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	5.97	56	124		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.16	75	152		N.D.	
38) Benzene	6.40	78	993		N.D.	
39) 1,2-Dichloroethane	6.38	62	310		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	268		N.D.	
42) Methylcyclohexane	7.30	83	145		N.D.	
43) 1,2-Dichloropropane	7.39	63	79		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.67	41	56		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.17	75	57		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	967		N.D.	
53) trans-1,3-Dichloropropene	8.75	75	59		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.02	166	201		N.D.	
57) 1,3-Dichloropropane	9.25	76	55		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	9.47	107	56		N.D.	
61) Chlorobenzene	9.90	112	622		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.04	131	70		N.D.	
64) Ethylbenzene	9.98	91	744		N.D.	
65) m,p-Xylene	10.10	106	566		N.D.	
66) o-xylene	10.51	106	58		N.D.	
67) Styrene	10.52	104	138		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419013.D  
 Acq On : 19 Apr 2007 17:11  
 Sample : JPL33-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 9:58 2007

Vial: 56  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
68) Bromoform	10.74	173	264	N.D.	
69) Isopropylbenzene	10.87	105	529	N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.	
73) Bromobenzene	11.18	156	66	N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	450	N.D.	
75) 1,2,3-Trichloropropane	11.04	110	246	N.D.	
76) n-Propylbenzene	11.27	120	109	N.D.	
77) 2-Chlorotoluene	11.38	91	358	N.D.	
78) 4-Chlorotoluene	11.48	91	555	N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	539	N.D.	
80) tert-Butylbenzene	11.77	119	522	N.D.	
81) 1,2,4-Trimethylbenzene	11.83	105	57	N.D.	
82) sec-butylbenzene	11.98	105	570	N.D.	
83) 1,3-Dichlorobenzene	12.12	146	386	N.D.	
84) 4-Isopropyltoluene	0.00	119	0	N.D. d	
85) 1,4-Dichlorobenzene	12.21	146	683	N.D.	
86) 1,2-Dichlorobenzene	12.58	146	158	N.D.	
87) n-Butylbenzene	12.54	91	669	<del>0.75 ug/l #</del> 87	
88) 1,2-Dibromo-3-chloropropan	13.26	157	62	N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0	N.D. d	
90) Hexachlorobutadiene	14.32	225	255	N.D.	
91) Naphthalene	0.00	128	0	N.D. d	
92) 1,2,3-Trichlorobenzene	0.00	180	0	N.D. d	

*WA 4/20/07*



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-13

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-005  
 Lab File ID: M0419021.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 20:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.58	
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.45	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.75	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.32	J

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-13

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-005  
 Lab File ID: M0419021.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 20:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	1.5	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-13

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-005  
 Lab File ID: M0419021.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 20:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

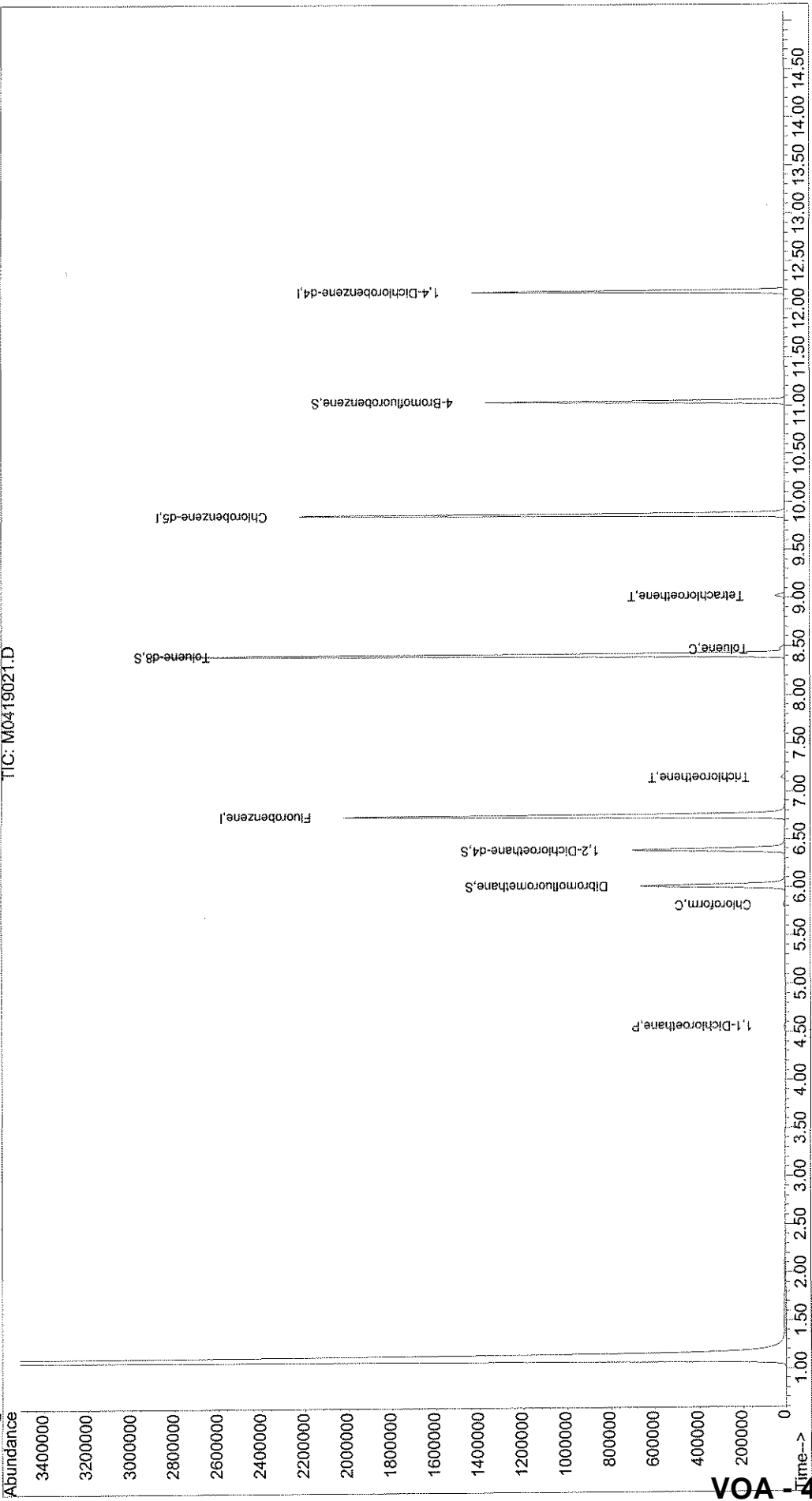
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419021.D Vial: 64  
Acq On : 19 Apr 2007 20:24 Operator: LH  
Sample : JPL33-005 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:15 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 46

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419021.D  
 Acq On : 19 Apr 2007 20:24  
 Sample : JPL33-005  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:15 2007

Vial: 64  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1951119	50.00	ug/l	0.00 99.96%
50) Chlorobenzene-d5	9.87	82	658454	50.00	ug/l	0.00 93.62%
70) 1,4-Dichlorobenzene-d4	12.18	152	352887	50.00	ug/l	0.00 90.66%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	462722	50.27	ug/l	0.00
Spiked Amount	50.000	Range 85 - 115	Recovery	=	100.54%	
37) 1,2-Dichloroethane-d4	6.39	65	524600	49.79	ug/l	0.00
Spiked Amount	50.000	Range 70 - 120	Recovery	=	99.58%	
51) Toluene-d8	8.41	98	1915051	52.66	ug/l	0.00
Spiked Amount	50.000	Range 85 - 120	Recovery	=	105.32%	
72) 4-Bromofluorobenzene	11.04	95	404992	55.64	ug/l	0.00
Spiked Amount	50.000	Range 75 - 120	Recovery	=	111.28%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.40	50	57	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.83	96	110	N.D.	
6) Chloroethane	1.99	64	58	N.D.	
7) Trichlorofluoromethane	2.19	101	1662	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.96	43	81	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.04	76	360	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.30	43	58	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	4.55	63	8743	0.58 ug/l	

(#) = qualifier out of range (m) = manual integration

M0419021.D 8260B.M Fri Apr 20 10:15:21 2007

97  
 LH 4/20/07

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419021.D  
 Acq On : 19 Apr 2007 20:24  
 Sample : JPL33-005  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:15 2007

Vial: 64  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	5.38	96	74		N.D.	
27) 2-Butanone	5.54	43	136		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.70	128	74		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	5.81	83	7070	0.45	ug/l	98
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.15	117	1332		N.D.	
36) 1,1-Dichloropropene	6.17	75	60		N.D.	
38) Benzene	6.41	78	796		N.D.	
39) 1,2-Dichloroethane	6.39	62	314		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.14	130	6588	0.75	ug/l	98
42) Methylcyclohexane	7.30	83	218		N.D.	
43) 1,2-Dichloropropane	7.40	63	61		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.52	41	65		N.D.	
46) Bromodichloromethane	7.73	83	311		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.27	75	57		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	7101	0.32	ug/l	93
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	8.97	69	56		N.D.	
55) 1,1,2-Trichloroethane	8.94	97	642		N.D.	
56) Tetrachloroethene	9.03	166	11734	1.48	ug/l	98
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	398		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	9.80	131	59		N.D.	
64) Ethylbenzene	10.01	91	344		N.D.	
65) m,p-Xylene	10.11	106	224		N.D.	
66) o-xylene	10.51	106	73		N.D.	
67) Styrene	10.39	104	63		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419021.D 8260B.M Fri Apr 20 10:15:22 2007

Quantitation Report

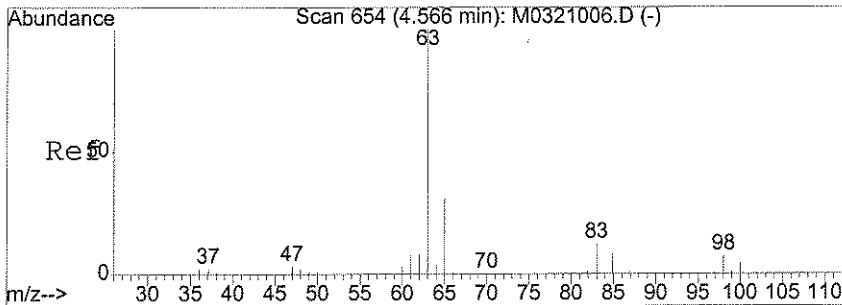
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 Acq On : 19 Apr 2007 20:24  
 Sample : JPL33-005  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:15 2007

Vial: 64  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

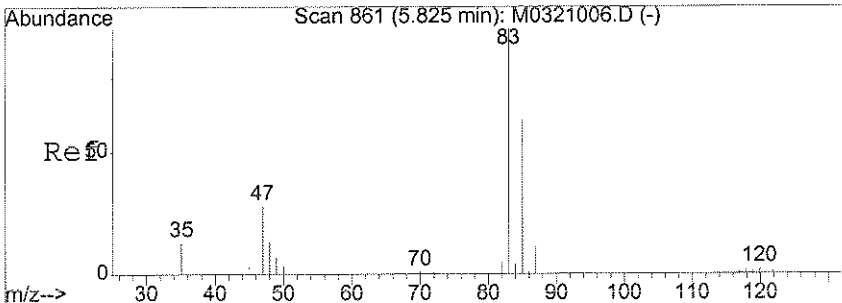
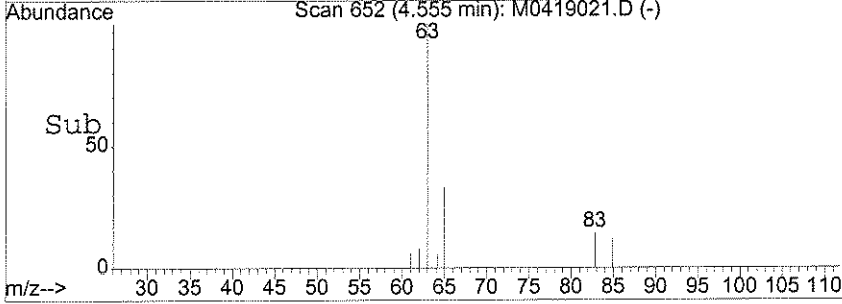
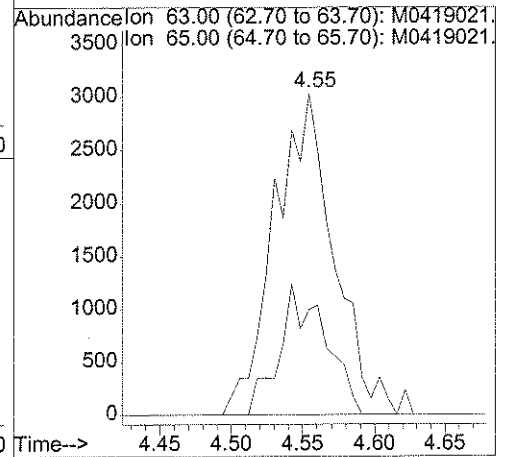
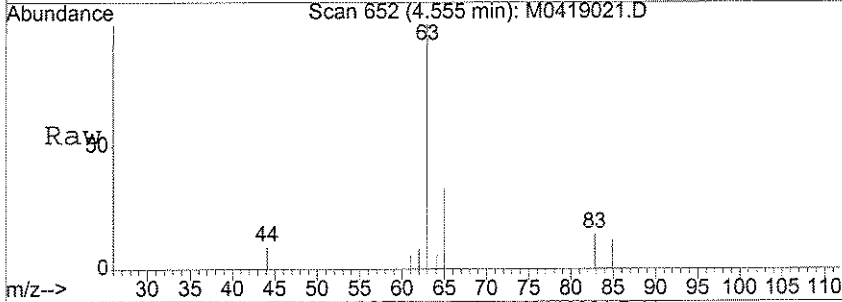
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	141		N.D.	
69) Isopropylbenzene	11.04	105	828		N.D.	
71) trans-1,4-Dichloro-2-buten	10.80	53	71		N.D.	
73) Bromobenzene	11.04	156	152		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.03	83	257		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	312		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.36	91	59		N.D.	
78) 4-Chlorotoluene	11.48	91	75		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.99	105	203		N.D.	
82) sec-butylbenzene	11.99	105	203		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	92		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.12	146	92		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	230		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.07	157	80		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



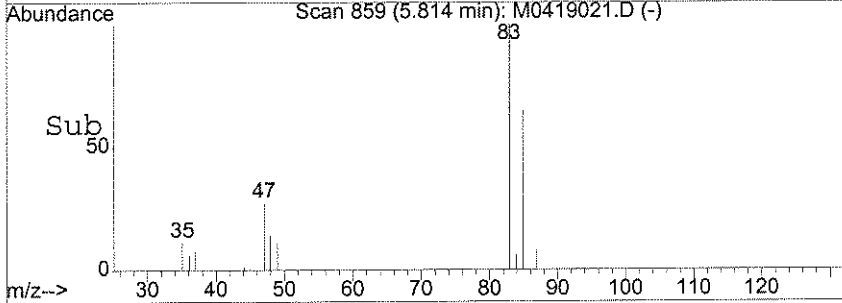
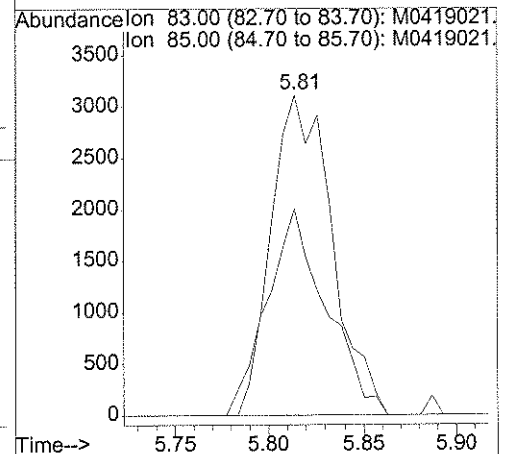
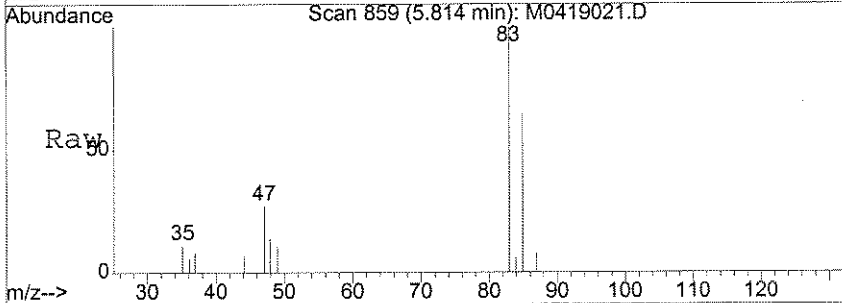
#22  
 1,1-Dichloroethane  
 Concen: 0.58 ug/l  
 RT: 4.55 min Scan# 652  
 Delta R.T. 0.01 min  
 Lab File: M0419021.D  
 Acq: 19 Apr 2007 20:24

Tgt Ion: 63 Resp: 8743  
 Ion Ratio Lower Upper  
 63 100  
 65 31.7 13.7 53.7

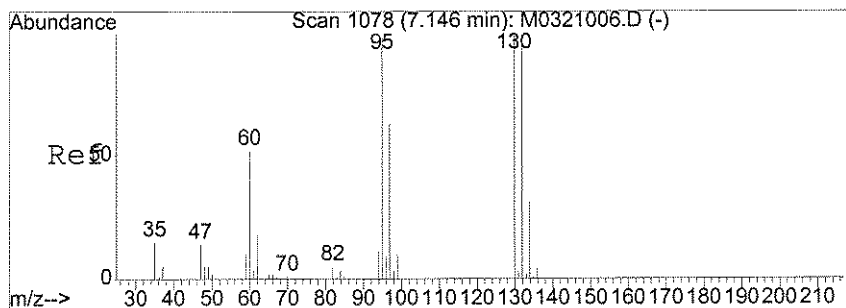


#31  
 Chloroform  
 Concen: 0.45 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419021.D  
 Acq: 19 Apr 2007 20:24

Tgt Ion: 83 Resp: 7070  
 Ion Ratio Lower Upper  
 83 100  
 85 59.6 41.2 81.2

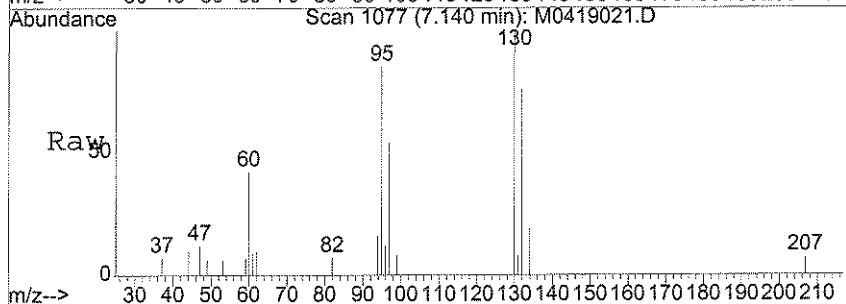




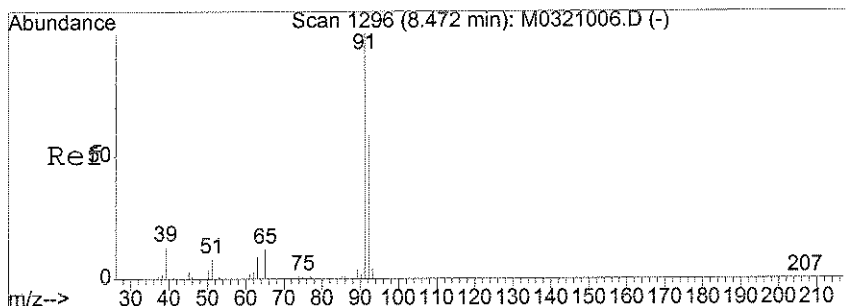
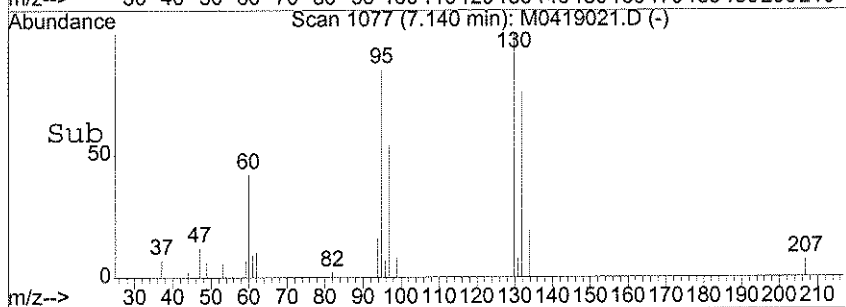
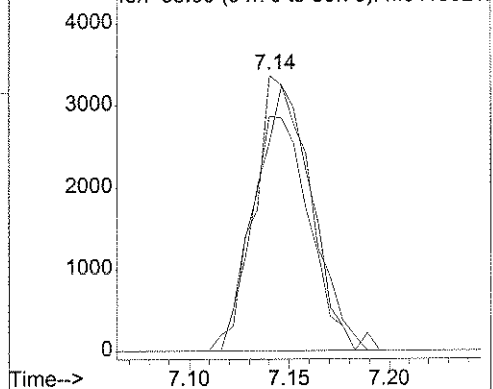


#41  
 Trichloroethene  
 Concen: 0.75 ug/l  
 RT: 7.14 min Scan# 1077  
 Delta R.T. -0.01 min  
 Lab File: M0419021.D  
 Acq: 19 Apr 2007 20:24

Tgt Ion	Resp	Lower	Upper
130	100		
132	96.2	75.0	115.0
95	87.2	69.4	109.4

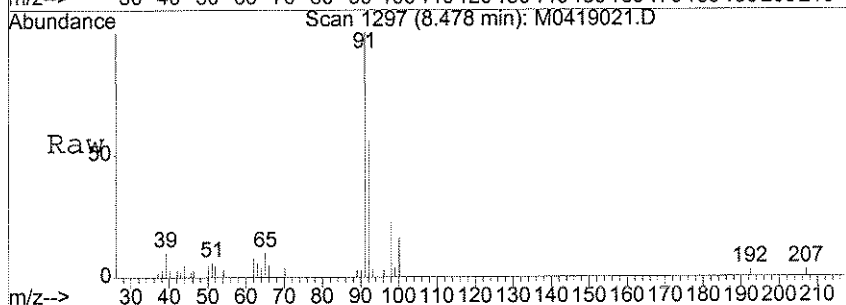


Abundance  
 Ion 130.00 (129.70 to 130.70): M04190  
 Ion 132.00 (131.70 to 132.70): M04190  
 Ion 95.00 (94.70 to 95.70): M0419021.

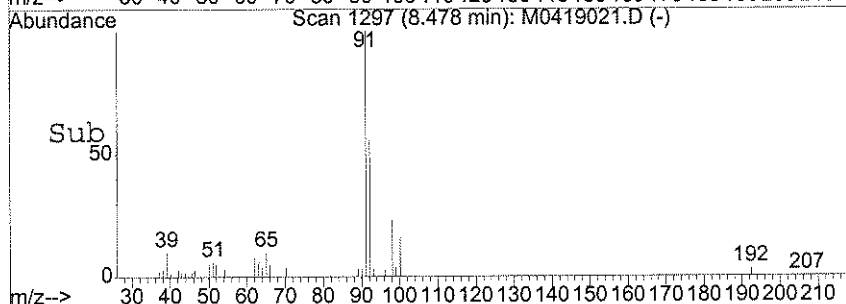
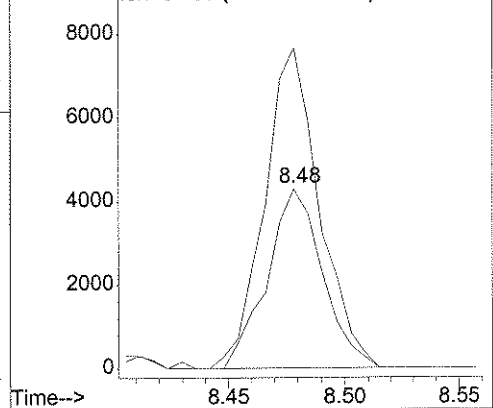


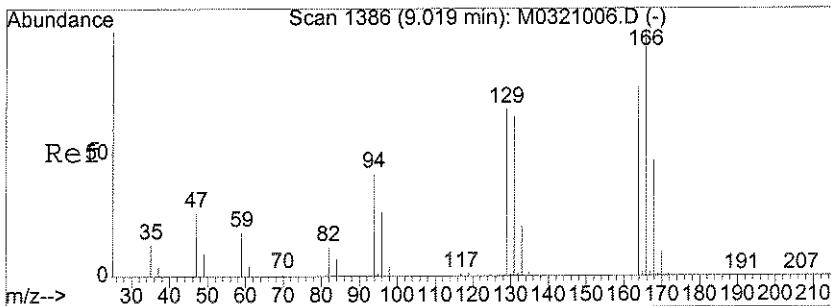
#52  
 Toluene  
 Concen: 0.32 ug/l  
 RT: 8.48 min Scan# 1297  
 Delta R.T. 0.00 min  
 Lab File: M0419021.D  
 Acq: 19 Apr 2007 20:24

Tgt Ion	Resp	Lower	Upper
92	100		
91	177.0	133.7	200.5



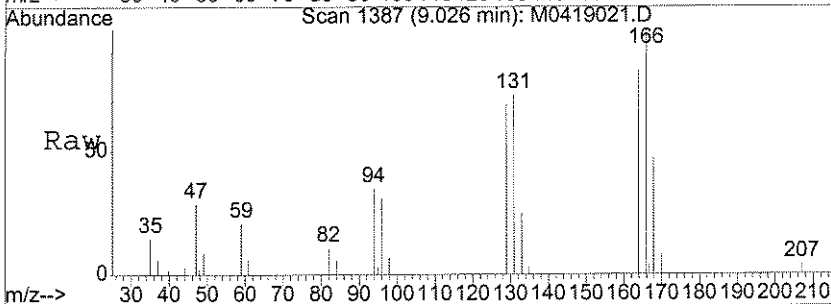
Abundance  
 Ion 92.05 (91.75 to 92.75): M0419021.  
 Ion 91.05 (90.75 to 91.75): M0419021.



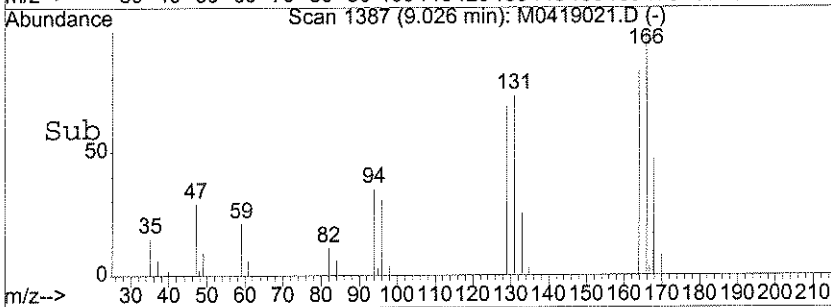
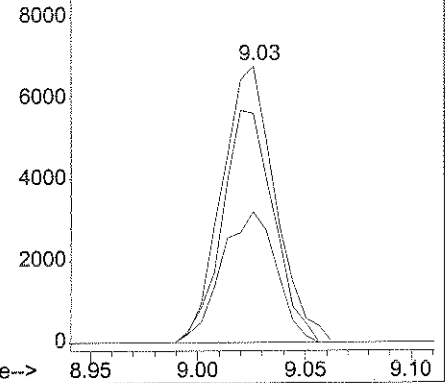


#56  
 Tetrachloroethene  
 Concen: 1.48 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0419021.D  
 Acq: 19 Apr 2007 20:24

Tgt Ion	Resp	Lower	Upper
166	11734		
166	100		
164	80.9	63.3	94.9
168	48.3	39.6	59.4



Abundance Ion 165.95 (165.65 to 166.65): M04190  
 Ion 163.95 (163.65 to 164.65): M04190  
 Ion 167.95 (167.65 to 168.65): M04190



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-16

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-006  
 Lab File ID: M0419022.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 20:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	1.2	
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	9.1	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	7.9	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.0	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	1.1	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-16

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-006

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419022.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/11/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 20:48

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	2.8	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-16

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-006  
 Lab File ID: M0419022.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 20:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

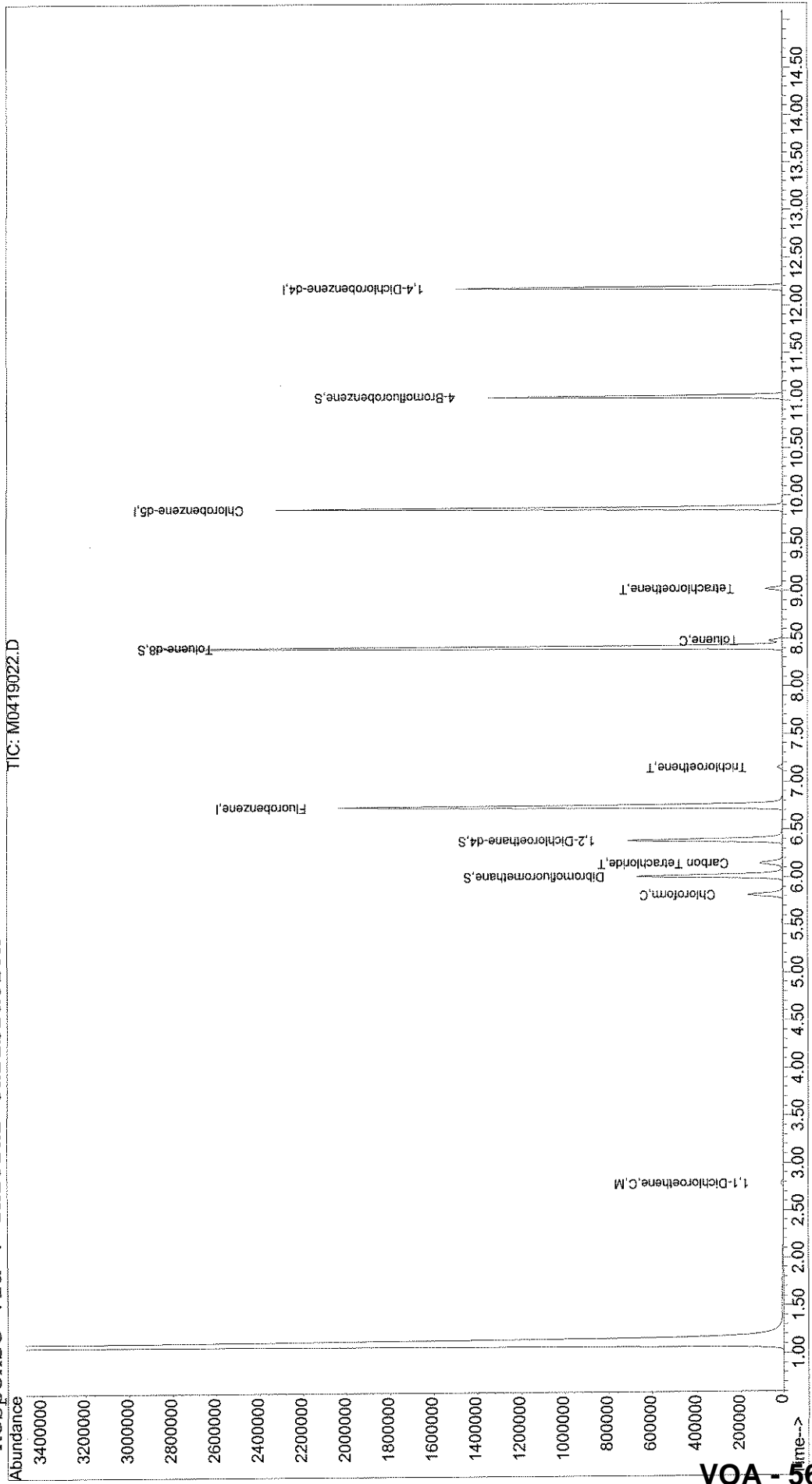
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419022.D  
Acq On : 19 Apr 2007 20:48  
Sample : JPL33-006  
Misc : 5ml +IS/SS #3  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:16 2007  
Vial: 65  
Operator: LH  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419022.D  
 Acq On : 19 Apr 2007 20:48  
 Sample : JPL33-006  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:16 2007

Vial: 65  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	1954436	50.00	ug/l	0.00 100.13%
50) Chlorobenzene-d5	9.87	82	661454	50.00	ug/l	0.00 94.05%
70) 1,4-Dichlorobenzene-d4	12.19	152	353818	50.00	ug/l	0.00 90.90%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	466023	50.54	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.08%
37) 1,2-Dichloroethane-d4	6.39	65	527243	49.95	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.90%
51) Toluene-d8	8.41	98	1915064	52.42	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.84%
72) 4-Bromofluorobenzene	11.04	95	399416	54.73	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.46%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.42	50	71	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.84	96	85	N.D.		
6) Chloroethane	1.90	64	57	N.D.		
7) Trichlorofluoromethane	2.18	101	632	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.79	96	6935	1.16	ug/l	95
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.98	43	300	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	360	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.88	96	63	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.55	63	1387	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0419022.D 8260B.M Fri Apr 20 10:16:54 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419022.D  
 Acq On : 19 Apr 2007 20:48  
 Sample : JPL33-006  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:16 2007

Vial: 65  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.56	43	75		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.59	128	74		N.D.	
30) Methacrylonitrile	5.80	41	281		N.D.	
31) Chloroform	5.82	83	142600	9.09	ug/l	96
32) 1,1,1-Trichloroethane	5.96	97	69		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.14	117	74033	7.94	ug/l	99
36) 1,1-Dichloropropene	6.17	75	60		N.D.	
38) Benzene	6.40	78	517		N.D.	
39) 1,2-Dichloroethane	6.47	62	263		N.D.	
40) Isobutanol	0.00	43	0		N.D.	
41) Trichloroethene	7.14	130	8872	1.01	ug/l	93
42) Methylcyclohexane	7.15	83	59		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.60	93	66		N.D.	
45) Methyl methacrylate	7.70	41	56		N.D.	
46) Bromodichloromethane	7.72	83	61		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	23779	1.07	ug/l	96
53) trans-1,3-Dichloropropene	8.90	75	134		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.93	97	480		N.D.	
56) Tetrachloroethene	9.02	166	21881	2.76	ug/l	98
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	274		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.98	91	271		N.D.	
65) m,p-Xylene	10.10	106	148		N.D.	
66) o-xylene	10.37	106	58		N.D.	
67) Styrene	0.00	104	0		N.D.	

*W 4/20/07*



Quantitation Report

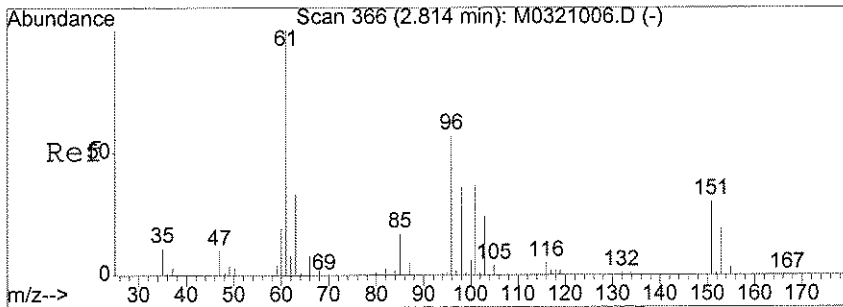
Data File : X:\MSVOA\MOBY\041907\M0419022.D  
 Acq On : 19 Apr 2007 20:48  
 Sample : JPL33-006  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:16 2007

Vial: 65  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

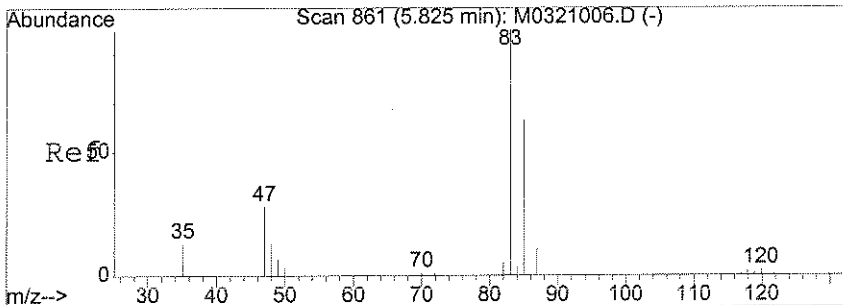
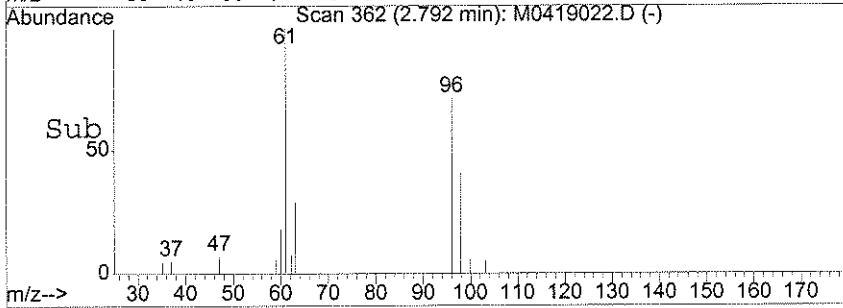
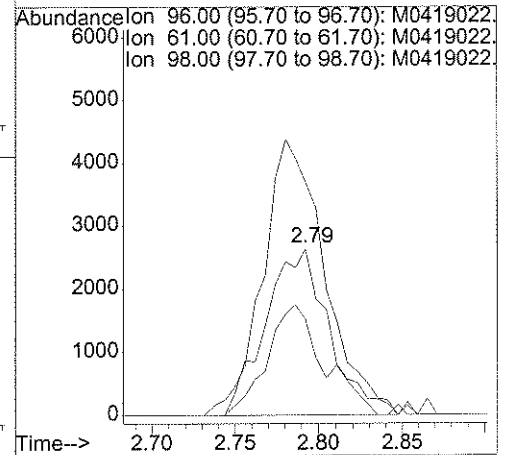
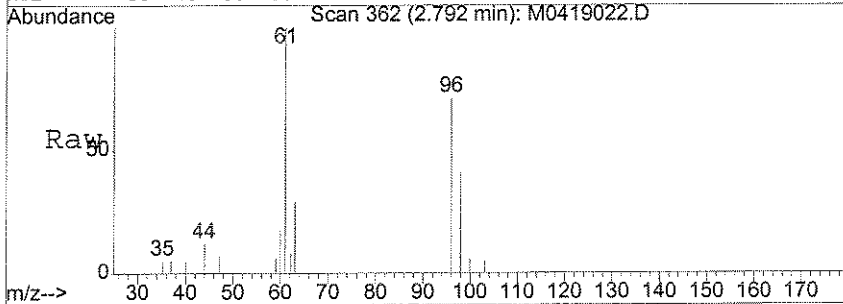
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	224		N.D.	
69) Isopropylbenzene	11.04	105	476		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	138		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	420		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	274		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.36	91	113		N.D.	
78) 4-Chlorotoluene	11.46	91	64		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.88	119	56		N.D.	
81) 1,2,4-Trimethylbenzene	11.97	105	80		N.D.	
82) sec-butylbenzene	11.97	105	80		N.D.	
83) 1,3-Dichlorobenzene	12.17	146	63		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.17	146	63		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.33	225	59		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



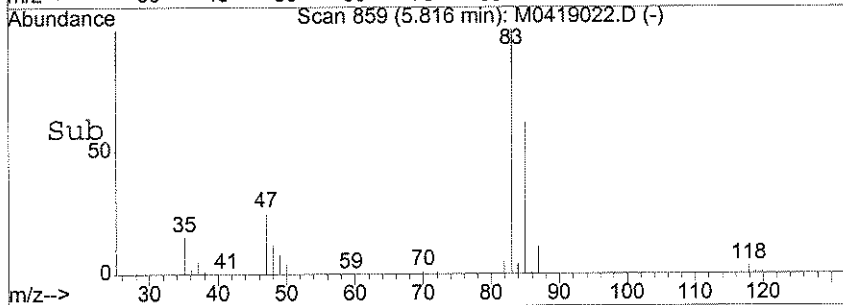
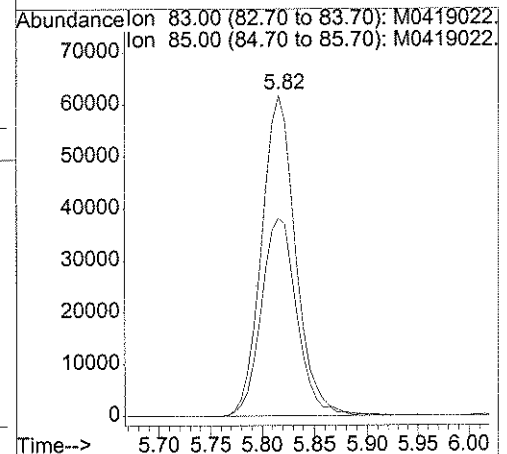
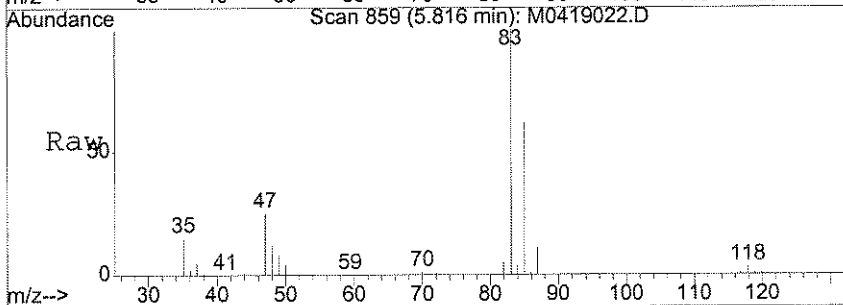
#9  
 1,1-Dichloroethene  
 Concen: 1.16 ug/l  
 RT: 2.79 min Scan# 362  
 Delta R.T. 0.01 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

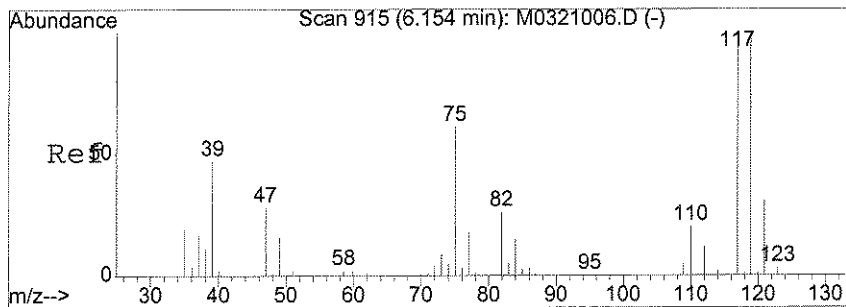
Tgt Ion	Resp	Lower	Upper
96	6935		
96	100		
61	162.3	138.6	178.6
98	59.4	32.2	72.2



#31  
 Chloroform  
 Concen: 9.09 ug/l  
 RT: 5.82 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

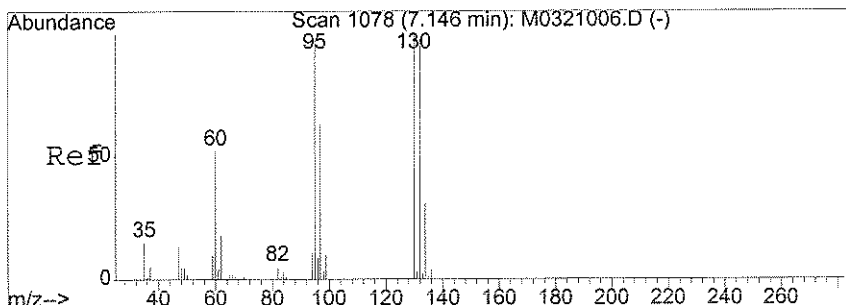
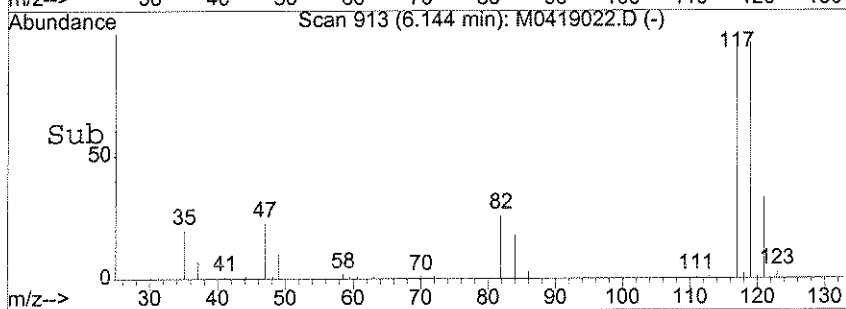
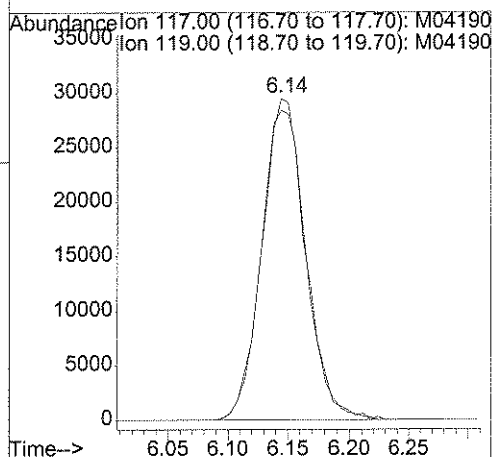
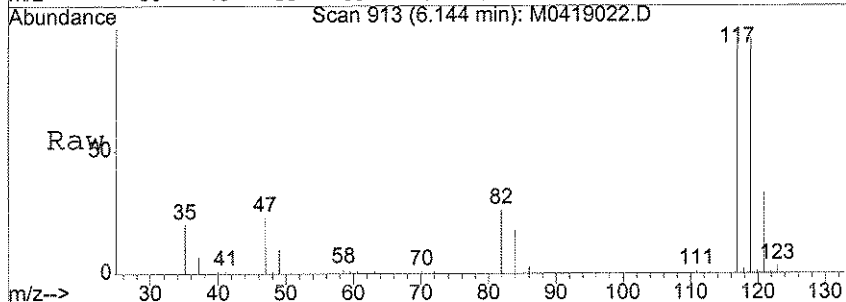
Tgt Ion	Resp	Lower	Upper
83	142600		
83	100		
85	64.1	41.2	81.2





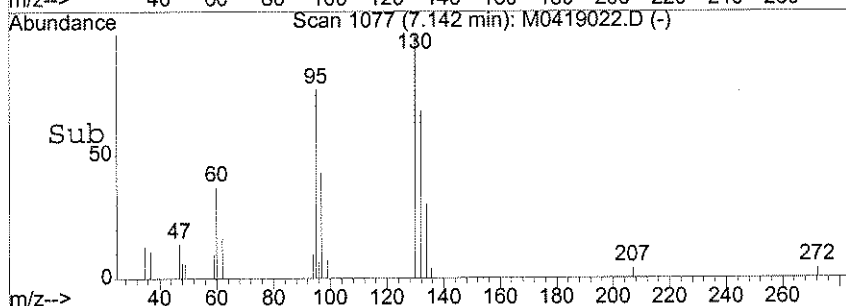
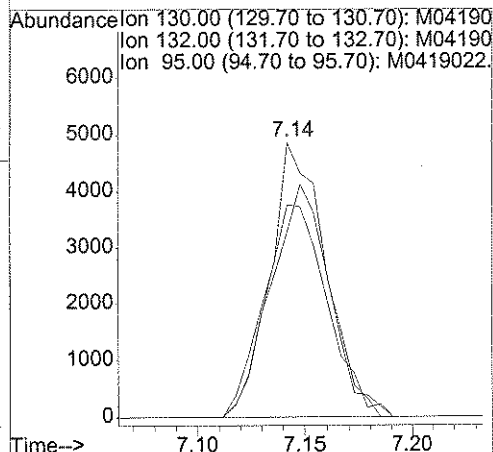
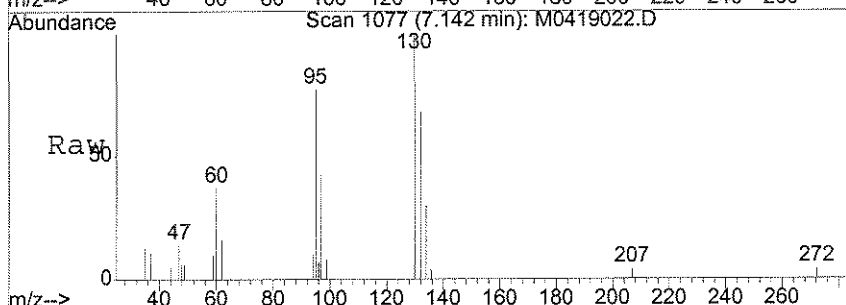
#35  
 Carbon Tetrachloride  
 Concen: 7.94 ug/l  
 RT: 6.14 min Scan# 913  
 Delta R.T. 0.00 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

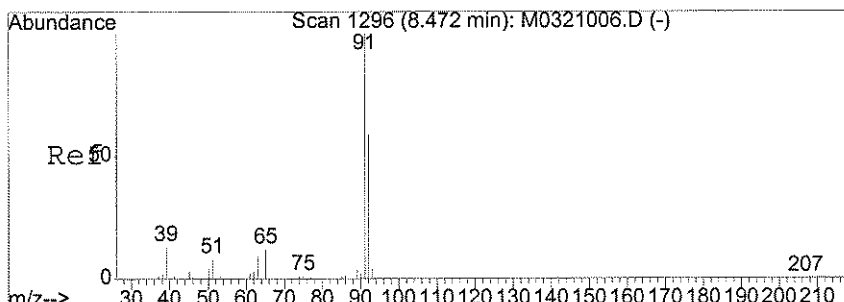
Tgt Ion:117 Resp: 74033  
 Ion Ratio Lower Upper  
 117 100  
 119 99.1 78.2 118.2



#41  
 Trichloroethene  
 Concen: 1.01 ug/l  
 RT: 7.14 min Scan# 1077  
 Delta R.T. -0.00 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

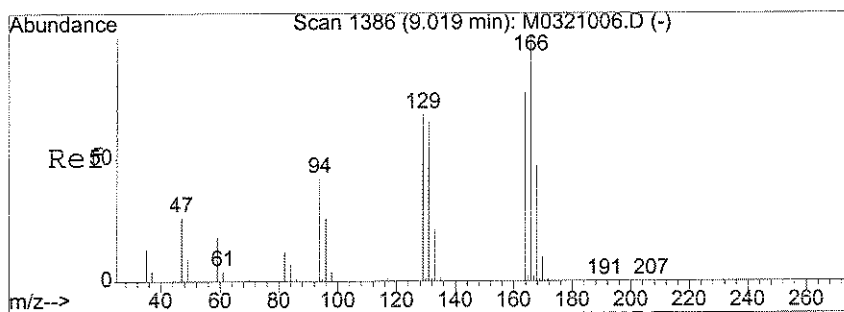
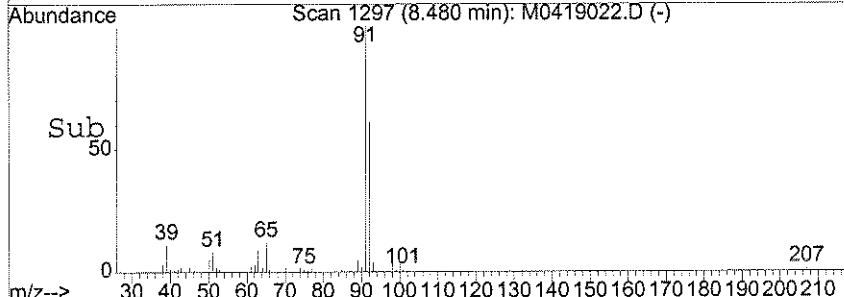
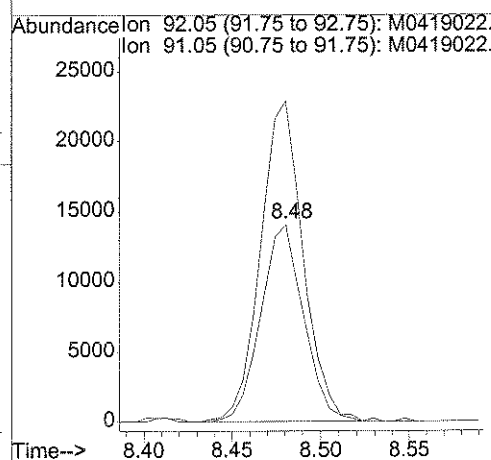
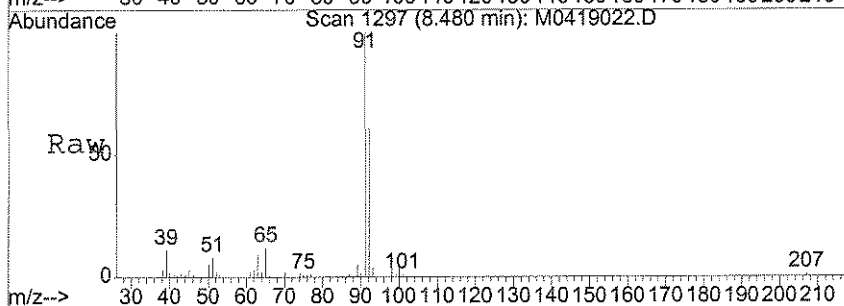
Tgt Ion:130 Resp: 8872  
 Ion Ratio Lower Upper  
 130 100  
 132 87.0 75.0 115.0  
 95 83.4 69.4 109.4





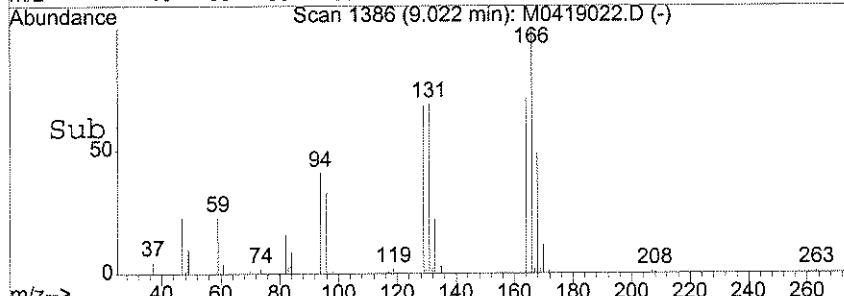
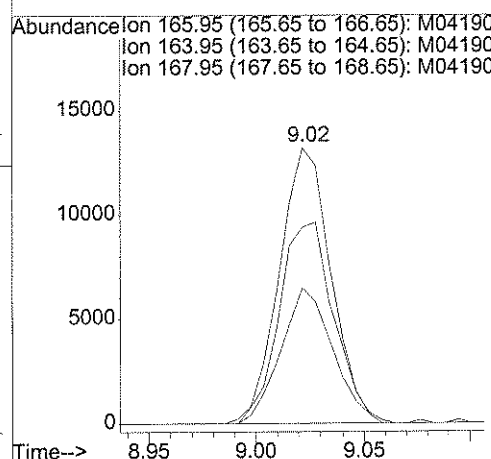
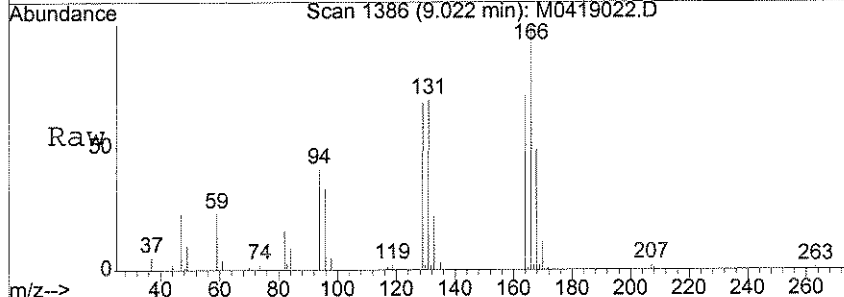
#52  
 Toluene  
 Concen: 1.07 ug/l  
 RT: 8.48 min Scan# 1297  
 Delta R.T. 0.00 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

Tgt Ion:	92	Resp:	23779
Ion Ratio	Lower	Upper	
92	100		
91	161.5	133.7	200.5



#56  
 Tetrachloroethene  
 Concen: 2.76 ug/l  
 RT: 9.02 min Scan# 1386  
 Delta R.T. -0.00 min  
 Lab File: M0419022.D  
 Acq: 19 Apr 2007 20:48

Tgt Ion:	166	Resp:	21881
Ion Ratio	Lower	Upper	
166	100		
164	77.0	63.3	94.9
168	49.2	39.6	59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-12-4/11/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-007  
 Lab File ID: M0419014.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 17:35  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-12-4/11/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-007  
 Lab File ID: M0419014.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 17:35  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-12-4/11/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-007  
 Lab File ID: M0419014.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 17:35  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

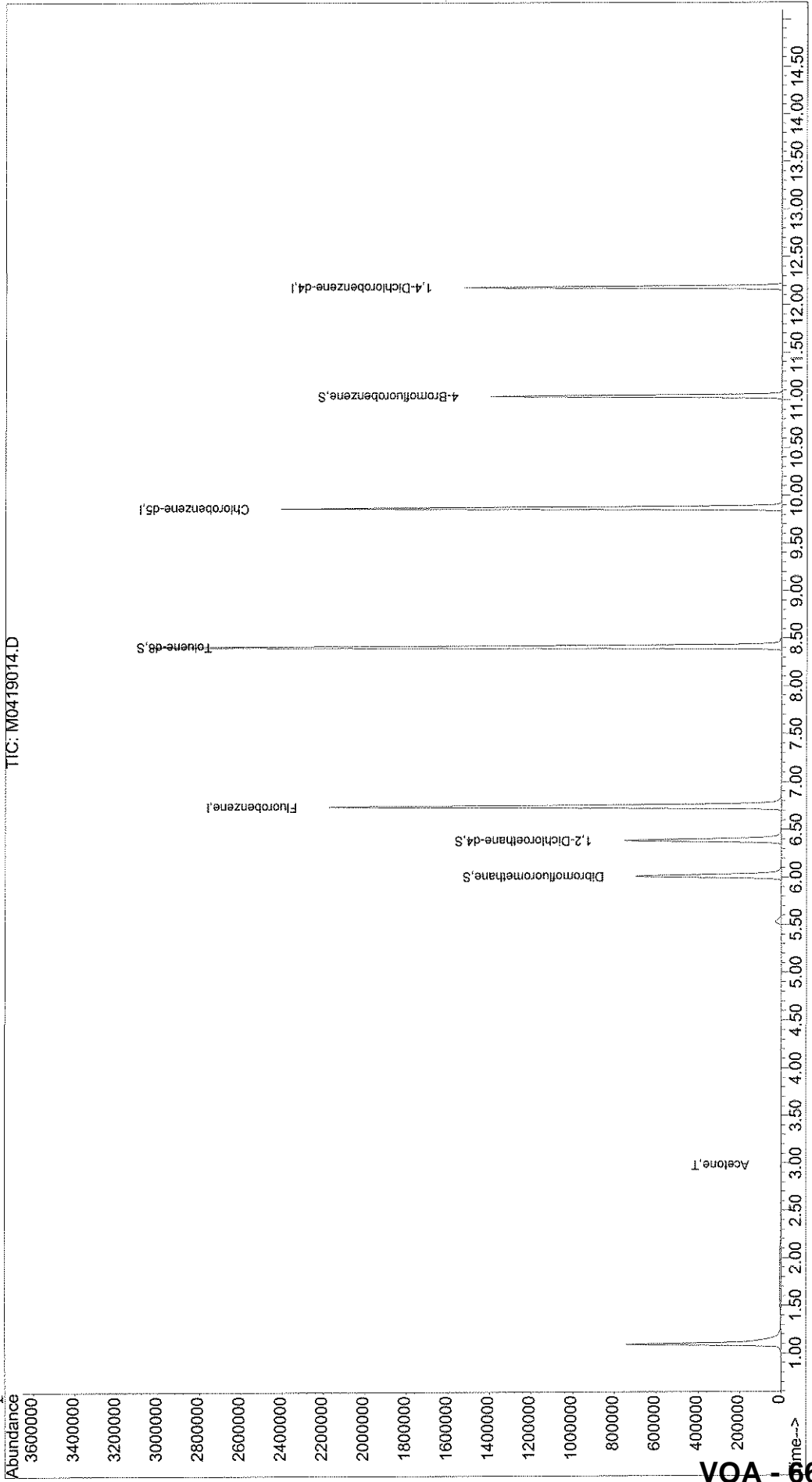
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419014.D Vial: 57  
Acq On : 19 Apr 2007 17:35 Operator: LH  
Sample : JPL33-007 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:02 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration





Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419014.D  
 Acq On : 19 Apr 2007 17:35  
 Sample : JPL33-007 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:02 2007

Vial: 57  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	2058498	50.00	ug/l	0.00 105.46%
50) Chlorobenzene-d5	9.87	82	701245	50.00	ug/l	0.00 99.71%
70) 1,4-Dichlorobenzene-d4	12.19	152	376296	50.00	ug/l	0.00 96.68%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	493525	50.82	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.64%
37) 1,2-Dichloroethane-d4	6.39	65	558488	50.24	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.48%
51) Toluene-d8	8.41	98	2030752	52.43	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.86%
72) 4-Bromofluorobenzene	11.04	95	430688	55.49	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.98%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	1.41	50	60		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	0.00	96	0		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	0.00	101	0		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	2.75	96	56		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D. d	
11) Acetone	2.97	43	3770	2.01	ug/l #	78
12) Iodomethane	2.90	142	55		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.04	76	757		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D. d	
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D. d	
19) trans-1,2-Dichloroethene	3.88	96	61		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	0.00	63	0		N.D.	

*Handwritten:* 78  
 LW 4/20/07

(#) = qualifier out of range (m) = manual integration  
 M0419014.D 8260B.M Fri Apr 20 10:03:06 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419014.D  
 Acq On : 19 Apr 2007 17:35  
 Sample : JPL33-007 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:02 2007

Vial: 57  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	4.67	53	58		N.D.	
25) 2,2-Dichloropropane	5.35	77	82		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	0.00	43	0		N.D.	d
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.69	41	76		N.D.	
31) Chloroform	5.83	83	266		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.16	75	387		N.D.	
38) Benzene	6.40	78	1067		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.14	130	134		N.D.	
42) Methylcyclohexane	7.29	83	192		N.D.	
43) 1,2-Dichloropropane	7.28	63	65		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.27	75	66		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	1690		N.D.	
53) trans-1,3-Dichloropropene	8.72	75	74		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.03	166	230		N.D.	
57) 1,3-Dichloropropane	9.13	76	87		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	9.51	107	88		N.D.	
61) Chlorobenzene	9.90	112	334		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	10.03	131	59		N.D.	
64) Ethylbenzene	10.00	91	893		N.D.	
65) m,p-Xylene	10.10	106	876		N.D.	
66) o-xylene	10.51	106	248		N.D.	
67) Styrene	10.52	104	186		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419014.D  
 Acq On : 19 Apr 2007 17:35  
 Sample : JPL33-007 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:02 2007

Vial: 57  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.76	173	349		N.D.	
69) Isopropylbenzene	10.87	105	408		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	412		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	283		N.D.	
76) n-Propylbenzene	11.27	120	166		N.D.	
77) 2-Chlorotoluene	11.27	91	687		N.D.	
78) 4-Chlorotoluene	11.48	91	408		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	307		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	435		N.D.	
82) sec-butylbenzene	11.98	105	413		N.D.	
83) 1,3-Dichlorobenzene	12.11	146	203		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	150		N.D.	
86) 1,2-Dichlorobenzene	12.57	146	58		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	13.54	157	56		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.33	225	357		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-7-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-008  
 Lab File ID: M0419023.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 21:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.84	
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	9.2	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	8.0	
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.94	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.98	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-7-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-008  
 Lab File ID: M0419023.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 21:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	2.7	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

DUPE-7-1Q07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-008  
 Lab File ID: M0419023.D  
 Date Collected: 04/11/2007  
 Date/Time Analyzed: 04/19/2007 21:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

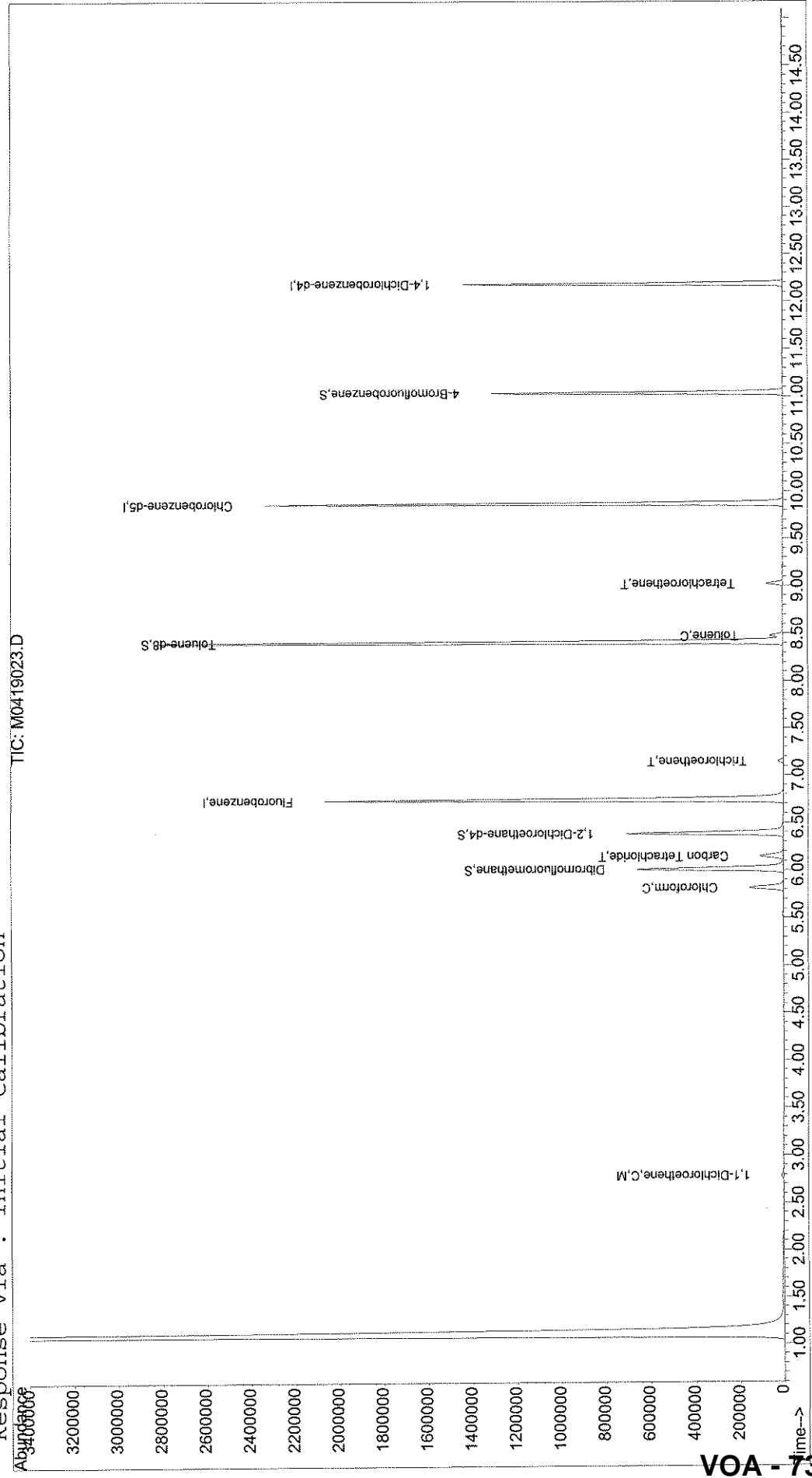
CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419023.D  
Acq On : 19 Apr 2007 21:12  
Sample : JPL33-008  
Misc : 5ml +IS/SS #4  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:18 2007  
Vial: 66  
Operator: LH  
Inst : MOBY  
Multiplr: 1.00  
Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 73

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419023.D  
 Acq On : 19 Apr 2007 21:12  
 Sample : JPL33-008  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:18 2007

Vial: 66  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.74	96	1943605	50.00	ug/l	0.00 99.57%
50) Chlorobenzene-d5	9.87	82	661314	50.00	ug/l	0.00 94.03%
70) 1,4-Dichlorobenzene-d4	12.19	152	353680	50.00	ug/l	0.00 90.87%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	460435	50.21	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.42%
37) 1,2-Dichloroethane-d4	6.39	65	521850	49.72	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.44%
51) Toluene-d8	8.41	98	1916455	52.47	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.94%
72) 4-Bromofluorobenzene	11.04	95	400959	54.96	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.92%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.79	96	88	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	2.20	101	726	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.79	96	4985	0.84 ug/l		94
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d	
11) Acetone	2.93	43	131	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.06	76	58	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	3.34	43	56	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.53	63	1146	N.D.		

*MS 4/20/07*

(#) = qualifier out of range (m) = manual integration  
 M0419023.D 8260B.M Fri Apr 20 10:18:34 2007



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419023.D  
 Acq On : 19 Apr 2007 21:12  
 Sample : JPL33-008  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:18 2007

Vial: 66  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.48	43	122		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.82	41	388		N.D.	
31) Chloroform	5.82	83	142784	9.15	ug/l	94
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	6.00	56	120		N.D.	
35) Carbon Tetrachloride	6.15	117	73925	7.97	ug/l	99
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	1180		N.D.	
39) 1,2-Dichloroethane	6.48	62	422		N.D.	
40) Isobutanol	0.00	43	0		N.D. d	
41) Trichloroethene	7.15	130	8180	0.94	ug/l	96
42) Methylcyclohexane	7.15	83	165		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.58	83	55		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.12	75	58		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	21634	0.98	ug/l	89
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.94	97	539		N.D.	
56) Tetrachloroethene	9.03	166	21488	2.71	ug/l	98
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	282		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	298		N.D.	
65) m,p-Xylene	10.11	106	295		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.76	104	77		N.D.	

LW 4/20/07

(#) = qualifier out of range (m) = manual integration  
 M0419023.D 8260B.M Fri Apr 20 10:18:34 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419023.D  
 Acq On : 19 Apr 2007 21:12  
 Sample : JPL33-008  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:18 2007

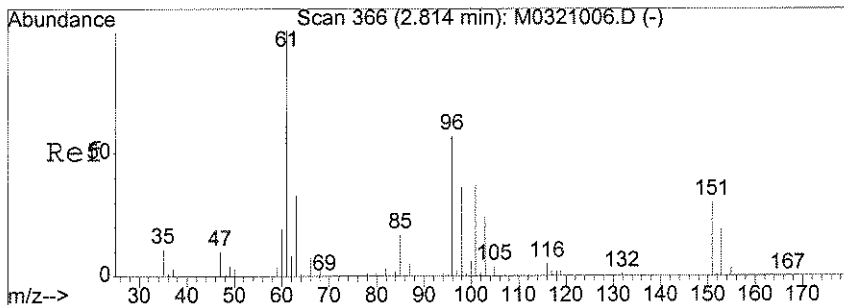
Vial: 66  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

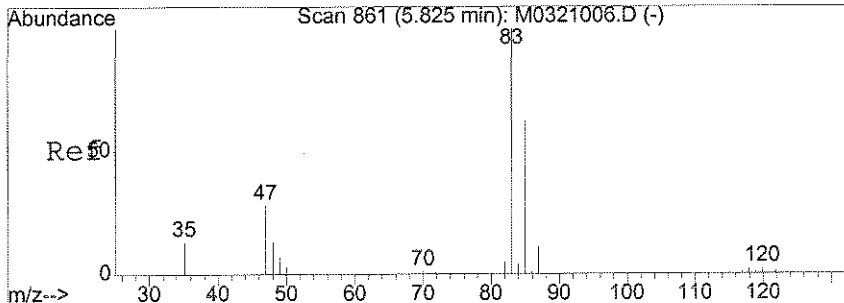
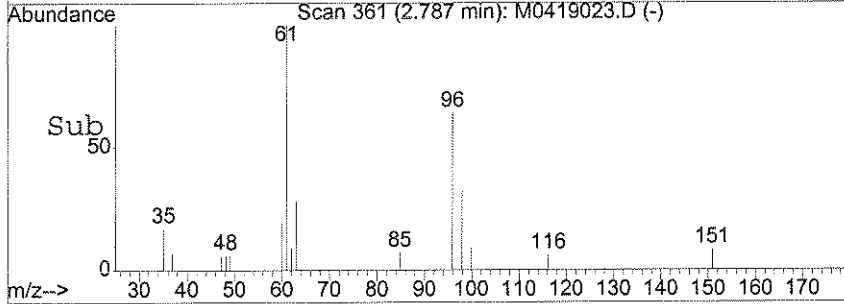
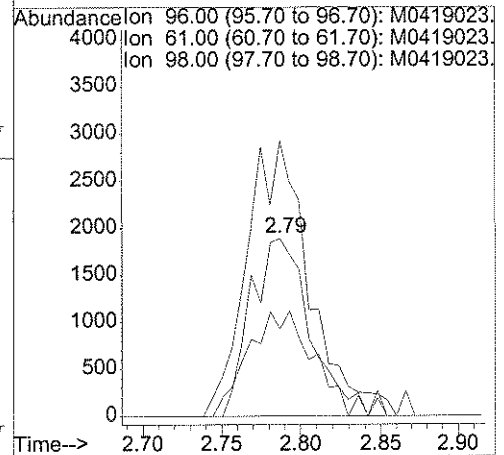
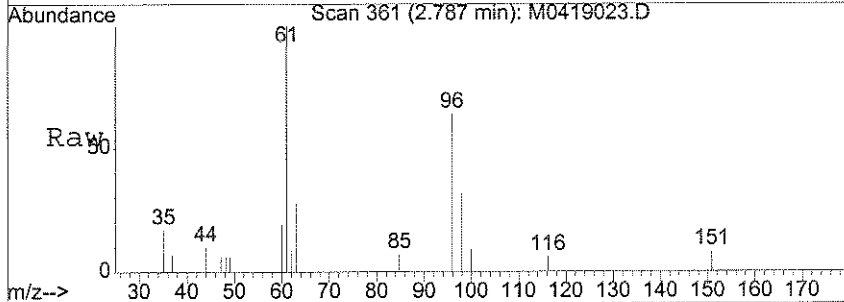
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	154		N.D.	
69) Isopropylbenzene	10.86	105	105		N.D.	
71) trans-1,4-Dichloro-2-buten	11.13	53	55		N.D.	
73) Bromobenzene	11.04	156	84		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.10	83	59		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	148		N.D.	
76) n-Propylbenzene	11.44	120	62		N.D.	
77) 2-Chlorotoluene	11.36	91	66		N.D.	
78) 4-Chlorotoluene	11.36	91	66		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	67		N.D.	
80) tert-Butylbenzene	11.88	119	55		N.D.	
81) 1,2,4-Trimethylbenzene	11.98	105	87		N.D.	
82) sec-butylbenzene	11.98	105	87		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	61		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.20	146	57		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

(#) = qualifier out of range (m) = manual integration  
 M0419023.D 8260B.M Fri Apr 20 10:18:35 2007



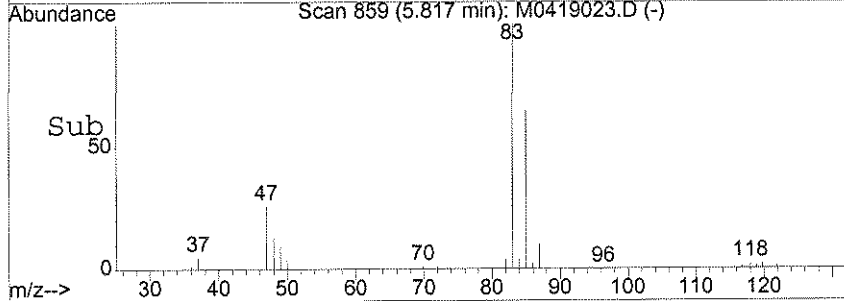
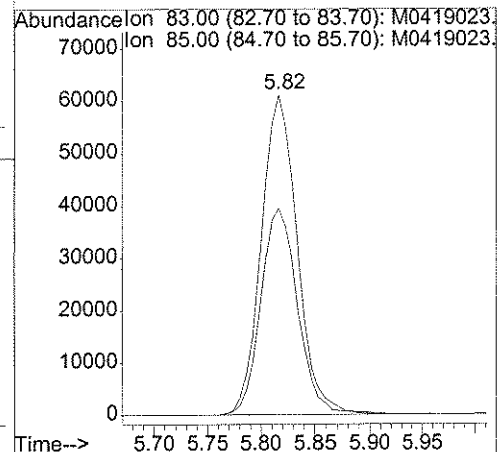
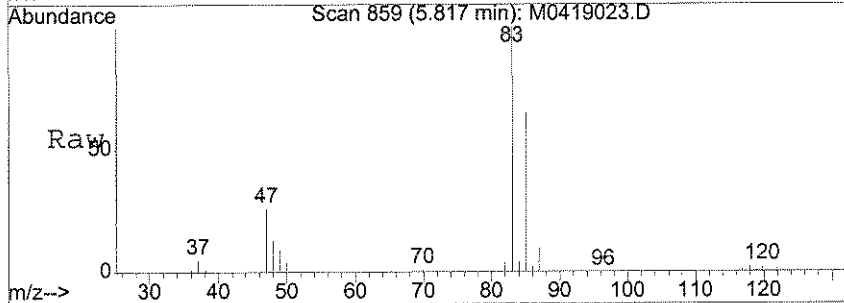
#9  
 1,1-Dichloroethene  
 Concen: 0.84 ug/l  
 RT: 2.79 min Scan# 361  
 Delta R.T. 0.01 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

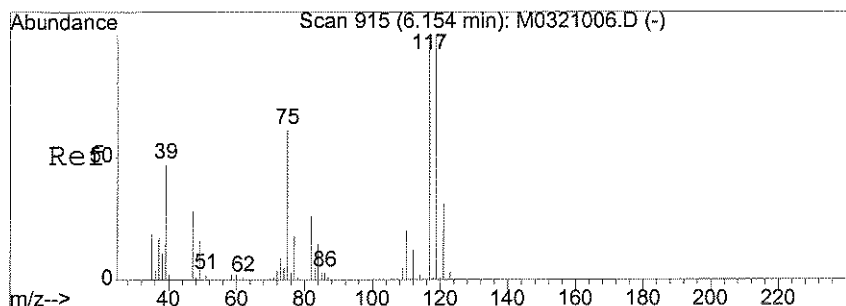
Tgt Ion	Resp	Lower	Upper
96	4985		
61	161.8	138.6	178.6
98	65.2	32.2	72.2



#31  
 Chloroform  
 Concen: 9.15 ug/l  
 RT: 5.82 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

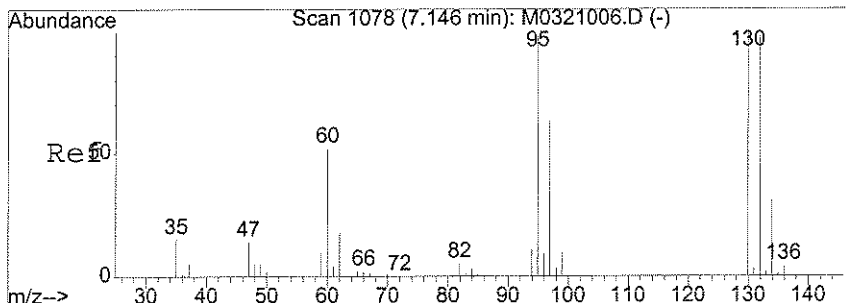
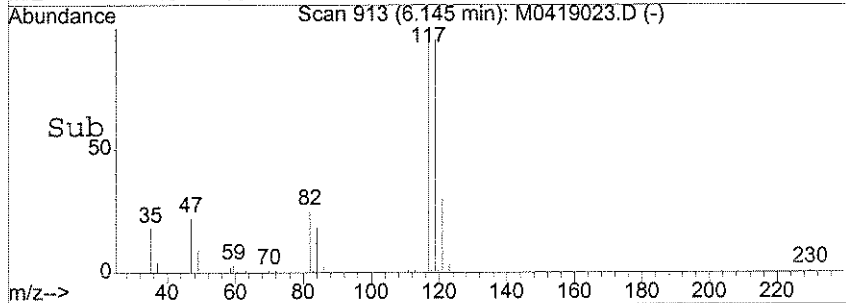
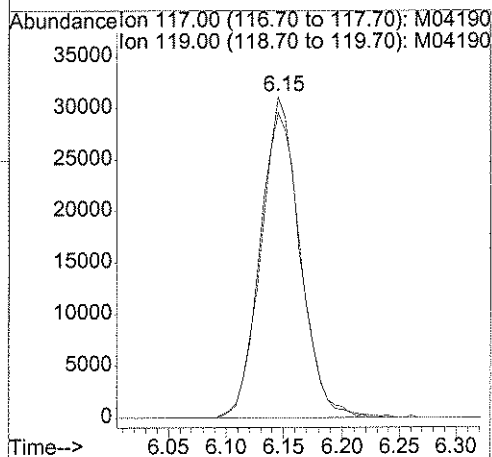
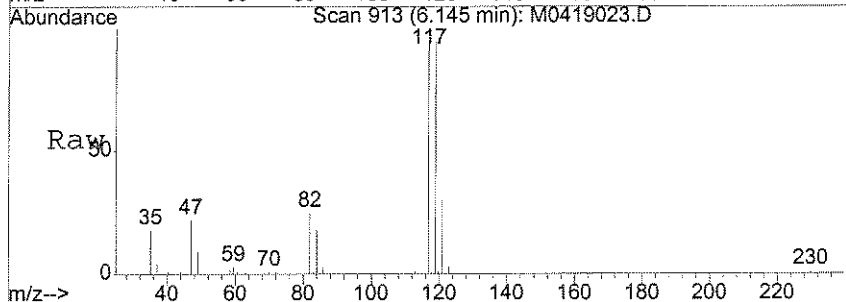
Tgt Ion	Resp	Lower	Upper
83	142784		
85	65.9	41.2	81.2





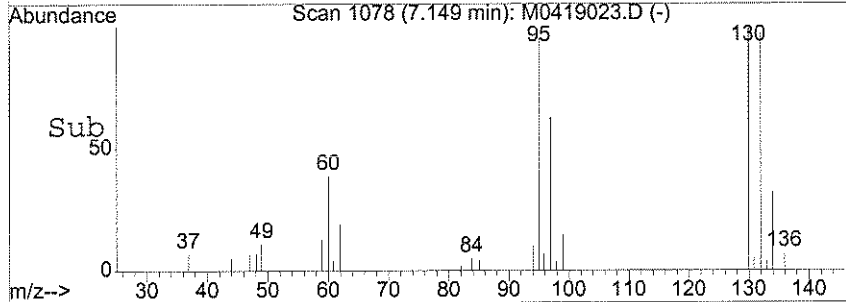
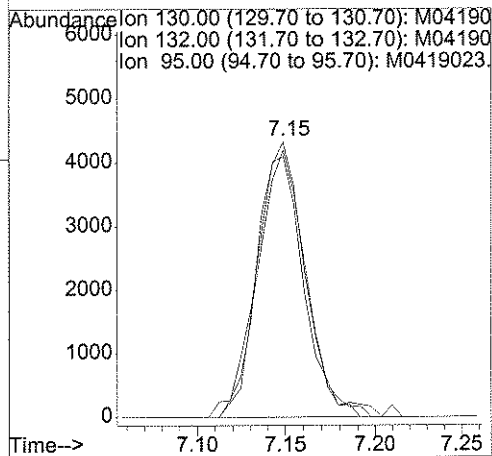
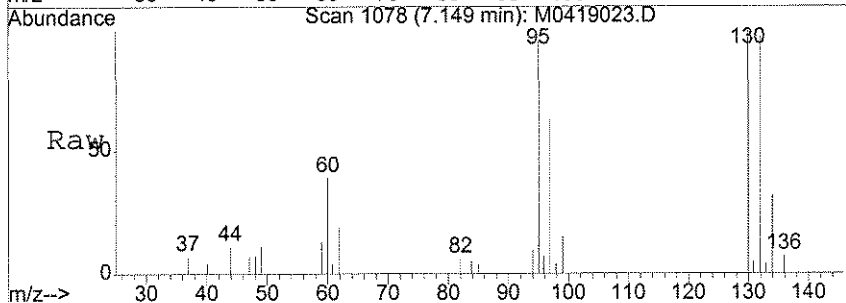
#35  
 Carbon Tetrachloride  
 Concen: 7.97 ug/l  
 RT: 6.15 min Scan# 913  
 Delta R.T. 0.00 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

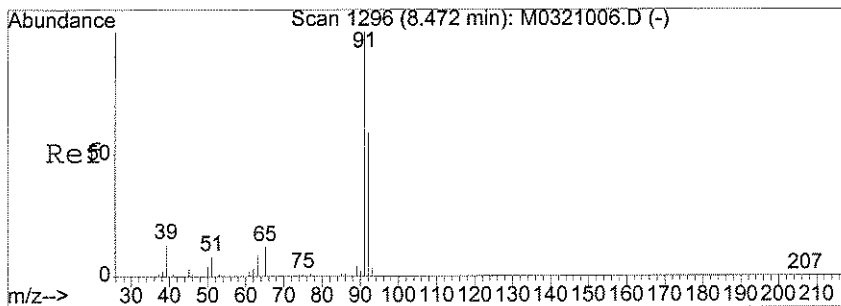
Tgt Ion:117 Resp: 73925  
 Ion Ratio Lower Upper  
 117 100  
 119 97.4 78.2 118.2



#41  
 Trichloroethene  
 Concen: 0.94 ug/l  
 RT: 7.15 min Scan# 1078  
 Delta R.T. 0.00 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

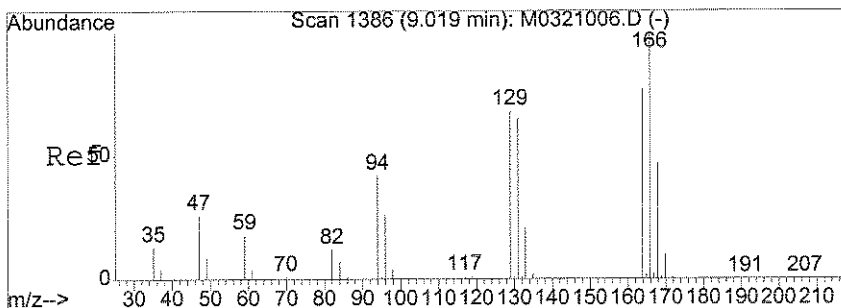
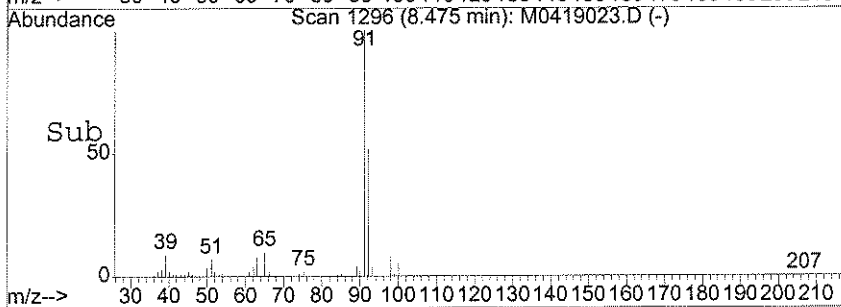
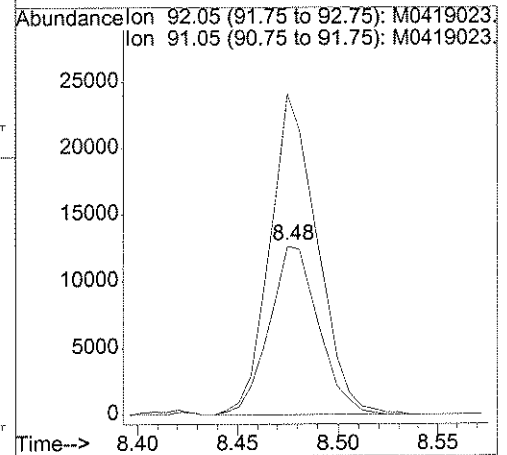
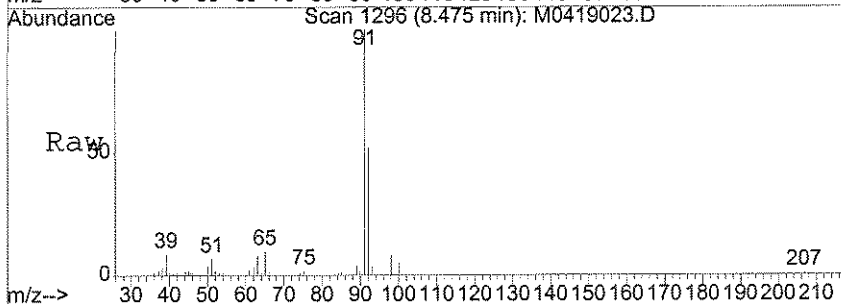
Tgt Ion:130 Resp: 8180  
 Ion Ratio Lower Upper  
 130 100  
 132 95.6 75.0 115.0  
 95 95.6 69.4 109.4





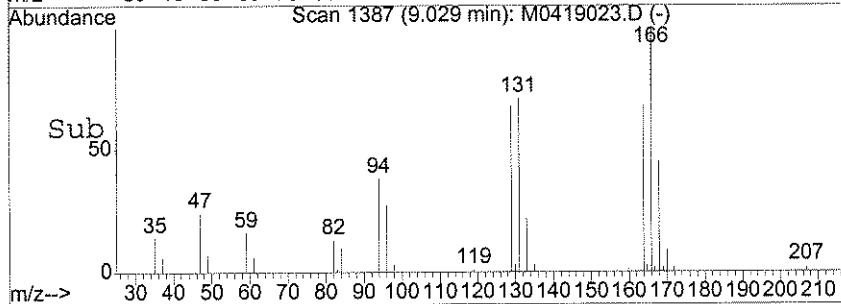
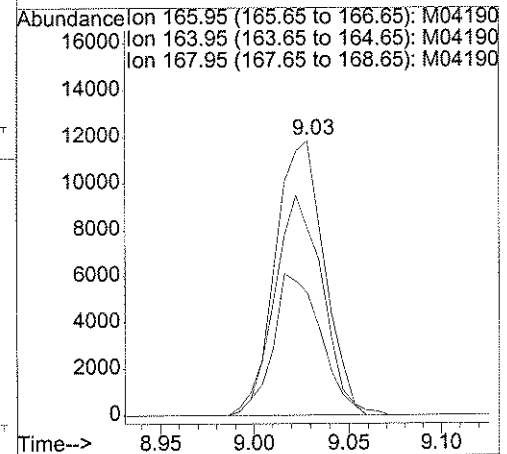
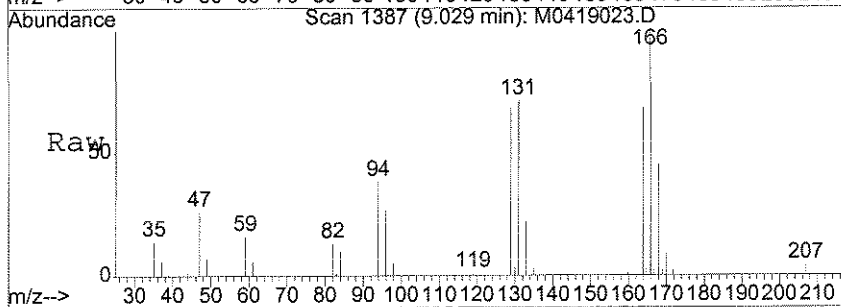
#52  
 Toluene  
 Concen: 0.98 ug/l  
 RT: 8.48 min Scan# 1296  
 Delta R.T. -0.00 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

Tgt Ion: 92 Resp: 21634  
 Ion Ratio Lower Upper  
 92 100  
 91 182.1 133.7 200.5



#56  
 Tetrachloroethene  
 Concen: 2.71 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0419023.D  
 Acq: 19 Apr 2007 21:12

Tgt Ion: 166 Resp: 21488  
 Ion Ratio Lower Upper  
 166 100  
 164 76.5 63.3 94.9  
 168 49.6 39.6 59.4



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-7

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-009  
 Lab File ID: M0419024.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 21:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	2.5	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	5.4	
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.56	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-7

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-009  
 Lab File ID: M0419024.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 21:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:		Q
		(ug/L or ug/kg)	ug/L	
10061-02-	trans-1,3-Dichloropropene	0.50		U
79-00-5	1,1,2-Trichloroethane	0.50		U
127-18-4	Tetrachloroethene	0.50		U
142-28-9	1,3-Dichloropropane	0.50		U
124-48-1	Dibromochloromethane	7.6		
106-93-4	1,2-Dibromoethane	0.50		U
108-90-7	Chlorobenzene	0.50		U
100-41-4	Ethylbenzene	0.50		U
630-20-6	1,1,1,2-Tetrachloroethane	0.50		U
179601-23	m,p-Xylene	1.0		U
95-47-6	o-Xylene	0.50		U
100-42-5	Styrene	0.50		U
75-25-2	Bromoform	5.7		
98-82-8	Isopropylbenzene	0.50		U
79-34-5	1,1,2,2-Tetrachloroethane	0.50		U
103-65-1	n-Propylbenzene	0.50		U
108-86-1	Bromobenzene	0.50		U
96-18-4	1,2,3-Trichloropropane	0.50		U
95-49-8	2-Chlorotoluene	0.50		U
108-67-8	1,3,5-Trimethylbenzene	0.50		U
106-43-4	4-Chlorotoluene	0.50		U
98-06-6	tert-Butylbenzene	0.50		U
95-63-6	1,2,4-Trimethylbenzene	0.50		U
135-98-8	sec-Butylbenzene	0.50		U
99-87-6	4-Isopropyltoluene	0.50		U
541-73-1	1,3-Dichlorobenzene	0.50		U
106-46-7	1,4-Dichlorobenzene	0.50		U
104-51-8	n-Butylbenzene	0.50		U
95-50-1	1,2-Dichlorobenzene	0.50		U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-7

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-009  
 Lab File ID: M0419024.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 21:36  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS:	
		(ug/L or ug/kg)	ug/L
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

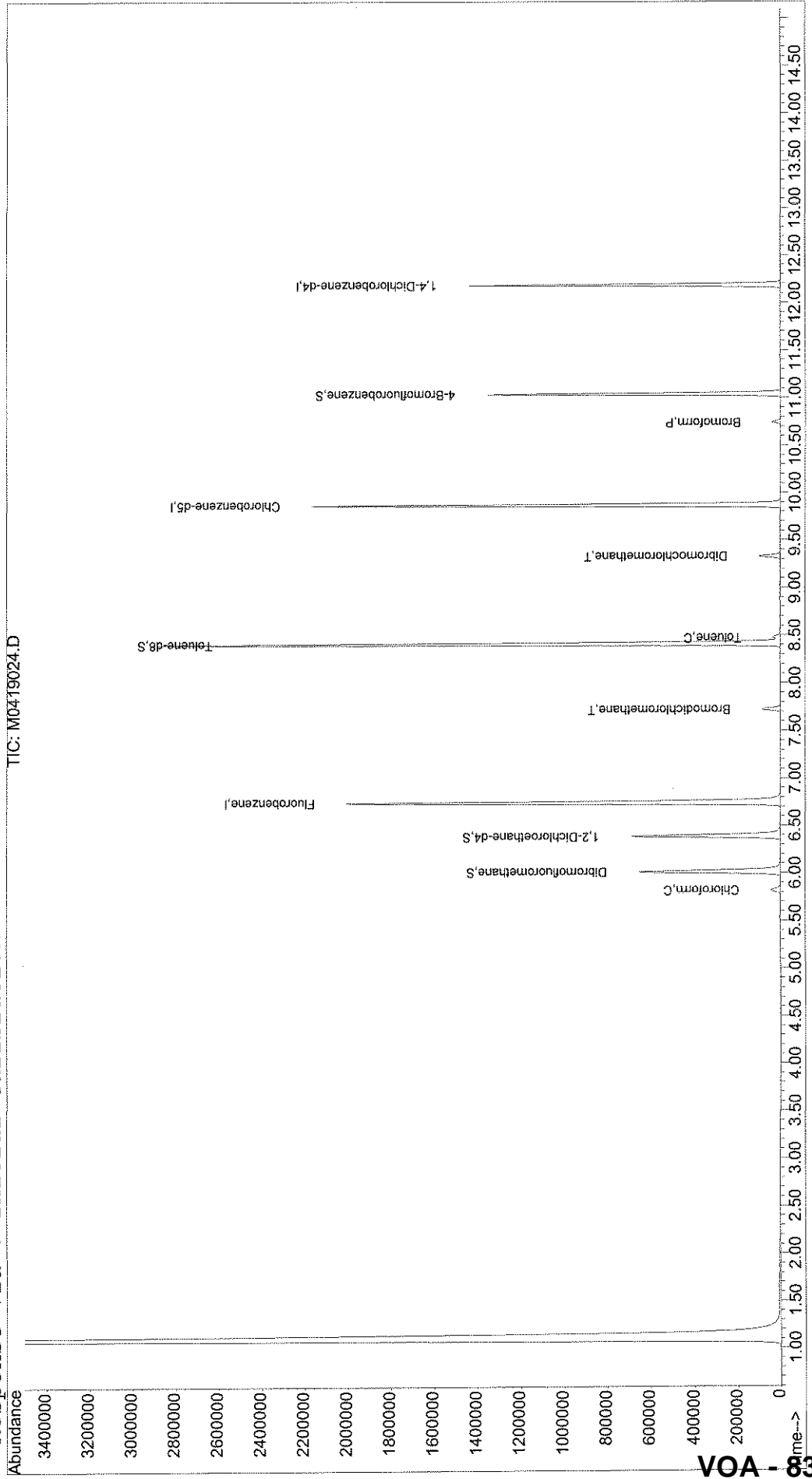
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419024.D Vial: 67  
Acq On : 19 Apr 2007 21:36 Operator: LH  
Sample : JPL33-009 Inst : MOBY  
Misc : 5ml +IS/SS #4 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:19 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 83

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419024.D  
 Acq On : 19 Apr 2007 21:36  
 Sample : JPL33-009  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:19 2007

Vial: 67  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
						Rcv (Ar )
1) Fluorobenzene	6.74	96	1932511	50.00	ug/l	0.00 99.01%
50) Chlorobenzene-d5	9.87	82	656618	50.00	ug/l	0.00 93.36%
70) 1,4-Dichlorobenzene-d4	12.18	152	347748	50.00	ug/l	0.00 89.34%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	455607	49.97	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	99.94%
37) 1,2-Dichloroethane-d4	6.39	65	515546	49.40	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	98.80%
51) Toluene-d8	8.41	98	1902500	52.46	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.92%
72) 4-Bromofluorobenzene	11.04	95	395198	55.10	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.20%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.42	50	81	N.D.		
4) Vinyl Chloride	1.52	62	55	N.D.		
5) Bromomethane	1.79	96	98	N.D.		
6) Chloroethane	1.95	64	55	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.77	96	58	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.98	43	55	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	338	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration  
 M0419024.D 8260B.M Fri Apr 20 10:20:08 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419024.D  
 Acq On : 19 Apr 2007 21:36  
 Sample : JPL33-009  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:19 2007

Vial: 67  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.54	43	57		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	5.69	128	77		N.D.	
30) Methacrylonitrile	5.81	41	58		N.D.	
31) Chloroform	5.81	83	39468	2.54	ug/l	96
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	0.00	56	0		N.D.	
35) Carbon Tetrachloride	6.14	117	285		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	1113		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D. d	
41) Trichloroethene	7.14	130	66		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.55	93	823		N.D.	
45) Methyl methacrylate	7.42	41	55		N.D.	
46) Bromodichloromethane	7.72	83	56494	5.35	ug/l	99
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	12409	0.56	ug/l	93
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.93	97	56		N.D.	
56) Tetrachloroethene	9.03	166	201		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	9.32	129	48117	7.57	ug/l	98
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.91	112	62		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	10.00	91	233		N.D.	
65) m,p-Xylene	10.10	106	149		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.53	104	70		N.D.	

LW 4/20/07

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419024.D  
 Acq On : 19 Apr 2007 21:36  
 Sample : JPL33-009  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:19 2007

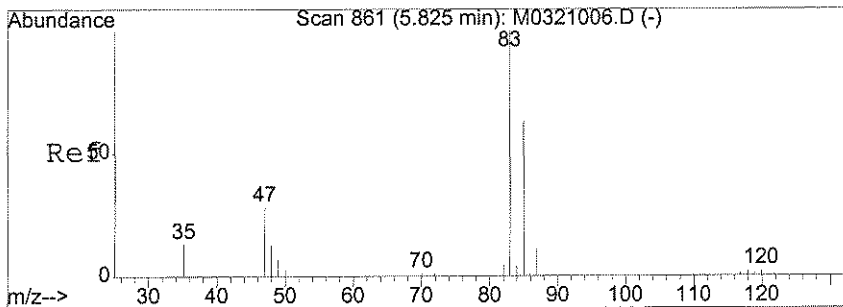
Vial: 67  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

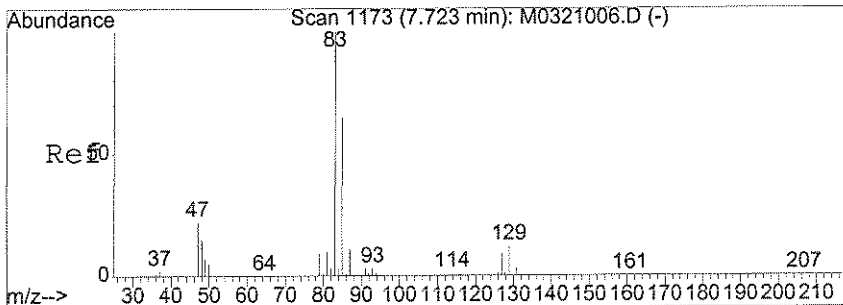
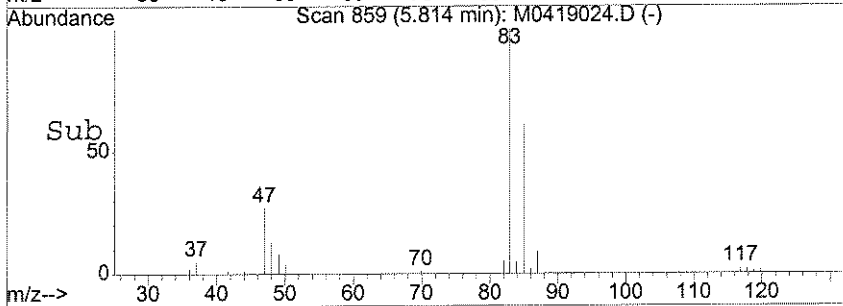
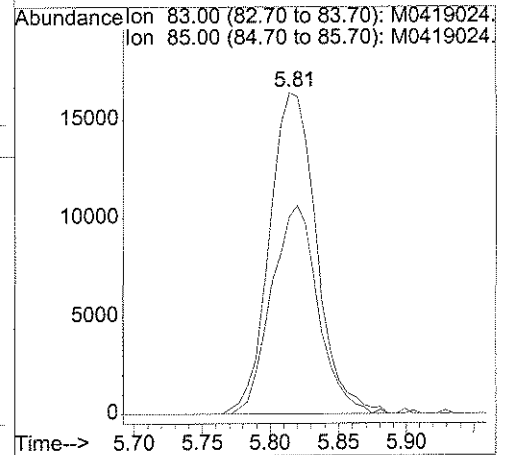
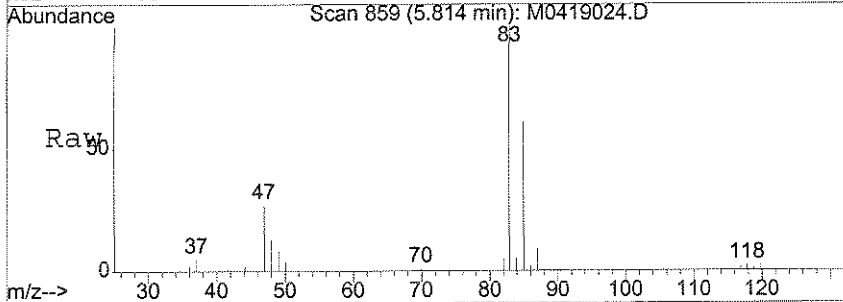
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	17870	5.66	ug/l	94
69) Isopropylbenzene	11.03	105	861	N.D.		
71) trans-1,4-Dichloro-2-buten	0.00	53	0	N.D.		
73) Bromobenzene	11.04	156	55	N.D.		
74) 1,1,2,2-Tetrachloroethane	11.03	83	215	N.D.		
75) 1,2,3-Trichloropropane	11.30	110	62	N.D.		
76) n-Propylbenzene	11.44	120	55	N.D.		
77) 2-Chlorotoluene	11.36	91	58	N.D.		
78) 4-Chlorotoluene	11.36	91	58	N.D.		
79) 1,3,5-Trimethylbenzene	0.00	105	0	N.D.		
80) tert-Butylbenzene	11.76	119	56	N.D.		
81) 1,2,4-Trimethylbenzene	11.82	105	160	N.D.		
82) sec-butylbenzene	11.98	105	77	N.D.		
83) 1,3-Dichlorobenzene	12.12	146	78	N.D.		
84) 4-Isopropyltoluene	0.00	119	0	N.D.		
85) 1,4-Dichlorobenzene	12.12	146	78	N.D.		
86) 1,2-Dichlorobenzene	0.00	146	0	N.D.		
87) n-Butylbenzene	0.00	91	0	N.D.	d	
88) 1,2-Dibromo-3-chloropropan	13.21	157	69	N.D.		
89) 1,2,4-Trichlorobenzene	0.00	180	0	N.D.		
90) Hexachlorobutadiene	14.33	225	71	N.D.		
91) Naphthalene	0.00	128	0	N.D.	d	
92) 1,2,3-Trichlorobenzene	0.00	180	0	N.D.	d	

LU 4/20/07



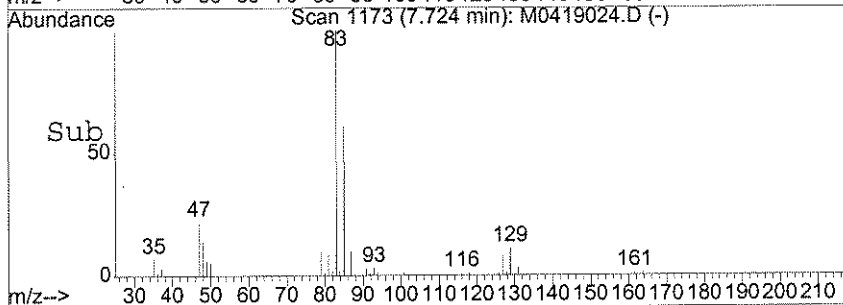
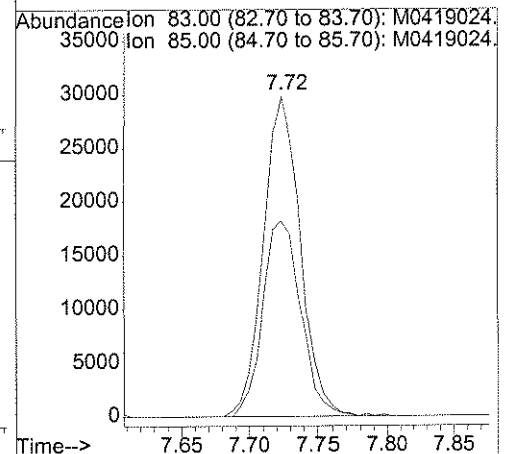
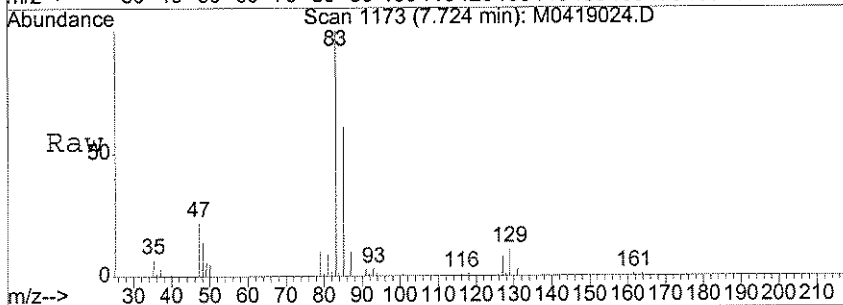
#31  
 Chloroform  
 Concen: 2.54 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419024.D  
 Acq: 19 Apr 2007 21:36

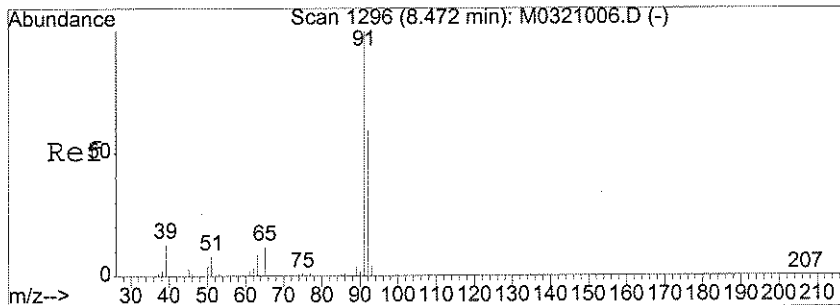
Tgt Ion: 83 Resp: 39468  
 Ion Ratio Lower Upper  
 83 100  
 85 64.5 41.2 81.2



#46  
 Bromodichloromethane  
 Concen: 5.35 ug/l  
 RT: 7.72 min Scan# 1173  
 Delta R.T. 0.00 min  
 Lab File: M0419024.D  
 Acq: 19 Apr 2007 21:36

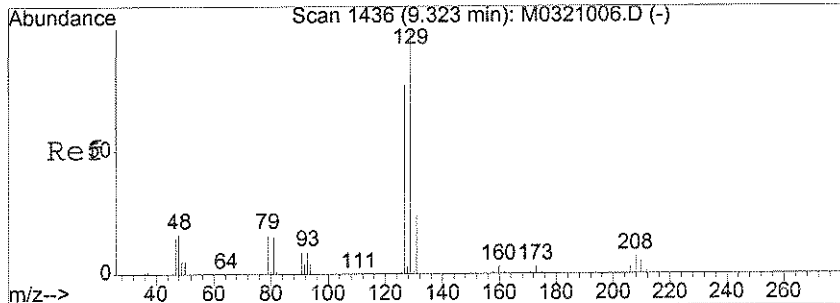
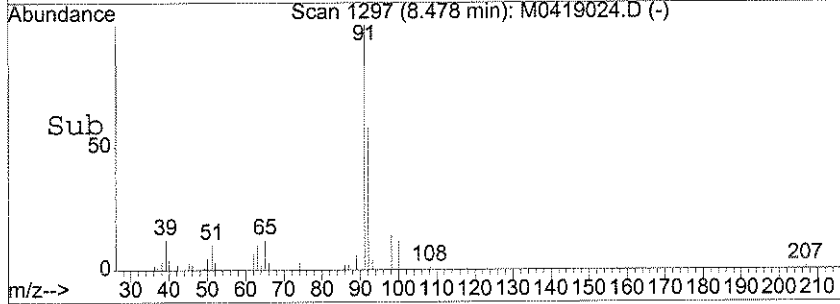
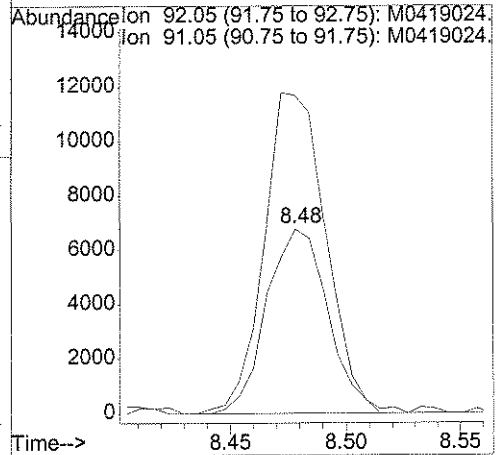
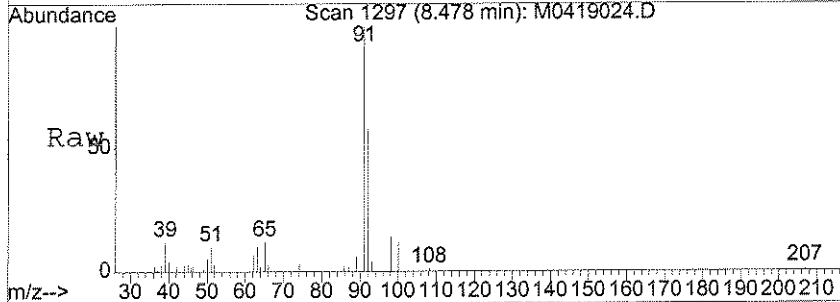
Tgt Ion: 83 Resp: 56494  
 Ion Ratio Lower Upper  
 83 100  
 85 63.0 44.2 84.2





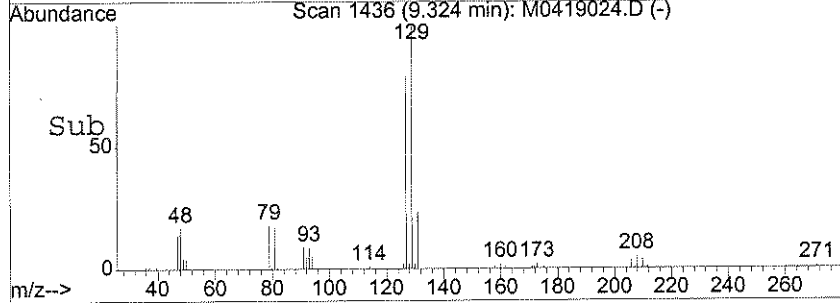
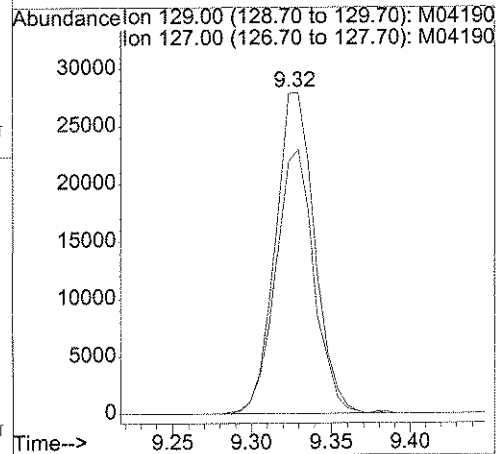
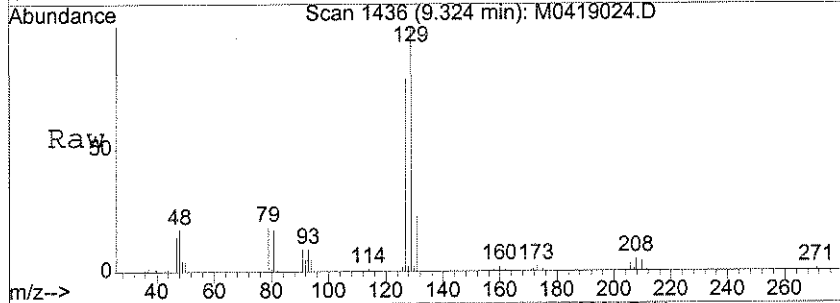
#52  
 Toluene  
 Concen: 0.56 ug/l  
 RT: 8.48 min Scan# 1297  
 Delta R.T. 0.00 min  
 Lab File: M0419024.D  
 Acq: 19 Apr 2007 21:36

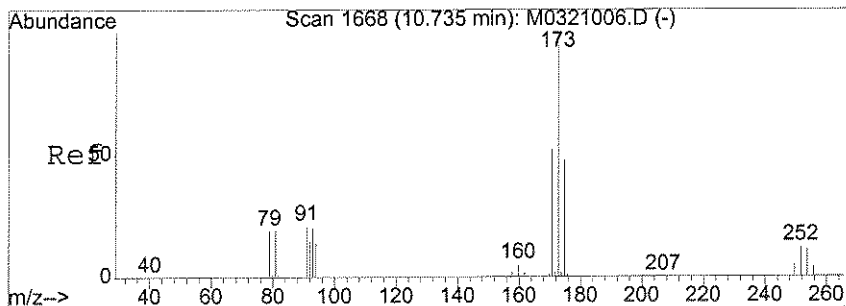
Tgt Ion: 92 Resp: 12409  
 Ion Ratio Lower Upper  
 92 100  
 91 176.0 133.7 200.5



#59  
 Dibromochloromethane  
 Concen: 7.57 ug/l  
 RT: 9.32 min Scan# 1436  
 Delta R.T. 0.00 min  
 Lab File: M0419024.D  
 Acq: 19 Apr 2007 21:36

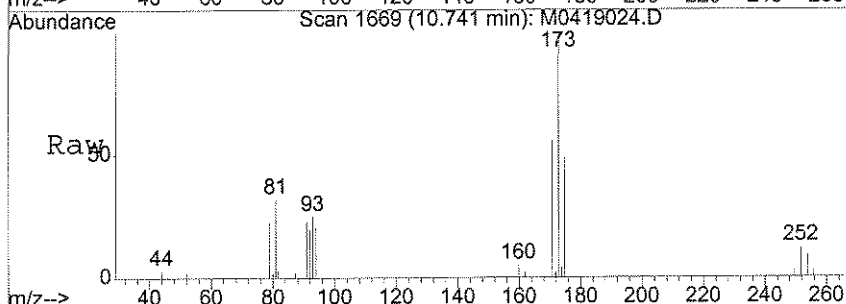
Tgt Ion: 129 Resp: 48117  
 Ion Ratio Lower Upper  
 129 100  
 127 80.4 58.9 98.9



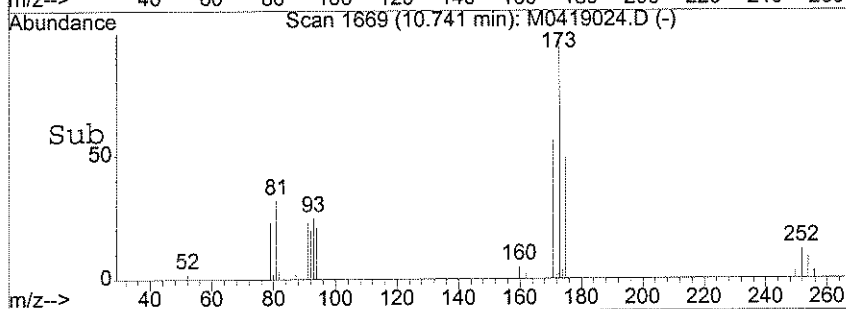
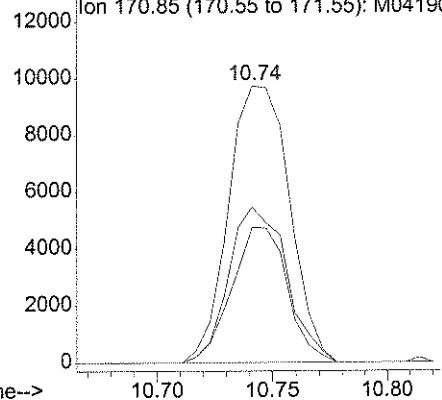


#68  
 Bromoform  
 Concen: 5.66 ug/l  
 RT: 10.74 min Scan# 1669  
 Delta R.T. 0.00 min  
 Lab File: M0419024.D  
 Acq: 19 Apr 2007 21:36

Tgt Ion	Resp	Lower	Upper
173	17870		
175	44.5	40.0	60.0
171	53.2	39.8	59.8



Abundance  
 Ion 172.85 (172.55 to 173.55): M04190  
 Ion 174.85 (174.55 to 175.55): M04190  
 Ion 170.85 (170.55 to 171.55): M04190



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-8

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-010  
 Lab File ID: M0419025.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 22:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.67	
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.26	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.71	



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-8

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-010  
 Lab File ID: M0419025.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 22:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-8

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-010  
 Lab File ID: M0419025.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 22:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

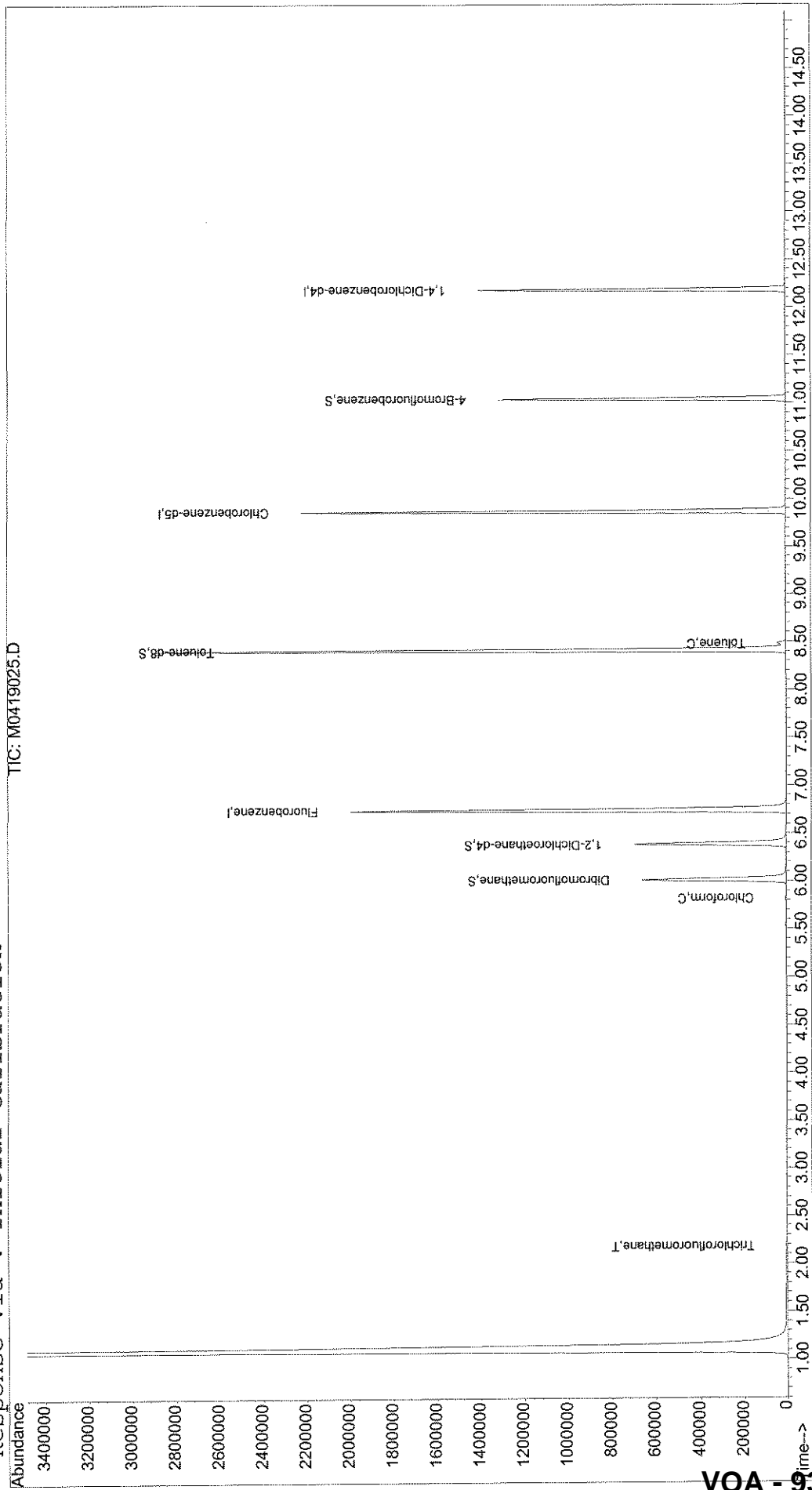
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419025.D Vial: 68  
Acq On : 19 Apr 2007 22:00 Operator: LH  
Sample : JPL33-010 Inst : MOBY  
Misc : 5ml +IS/SS #4 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:21 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 93

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419025.D  
 Acq On : 19 Apr 2007 22:00  
 Sample : JPL33-010  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:21 2007

Vial: 68  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
						Rcv (Ar )
1) Fluorobenzene	6.74	96	1925268	50.00	ug/l	0.00 98.63%
50) Chlorobenzene-d5	9.87	82	650221	50.00	ug/l	0.00 92.45%
70) 1,4-Dichlorobenzene-d4	12.18	152	350803	50.00	ug/l	0.00 90.13%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	461530	50.81	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.62%
37) 1,2-Dichloroethane-d4	6.39	65	517507	49.77	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.54%
51) Toluene-d8	8.41	98	1889202	52.61	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.22%
72) 4-Bromofluorobenzene	11.04	95	394161	54.47	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	108.94%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	1.97	64	73	N.D.		
7) Trichlorofluoromethane	2.18	101	6610mS	0.67 ug/l	#	34
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.84	96	65	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.94	43	80	N.D.		
12) Iodomethane	3.03	142	65	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.03	76	63	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	4.66	63	69	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419025.D  
 Acq On : 19 Apr 2007 22:00  
 Sample : JPL33-010  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:21 2007

Vial: 68  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	0.00	77	0	N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0	N.D.	
27) 2-Butanone	5.52	43	75	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	5.84	41	58	N.D.	
31) Chloroform	5.81	83	4032	0.26 ug/l	98
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	6.14	117	1006	N.D.	
36) 1,1-Dichloropropene	0.00	75	0	N.D.	
38) Benzene	6.40	78	1027	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D. d	
41) Trichloroethene	7.14	130	1033	N.D.	
42) Methylcyclohexane	7.30	83	73	N.D.	
43) 1,2-Dichloropropane	0.00	63	0	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	7.53	41	62	N.D.	
46) Bromodichloromethane	7.72	83	744	N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.	
48) cis-1,3-Dichloropropene	8.23	75	57	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	8.48	92	15391	0.71 ug/l	100
53) trans-1,3-Dichloropropene	8.67	75	57	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	8.94	97	60	N.D.	
56) Tetrachloroethene	9.03	166	480	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D. d	
59) Dibromochloromethane	9.34	129	73	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.91	112	85	N.D.	
62) 1-Chlorohexane	0.00	91	0	N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.99	91	233	N.D.	
65) m,p-Xylene	10.10	106	205	N.D.	
66) o-xylene	10.50	106	117	N.D.	
67) Styrene	0.00	104	0	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419025.D 8260B.M Fri Apr 20 10:21:38 2007

W 4/20/07

Quantitation Report

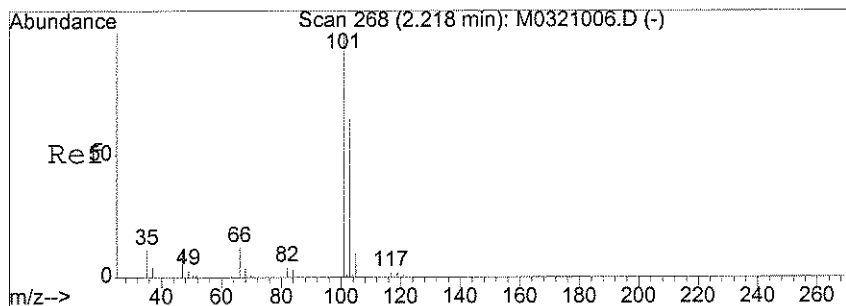
Data File : X:\MSVOA\MOBY\041907\M0419025.D  
 Acq On : 19 Apr 2007 22:00  
 Sample : JPL33-010  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:21 2007

Vial: 68  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

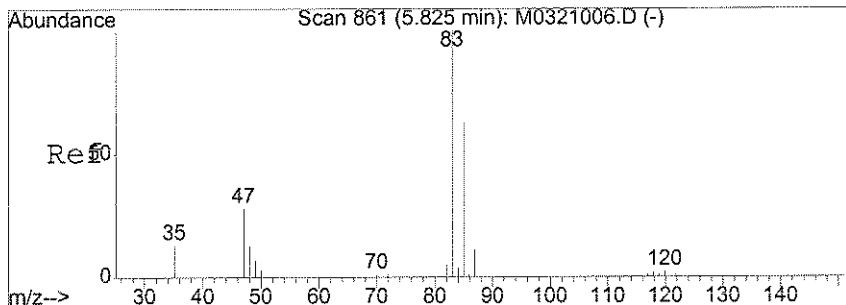
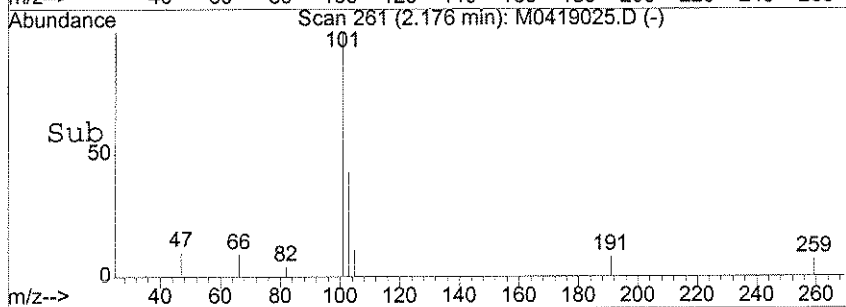
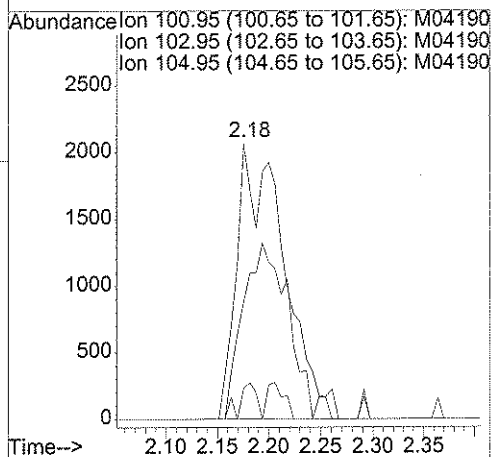
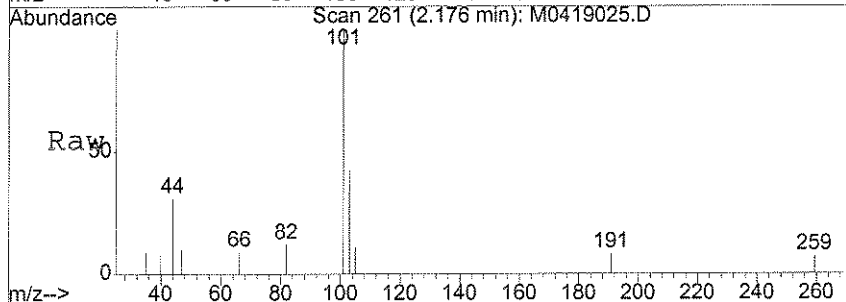
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	163		N.D.	
69) Isopropylbenzene	10.86	105	85		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.05	156	97		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	333		N.D.	
75) 1,2,3-Trichloropropane	11.21	110	61		N.D.	
76) n-Propylbenzene	0.00	120	0		N.D.	
77) 2-Chlorotoluene	11.26	91	249		N.D.	
78) 4-Chlorotoluene	11.48	91	59		N.D.	
79) 1,3,5-Trimethylbenzene	11.44	105	73		N.D.	
80) tert-Butylbenzene	11.90	119	67		N.D.	
81) 1,2,4-Trimethylbenzene	11.98	105	67		N.D.	
82) sec-butylbenzene	11.98	105	67		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	55		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	92		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.32	225	65		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



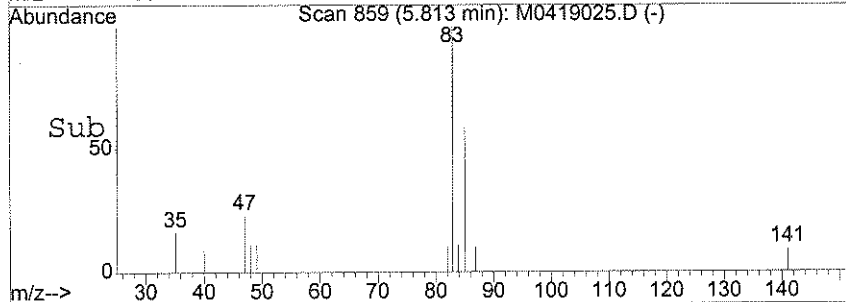
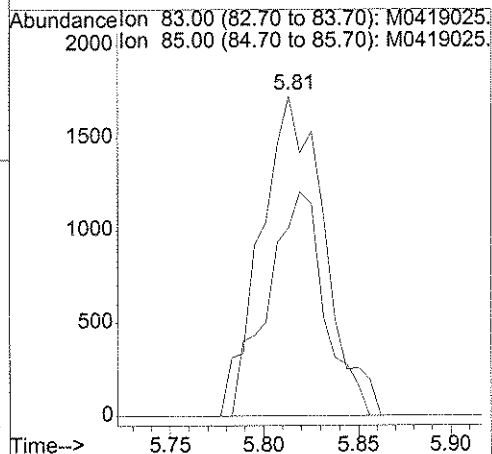
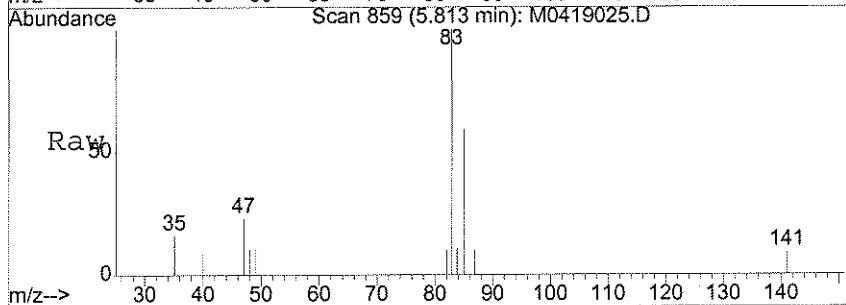
#7  
 Trichlorofluoromethane  
 Concen: 0.67 ug/l m  
 RT: 2.18 min Scan# 261  
 Delta R.T. 0.00 min  
 Lab File: M0419025.D  
 Acq: 19 Apr 2007 22:00

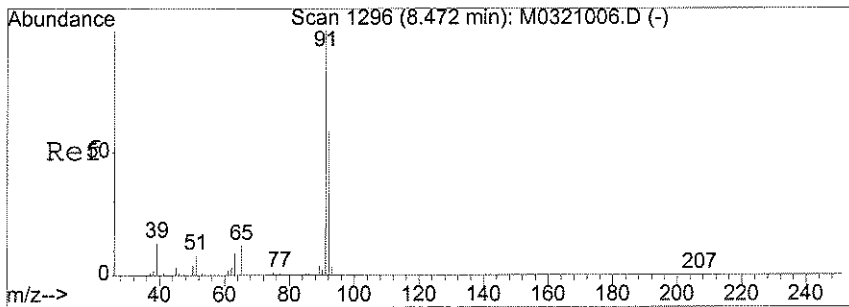
Tgt Ion	Resp	Lower	Upper
101	6610		
103	0.0	46.3	69.5#
105	4.8	9.0	13.4#



#31  
 Chloroform  
 Concen: 0.26 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419025.D  
 Acq: 19 Apr 2007 22:00

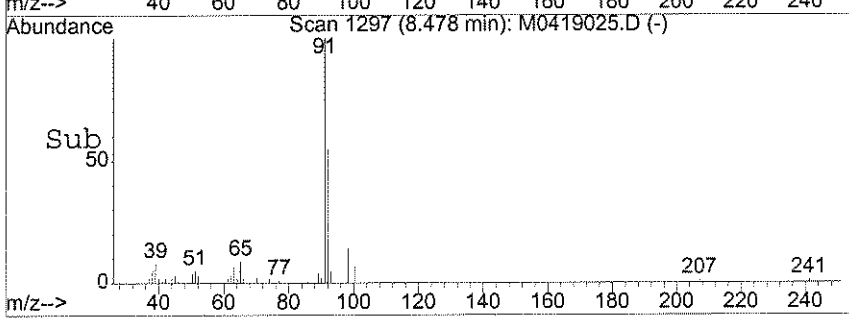
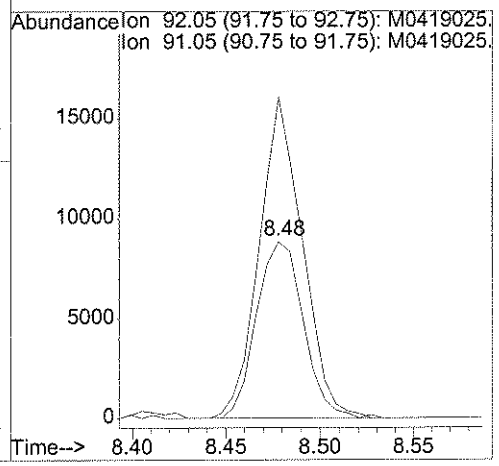
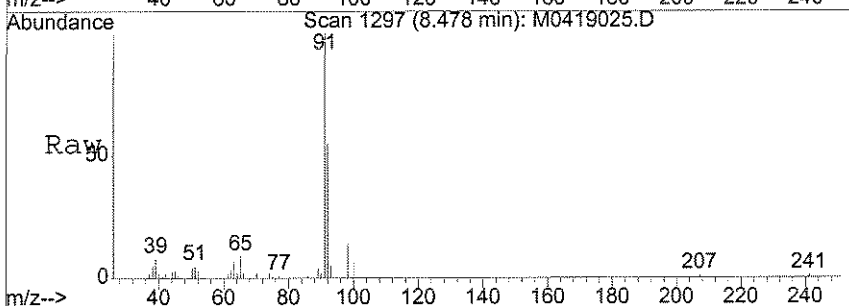
Tgt Ion	Resp	Lower	Upper
83	4032		
85	62.6	41.2	81.2





#52  
 Toluene  
 Concen: 0.71 ug/l  
 RT: 8.48 min Scan# 1297  
 Delta R.T. 0.00 min  
 Lab File: M0419025.D  
 Acq: 19 Apr 2007 22:00

Tgt Ion: 92 Resp: 15391  
 Ion Ratio Lower Upper  
 92 100  
 91 167.1 133.7 200.5





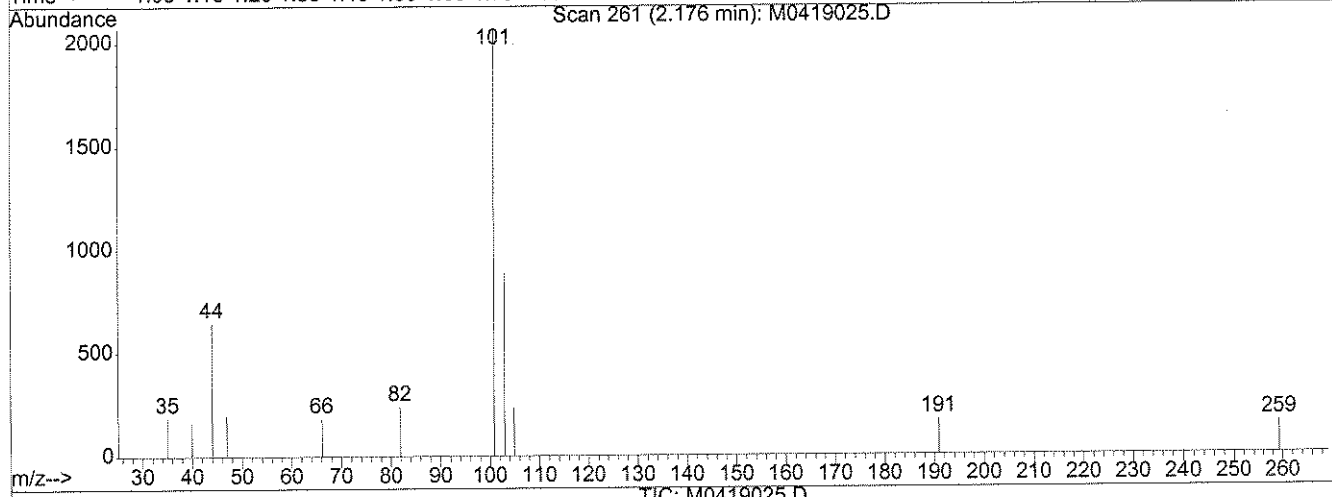
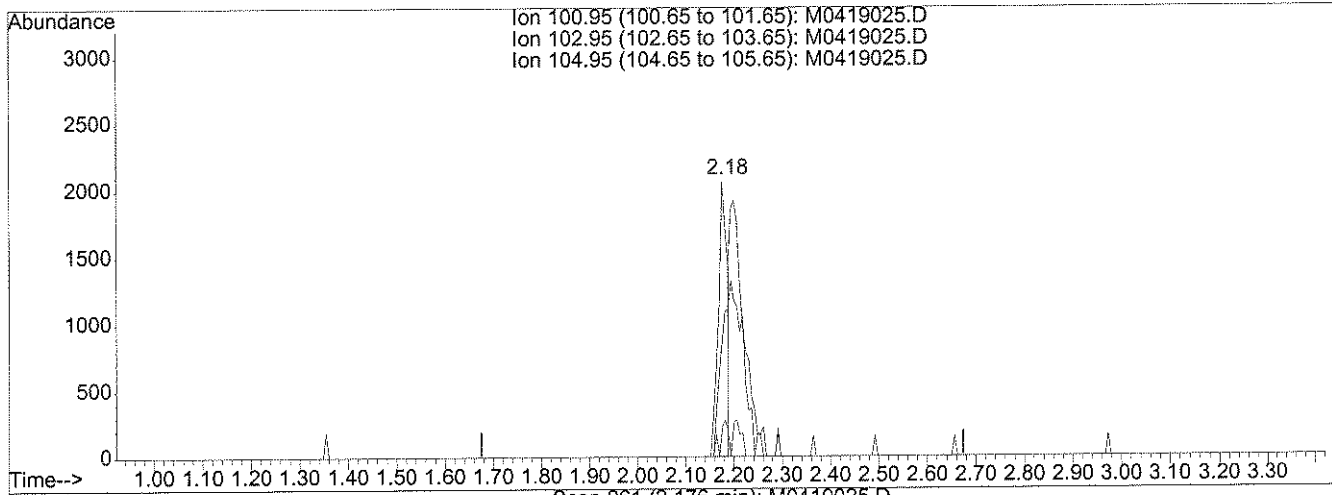
Quantitation Report (Qedit)

Data File : X:\MSVOA\MOBY\041907\M0419025.D  
 Acq On : 19 Apr 2007 22:00  
 Sample : JPL33-010  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:05 2007

Vial: 68  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: temp.res

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Multiple Level Calibration



(7) Trichlorofluoromethane (T)

2.18min 0.27ug/l

response 2705

Ion	Exp%	Act%
100.95	100	100
102.95	57.90	0.00#
104.95	11.20	11.76
0.00	0.00	0.00

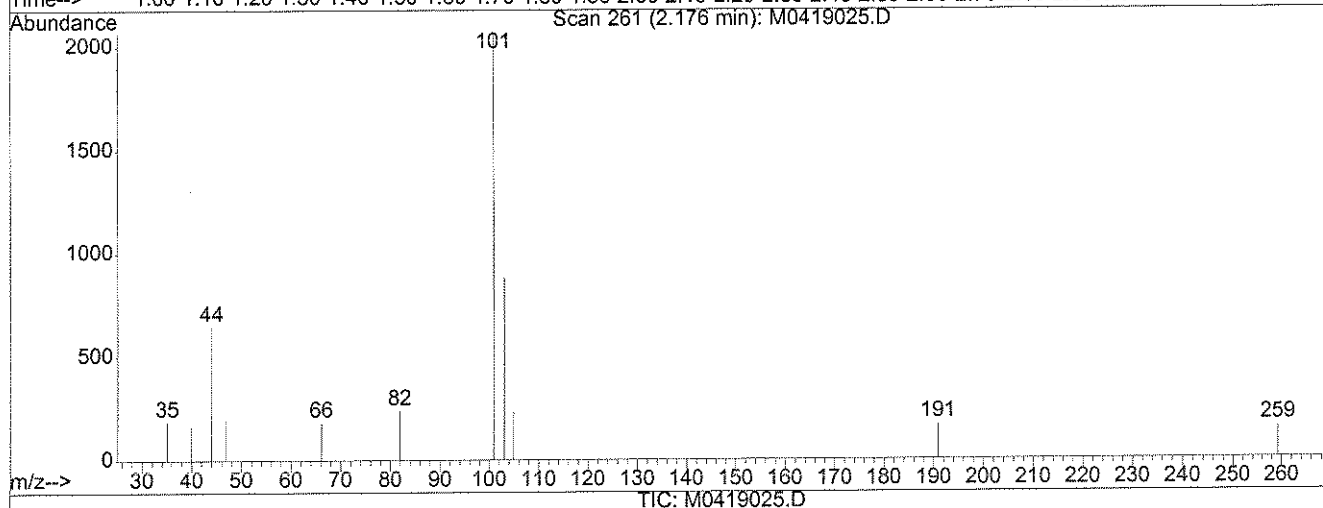
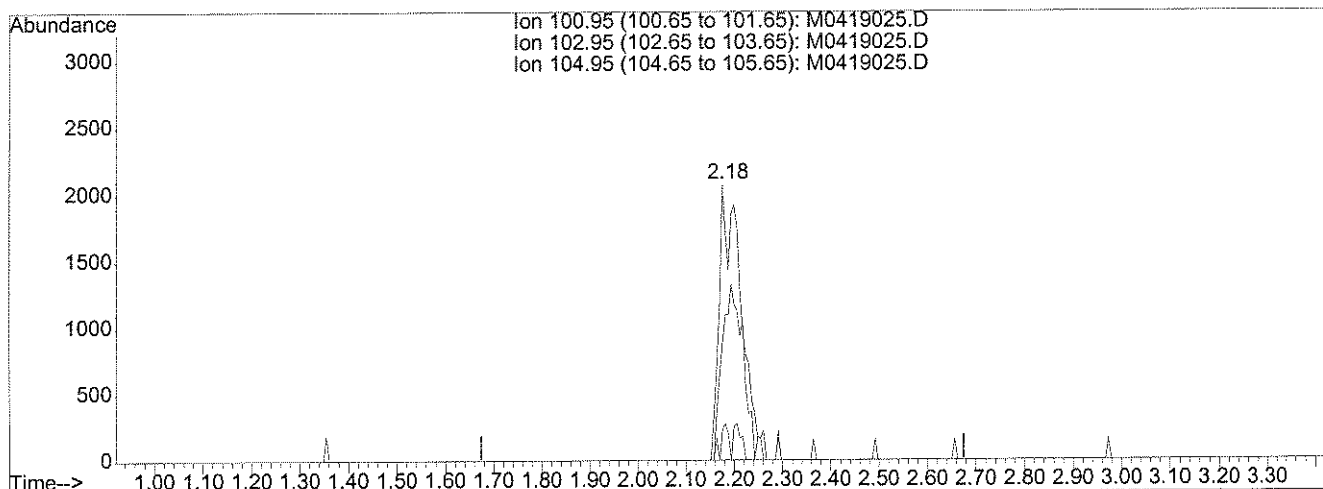
Quantitation Report (Qedit)

Data File : X:\MSVOA\MOBY\041907\M0419025.D  
 Acq On : 19 Apr 2007 22:00  
 Sample : JPL33-010  
 Misc : 5ml +IS/SS #4  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:20 2007

Vial: 68  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: temp.res

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Multiple Level Calibration



(7) Trichlorofluoromethane (T)

2.18min 0.67ug/l m

response 6610

Ion	Exp%	Act%
100.95	100	100
102.95	57.90	0.00#
104.95	11.20	4.81#
0.00	0.00	0.00

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-13-4/12/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-011  
 Lab File ID: M0419015.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 18:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-13-4/12/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-011  
 Lab File ID: M0419015.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 18:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-13-4/12/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-011  
 Lab File ID: M0419015.D  
 Date Collected: 04/12/2007  
 Date/Time Analyzed: 04/19/2007 18:00  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

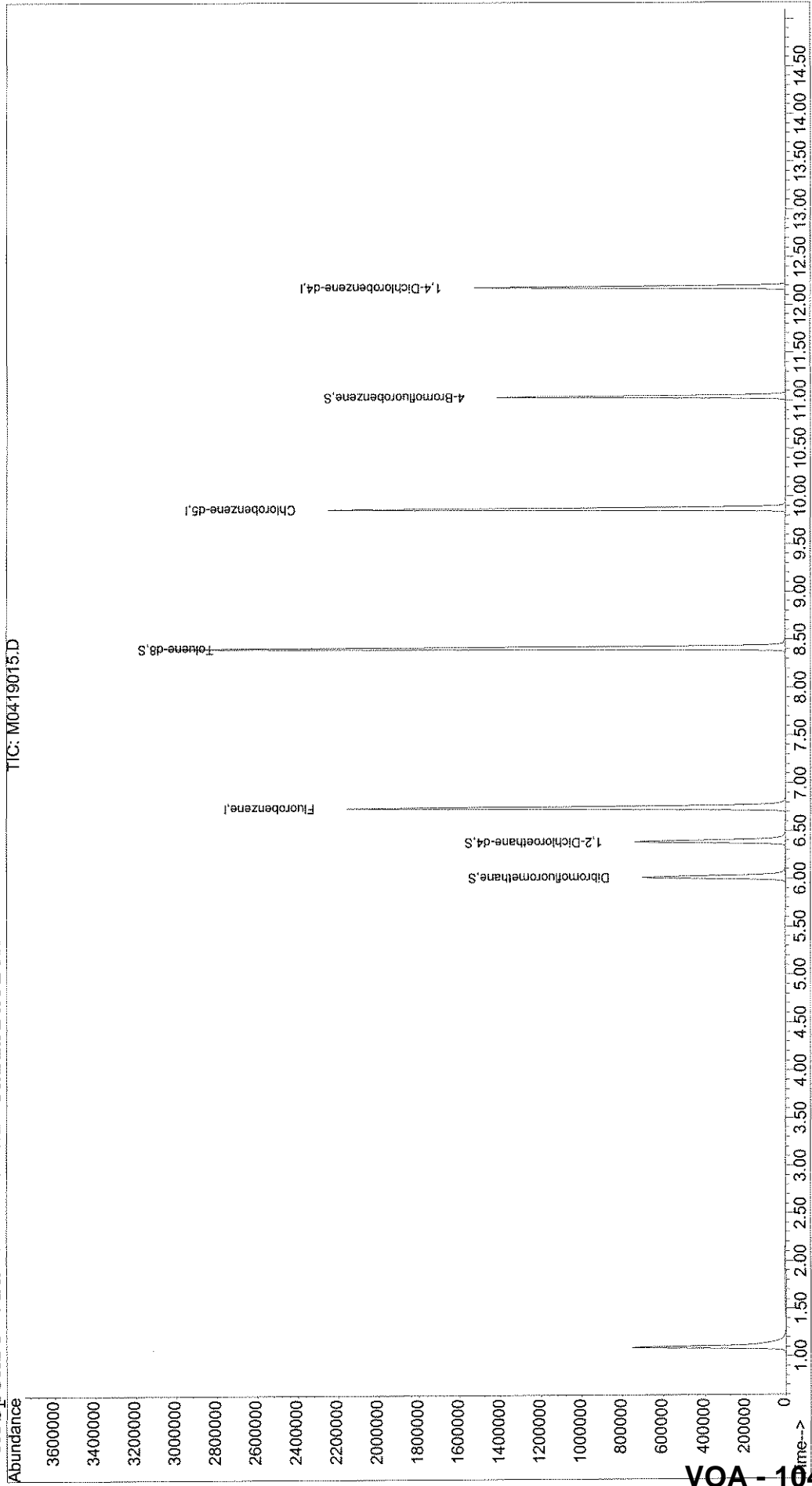
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419015.D Vial: 58  
Acq On : 19 Apr 2007 18:00 Operator: LH  
Sample : JPL33-011 TB Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:04 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 104

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419015.D  
 Acq On : 19 Apr 2007 18:00  
 Sample : JPL33-011 TB  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:04 2007

Vial: 58  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)

Title : VOA 8260- 5ML Calibration 5973M

Last Update : Tue Apr 17 16:33:44 2007

Response via : Initial Calibration

DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	2035144	50.00	ug/l	0.00	104.26%
50) Chlorobenzene-d5	9.87	82	687936	50.00	ug/l	0.00	97.82%
70) 1,4-Dichlorobenzene-d4	12.18	152	370073	50.00	ug/l	0.00	95.08%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	480888	50.09	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.18%	
37) 1,2-Dichloroethane-d4	6.39	65	550839	50.12	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.24%	
51) Toluene-d8	8.41	98	2004851	52.77	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.54%	
72) 4-Bromofluorobenzene	11.04	95	419650	54.97	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.94%	

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	1.43	50	57	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	0.00	96	0	N.D.	
6) Chloroethane	1.94	64	64	N.D.	
7) Trichlorofluoromethane	0.00	101	0	N.D.	
8) Acrolein	0.00	56	0	N.D.	
9) 1,1-Dichloroethene	0.00	96	0	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	
11) Acetone	2.98	43	186	N.D.	
12) Iodomethane	2.98	142	57	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.04	76	301	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	3.34	40	72	N.D.	
17) Methyl Acetate	3.41	43	81	N.D.	
18) Methylene Chloride	0.00	84	0	N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.	
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	0.00	73	0	N.D.	
22) 1,1-Dichloroethane	0.00	63	0	N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419015.D  
 Acq On : 19 Apr 2007 18:00  
 Sample : JPL33-011 TB  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:04 2007

Vial: 58  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.62	43	78		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.51	43	56		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	5.86	56	68		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	1131		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	72		N.D.	
42) Methylcyclohexane	7.30	83	55		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	7.71	93	63		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	7.73	83	56		N.D.	
47) 2-Chloroethyl vinyl ether	8.01	63	61		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.40	92	230		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.82	97	76		N.D.	
56) Tetrachloroethene	9.02	166	78		N.D.	
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	503		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	9.87	131	57		N.D.	
64) Ethylbenzene	9.99	91	457		N.D.	
65) m,p-Xylene	10.11	106	369		N.D.	
66) o-xylene	10.52	106	155		N.D.	
67) Styrene	0.00	104	0		N.D.	

(#) = qualifier out of range (m) = manual integration



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419015.D  
 Acq On : 19 Apr 2007 18:00  
 Sample : JPL33-011 TB  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:04 2007

Vial: 58  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	264		N.D.	
69) Isopropylbenzene	10.86	105	82		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.07	156	61		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	310		N.D.	
75) 1,2,3-Trichloropropane	11.40	110	68		N.D.	
76) n-Propylbenzene	11.03	120	57		N.D.	
77) 2-Chlorotoluene	11.37	91	246		N.D.	
78) 4-Chlorotoluene	11.49	91	248		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	226		N.D.	
80) tert-Butylbenzene	11.76	119	218		N.D.	
81) 1,2,4-Trimethylbenzene	11.81	105	58		N.D.	
82) sec-butylbenzene	11.98	105	448		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	79		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.22	146	64		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419015.D 8260B.M Fri Apr 20 10:04:50 2007

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-012  
 Lab File ID: M0419026.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.37	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	1.8	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-012  
 Lab File ID: M0419026.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-5

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-012  
 Lab File ID: M0419026.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

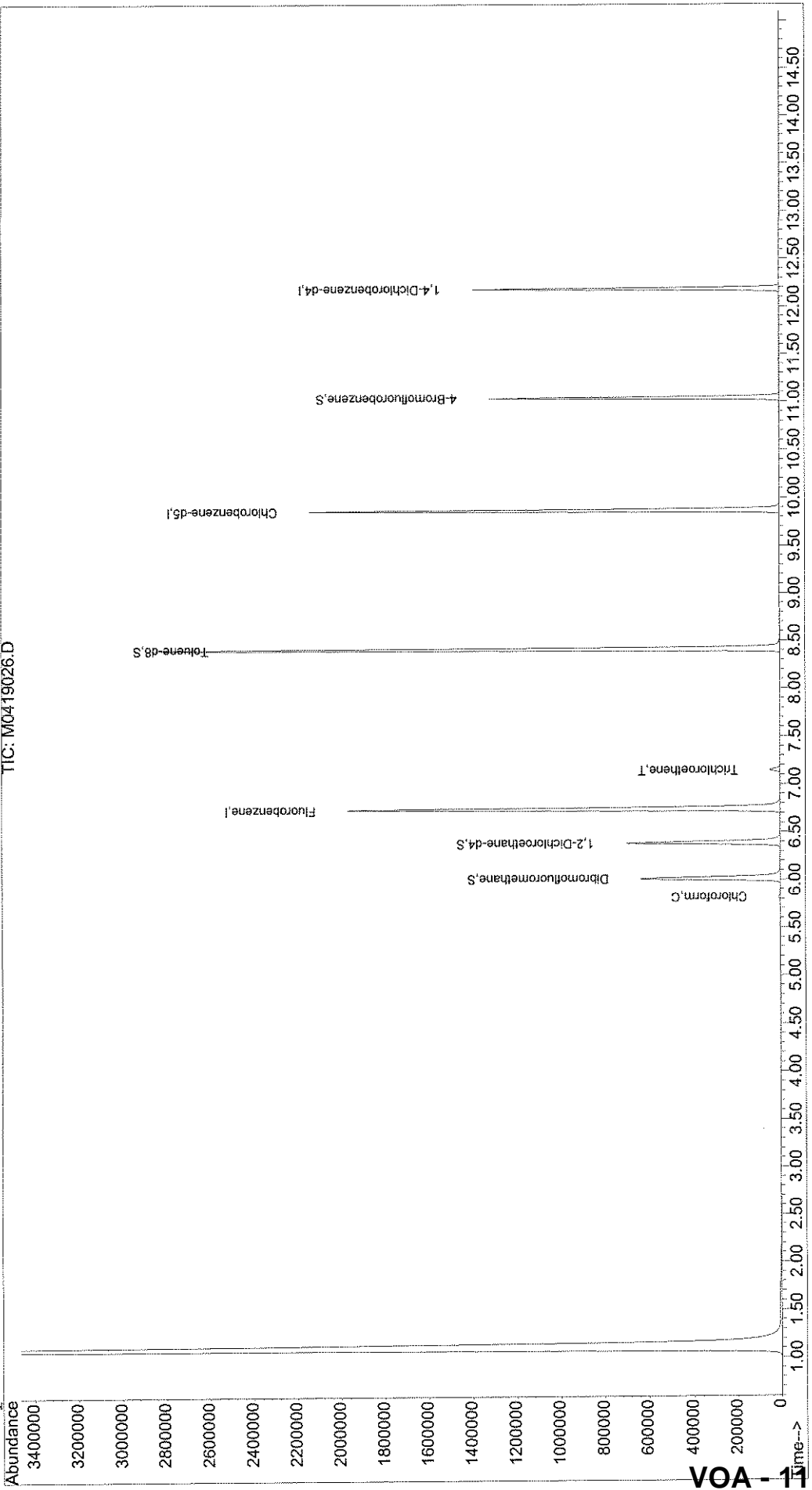
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419026.D Vial: 69  
Acq On : 19 Apr 2007 22:24 Operator: LH  
Sample : JPL33-012 Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:22 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA-111

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419026.D  
 Acq On : 19 Apr 2007 22:24  
 Sample : JPL33-012  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:22 2007

Vial: 69  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
						Rcv(Ar )
1) Fluorobenzene	6.75	96	1895786	50.00	ug/l	0.00 97.12%
50) Chlorobenzene-d5	9.87	82	643468	50.00	ug/l	0.00 91.49%
70) 1,4-Dichlorobenzene-d4	12.18	152	343217	50.00	ug/l	0.00 88.18%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	448460	50.14	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	100.28%
37) 1,2-Dichloroethane-d4	6.39	65	514622	50.26	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.52%
51) Toluene-d8	8.41	98	1865783	52.50	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	105.00%
72) 4-Bromofluorobenzene	11.04	95	387174	54.69	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.38%

Target Compounds

					Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.	
3) Chloromethane	0.00	50	0	N.D.	
4) Vinyl Chloride	0.00	62	0	N.D.	
5) Bromomethane	1.83	96	58	N.D.	
6) Chloroethane	1.91	64	56	N.D.	
7) Trichlorofluoromethane	2.19	101	168	N.D.	
8) Acrolein	2.79	56	55	N.D.	
9) 1,1-Dichloroethene	2.86	96	57	N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.	d
11) Acetone	2.90	43	73	N.D.	
12) Iodomethane	0.00	142	0	N.D.	
13) Bromoethane	0.00	108	0	N.D.	
14) Carbon Disulfide	3.03	76	67	N.D.	
15) Allyl chloride	0.00	76	0	N.D.	
16) Acetonitrile	0.00	40	0	N.D.	d
17) Methyl Acetate	3.33	43	72	N.D.	
18) Methylene Chloride	3.47	84	69	Below Cal	# 54
19) trans-1,2-Dichloroethene	3.92	96	66	N.D.	LW 4/20/07
20) Acrylonitrile	0.00	53	0	N.D.	
21) Methyl tert-butyl ether	3.92	73	57	N.D.	
22) 1,1-Dichloroethane	4.51	63	61	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419026.D 8260B.M Fri Apr 20 10:22:58 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419026.D  
 Acq On : 19 Apr 2007 22:24  
 Sample : JPL33-012  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:22 2007

Vial: 69  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
23) Vinyl acetate	0.00	43	0	N.D.	
24) Chloroprene	0.00	53	0	N.D.	
25) 2,2-Dichloropropane	5.41	77	55	N.D.	
26) cis-1,2-Dichloroethene	5.50	96	64	N.D.	
27) 2-Butanone	5.52	43	170	N.D.	
28) Propionitrile	0.00	54	0	N.D.	
29) Bromochloromethane	0.00	128	0	N.D.	
30) Methacrylonitrile	5.86	41	56	N.D.	
31) Chloroform	5.81	83	5578	0.37 ug/l	95
32) 1,1,1-Trichloroethane	5.97	97	61	N.D.	
34) Cyclohexane	0.00	56	0	N.D.	
35) Carbon Tetrachloride	6.14	117	1188	N.D.	
36) 1,1-Dichloropropene	6.19	75	66	N.D.	
38) Benzene	6.40	78	715	N.D.	
39) 1,2-Dichloroethane	0.00	62	0	N.D.	
40) Isobutanol	0.00	43	0	N.D.	
41) Trichloroethene	7.15	130	15028	1.76 ug/l	92
42) Methylcyclohexane	7.16	83	188	N.D.	
43) 1,2-Dichloropropane	7.49	63	62	N.D.	
44) Dibromomethane	0.00	93	0	N.D.	
45) Methyl methacrylate	0.00	41	0	N.D.	
46) Bromodichloromethane	7.72	83	178	N.D.	
47) 2-Chloroethyl vinyl ether	7.99	63	57	N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0	N.D. d	
52) Toluene	8.48	92	204	N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.	
54) Ethyl methacrylate	0.00	69	0	N.D.	
55) 1,1,2-Trichloroethane	8.94	97	80	N.D.	
56) Tetrachloroethene	9.02	166	1059	N.D.	
57) 1,3-Dichloropropane	0.00	76	0	N.D.	
58) 2-Hexanone	0.00	43	0	N.D. d	
59) Dibromochloromethane	9.43	129	69	N.D.	
60) 1,2-Dibromoethane	0.00	107	0	N.D.	
61) Chlorobenzene	9.90	112	175	N.D.	
62) 1-Chlorohexane	0.00	91	0	N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.	
64) Ethylbenzene	9.99	91	128	N.D.	
65) m,p-Xylene	10.11	106	67	N.D.	
66) o-xylene	0.00	106	0	N.D.	
67) Styrene	10.53	104	55	N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419026.D 8260B.M Fri Apr 20 10:22:58 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419026.D  
 Acq On : 19 Apr 2007 22:24  
 Sample : JPL33-012  
 Misc : 5ml +IS/SS #2  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:22 2007

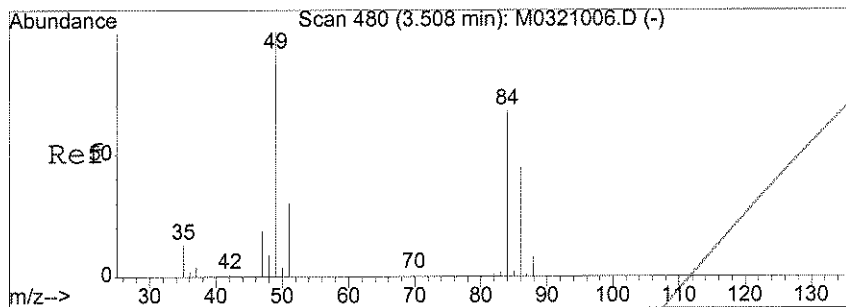
Vial: 69  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

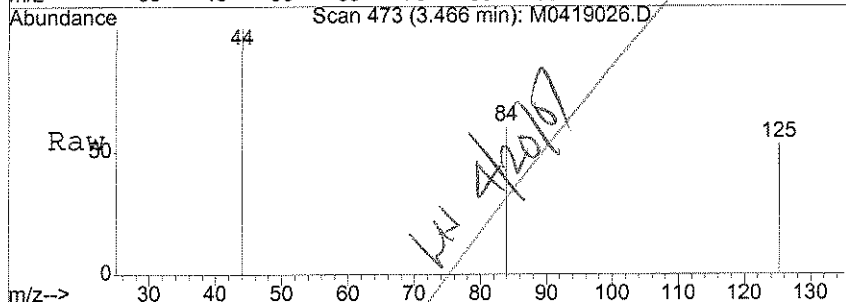
Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	164		N.D.	
69) Isopropylbenzene	10.85	105	57		N.D.	
71) trans-1,4-Dichloro-2-buten	11.13	53	55		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.13	83	61		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	173		N.D.	
76) n-Propylbenzene	11.03	120	59		N.D.	
77) 2-Chlorotoluene	0.00	91	0		N.D.	
78) 4-Chlorotoluene	11.69	91	100		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.77	119	58		N.D.	
81) 1,2,4-Trimethylbenzene	11.97	105	61		N.D.	
82) sec-butylbenzene	11.97	105	61		N.D.	
83) 1,3-Dichlorobenzene	12.12	146	65		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.12	146	65		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



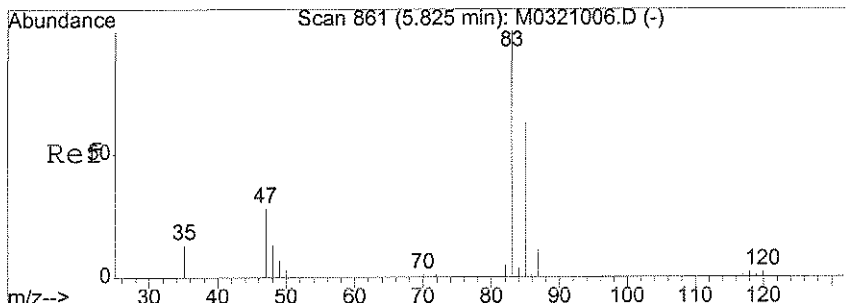
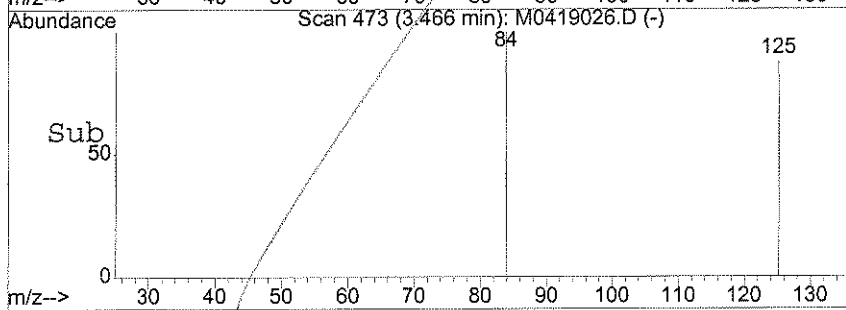
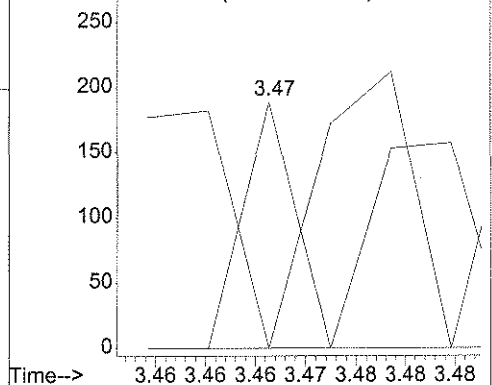


#18  
 Methylene Chloride  
 Concen: Below Cal  
 RT: 3.47 min Scan# 473  
 Delta R.T. -0.00 min  
 Lab File: M0419026.D  
 Acq: 19 Apr 2007 22:24

Tgt Ion	Resp	Lower	Upper
84	100		
49	189.9	133.5	173.5#
86	0.0	49.1	89.1#

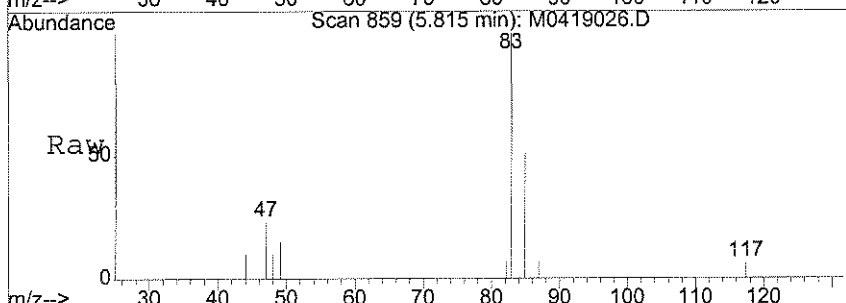


Abundance Ion 84.00 (83.70 to 84.70): M0419026  
 Ion 49.00 (48.70 to 49.70): M0419026  
 Ion 86.00 (85.70 to 86.70): M0419026

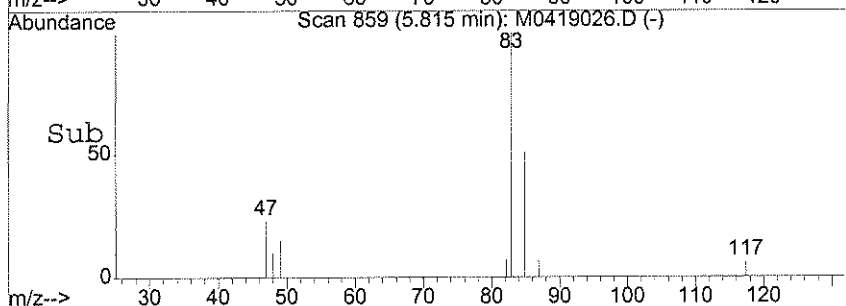
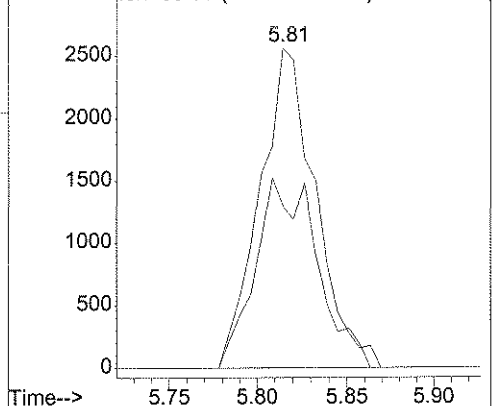


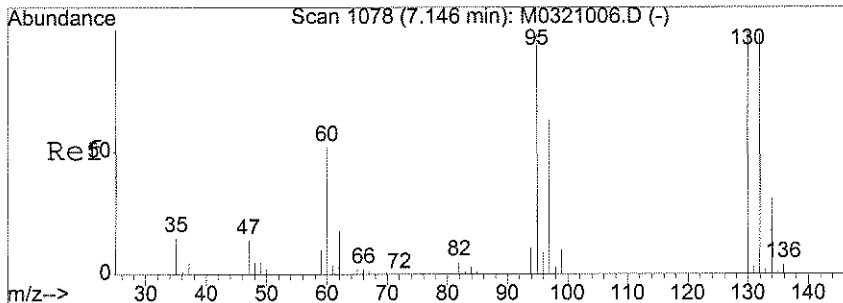
#31  
 Chloroform  
 Concen: 0.37 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419026.D  
 Acq: 19 Apr 2007 22:24

Tgt Ion	Resp	Lower	Upper
83	100		
85	65.3	41.2	81.2



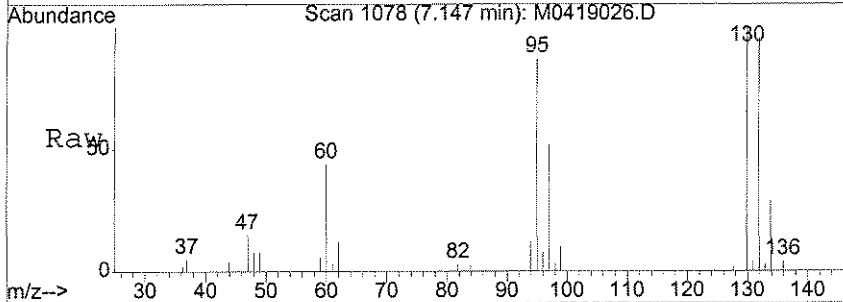
Abundance Ion 83.00 (82.70 to 83.70): M0419026  
 Ion 85.00 (84.70 to 85.70): M0419026



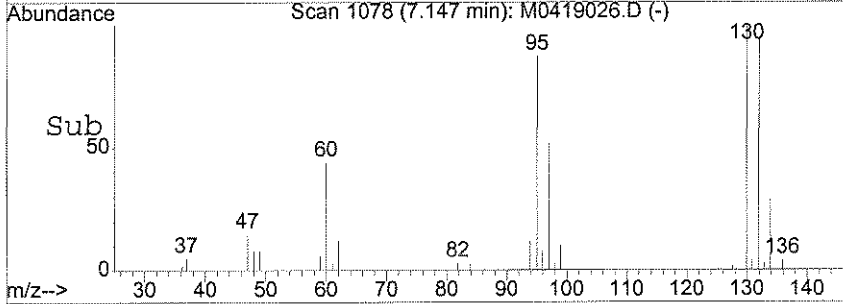
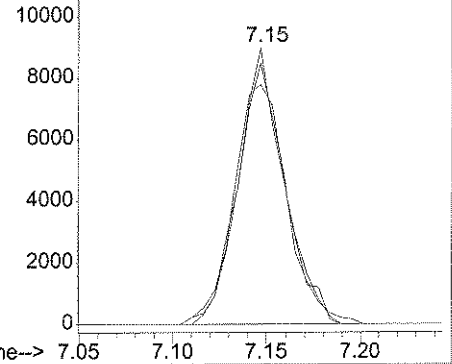


#41  
 Trichloroethene  
 Concen: 1.76 ug/l  
 RT: 7.15 min Scan# 1078  
 Delta R.T. 0.00 min  
 Lab File: M0419026.D  
 Acq: 19 Apr 2007 22:24

Tgt Ion	Resp	Lower	Upper
130	15028		
130	100		
132	98.8	75.0	115.0
95	101.6	69.4	109.4



Abundance  
 Ion 130.00 (129.70 to 130.70): M04190  
 Ion 132.00 (131.70 to 132.70): M04190  
 Ion 95.00 (94.70 to 95.70): M0419026



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-10

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-013  
 Lab File ID: M0419027.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.68	
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.61	
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	3.1	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.76	

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-10

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-013  
 Lab File ID: M0419027.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	1.2	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-10

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-013  
 Lab File ID: M0419027.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 22:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

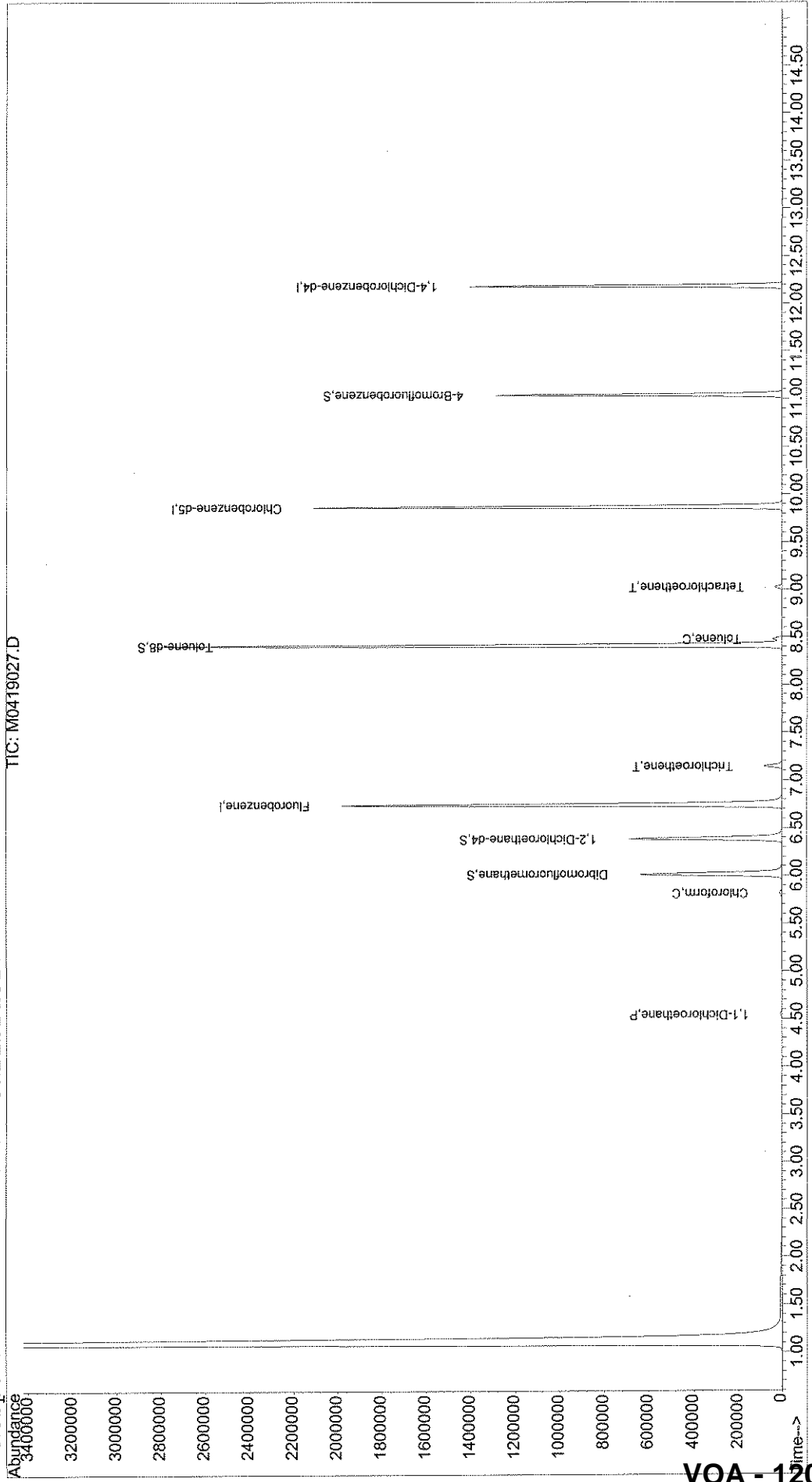
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419027.D Vial: 70  
Acq On : 19 Apr 2007 22:48 Operator: LH  
Sample : JPL33-013 Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:24 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA-120

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419027.D  
 Acq On : 19 Apr 2007 22:48  
 Sample : JPL33-013  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:24 2007

Vial: 70  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)	Rcv (Ar)
1) Fluorobenzene	6.74	96	1877662	50.00	ug/l	0.00	96.20%
50) Chlorobenzene-d5	9.86	82	629978	50.00	ug/l	0.00	89.57%
70) 1,4-Dichlorobenzene-d4	12.18	152	342366	50.00	ug/l	0.00	87.96%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	449032	50.69	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.38%	
37) 1,2-Dichloroethane-d4	6.39	65	511246	50.42	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.84%	
51) Toluene-d8	8.41	98	1857767	53.39	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	106.78%	
72) 4-Bromofluorobenzene	11.04	95	384269	54.41	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	108.82%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	1.87	96	66		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	2.18	101	943		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	2.77	96	95		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D. d	
11) Acetone	0.00	43	0		N.D. d	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.03	76	207		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D. d	
17) Methyl Acetate	3.34	43	62		N.D.	
18) Methylene Chloride	0.00	84	0		N.D. d	
19) trans-1,2-Dichloroethene	3.87	96	415		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	0.00	73	0		N.D.	
22) 1,1-Dichloroethane	4.54	63	9830	0.68	ug/l	91

(#) = qualifier out of range (m) = manual integration  
 M0419027.D 8260B.M Fri Apr 20 10:24:47 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419027.D  
 Acq On : 19 Apr 2007 22:48  
 Sample : JPL33-013  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:24 2007

Vial: 70  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0	N.D.		
24) Chloroprene	0.00	53	0	N.D.		
25) 2,2-Dichloropropane	0.00	77	0	N.D.		
26) cis-1,2-Dichloroethene	5.38	96	581	N.D.		
27) 2-Butanone	5.50	43	62	N.D.		
28) Propionitrile	0.00	54	0	N.D.		
29) Bromochloromethane	0.00	128	0	N.D.		
30) Methacrylonitrile	0.00	41	0	N.D.		
31) Chloroform	5.82	83	9126	0.61	ug/l	94
32) 1,1,1-Trichloroethane	0.00	97	0	N.D.		
34) Cyclohexane	6.00	56	57	N.D.		
35) Carbon Tetrachloride	6.15	117	179	N.D.		
36) 1,1-Dichloropropene	0.00	75	0	N.D.		
38) Benzene	6.40	78	715	N.D.		
39) 1,2-Dichloroethane	0.00	62	0	N.D.		
40) Isobutanol	0.00	43	0	N.D.	d	
41) Trichloroethene	7.15	130	26383	3.12	ug/l	90
42) Methylcyclohexane	7.33	83	70	N.D.		
43) 1,2-Dichloropropane	0.00	63	0	N.D.		
44) Dibromomethane	0.00	93	0	N.D.		
45) Methyl methacrylate	7.49	41	63	N.D.		
46) Bromodichloromethane	7.74	83	654	N.D.		
47) 2-Chloroethyl vinyl ether	0.00	63	0	N.D.		
48) cis-1,3-Dichloropropene	0.00	75	0	N.D.		
49) 4-Methyl-2-pentanone	0.00	43	0	N.D.	d	
52) Toluene	8.48	92	16061	0.76	ug/l	98
53) trans-1,3-Dichloropropene	0.00	75	0	N.D.		
54) Ethyl methacrylate	0.00	69	0	N.D.		
55) 1,1,2-Trichloroethane	0.00	97	0	N.D.		
56) Tetrachloroethene	9.03	166	9015	1.19	ug/l	95
57) 1,3-Dichloropropane	0.00	76	0	N.D.		
58) 2-Hexanone	0.00	43	0	N.D.	d	
59) Dibromochloromethane	0.00	129	0	N.D.		
60) 1,2-Dibromoethane	0.00	107	0	N.D.		
61) Chlorobenzene	9.91	112	284	N.D.		
62) 1-Chlorohexane	0.00	91	0	N.D.	d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0	N.D.		
64) Ethylbenzene	9.99	91	339	N.D.		
65) m,p-Xylene	10.11	106	363	N.D.		
66) o-xylene	0.00	106	0	N.D.		
67) Styrene	0.00	104	0	N.D.		

LH 4/20/07



Quantitation Report

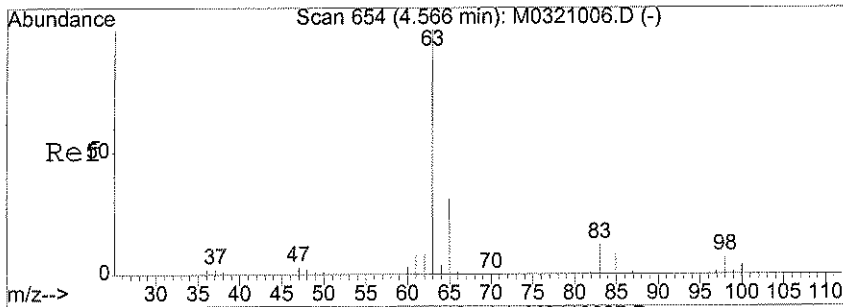
Data File : X:\MSVOA\MOBY\041907\M0419027.D  
 Acq On : 19 Apr 2007 22:48  
 Sample : JPL33-013  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:24 2007

Vial: 70  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

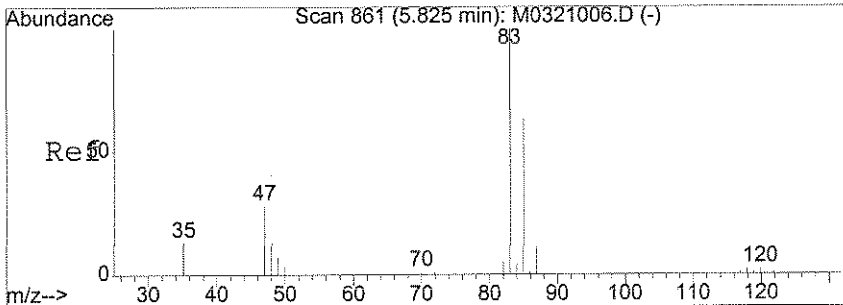
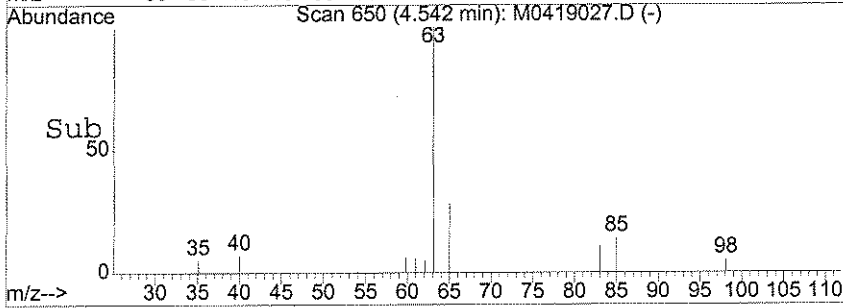
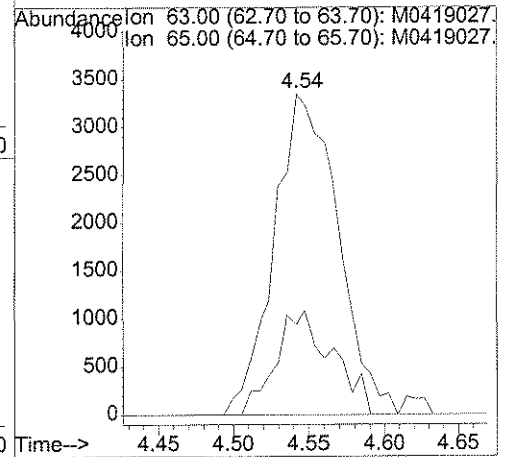
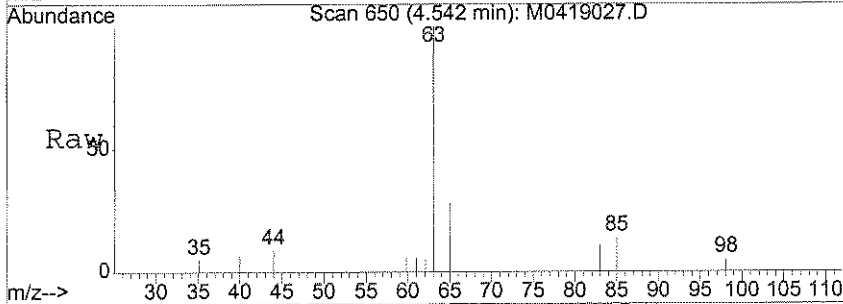
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	156		N.D.	
69) Isopropylbenzene	11.04	105	691		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	11.03	156	57		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.05	83	181		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	117		N.D.	
76) n-Propylbenzene	11.28	120	72		N.D.	
77) 2-Chlorotoluene	11.43	91	57		N.D.	
78) 4-Chlorotoluene	11.48	91	66		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.99	105	71		N.D.	
82) sec-butylbenzene	11.99	105	71		N.D.	
83) 1,3-Dichlorobenzene	12.19	146	56		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.19	146	56		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	141		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	0.00	225	0		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d



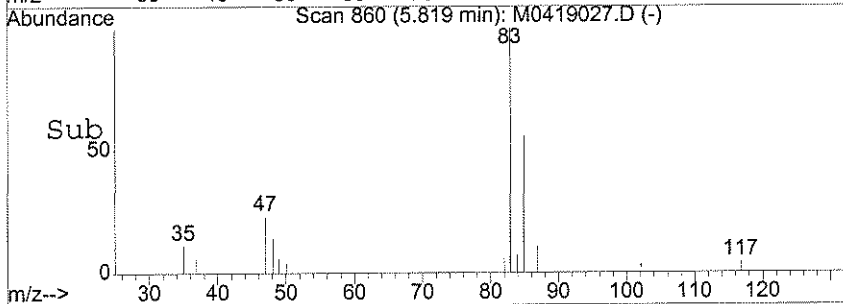
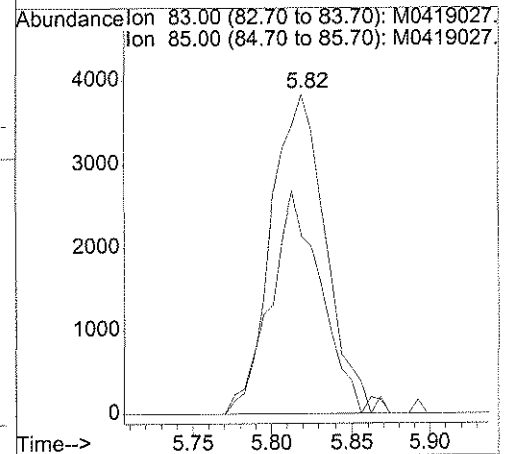
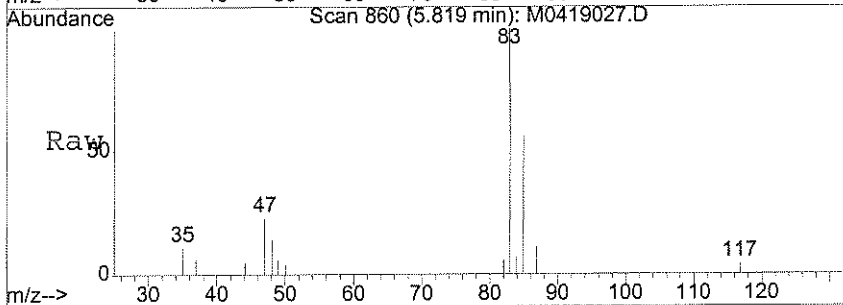
#22  
 1,1-Dichloroethane  
 Concen: 0.68 ug/l  
 RT: 4.54 min Scan# 650  
 Delta R.T. 0.00 min  
 Lab File: M0419027.D  
 Acq: 19 Apr 2007 22:48

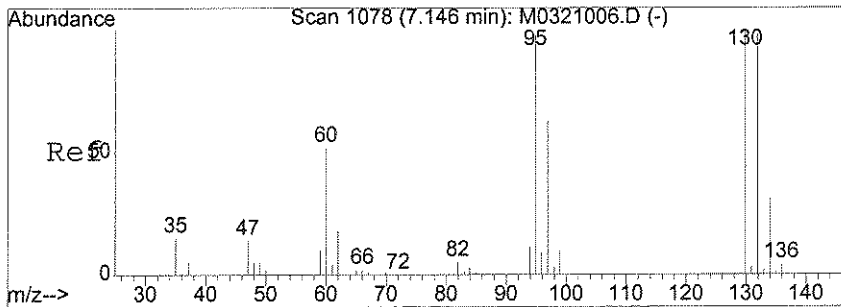
Tgt Ion: 63 Resp: 9830  
 Ion Ratio Lower Upper  
 63 100  
 65 28.7 13.7 53.7



#31  
 Chloroform  
 Concen: 0.61 ug/l  
 RT: 5.82 min Scan# 860  
 Delta R.T. 0.01 min  
 Lab File: M0419027.D  
 Acq: 19 Apr 2007 22:48

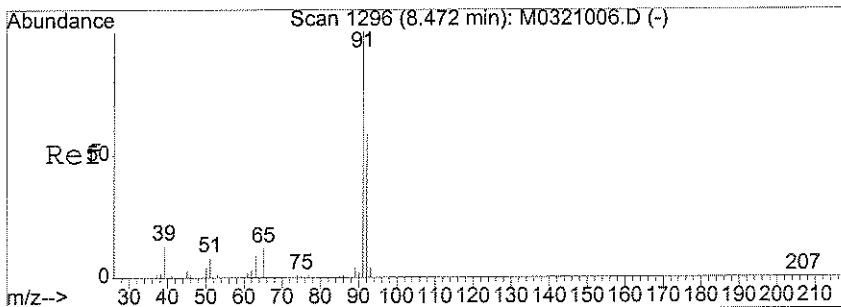
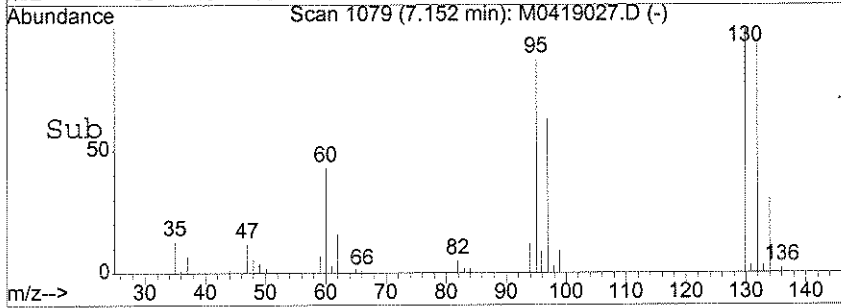
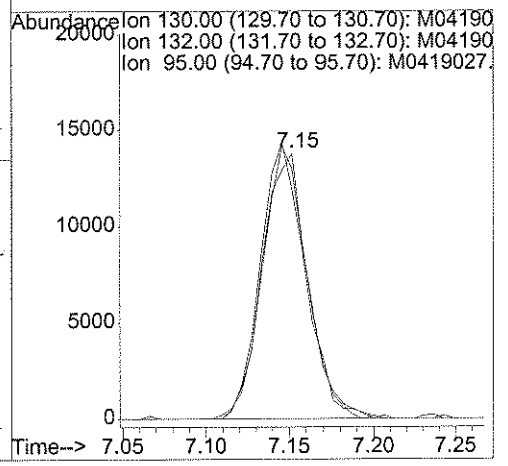
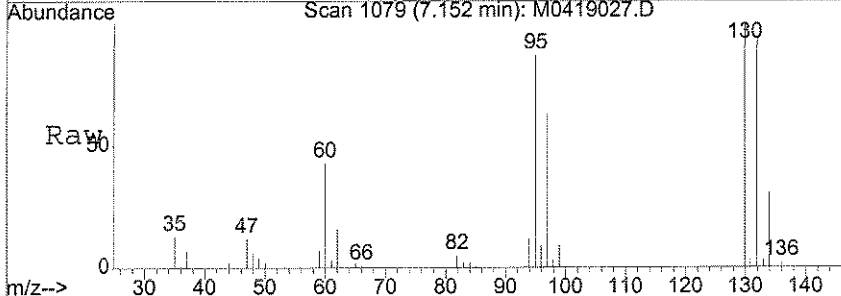
Tgt Ion: 83 Resp: 9126  
 Ion Ratio Lower Upper  
 83 100  
 85 66.0 41.2 81.2





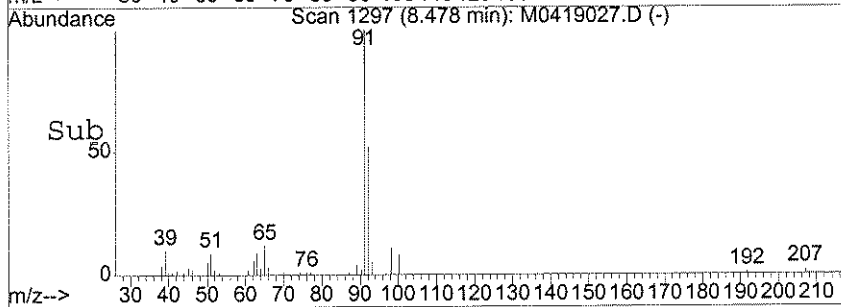
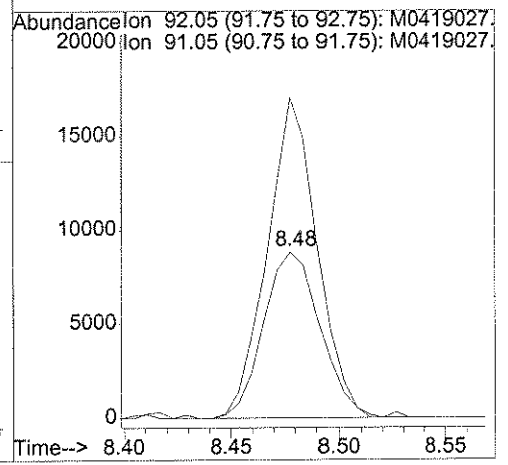
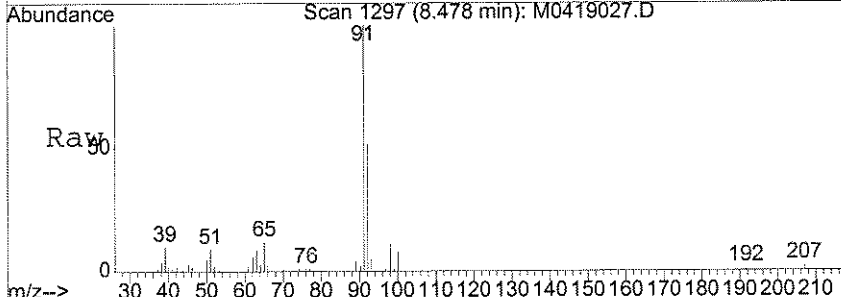
#41  
 Trichloroethene  
 Concen: 3.12 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0419027.D  
 Acq: 19 Apr 2007 22:48

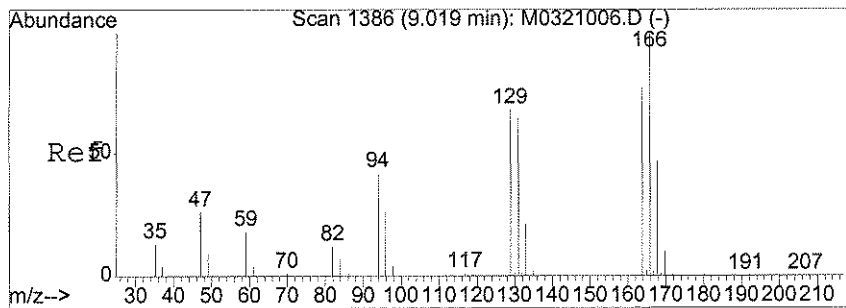
Tgt Ion	Resp	Lower	Upper
130	26383		
130	100		
132	100.7	75.0	115.0
95	102.1	69.4	109.4



#52  
 Toluene  
 Concen: 0.76 ug/l  
 RT: 8.48 min Scan# 1297  
 Delta R.T. 0.00 min  
 Lab File: M0419027.D  
 Acq: 19 Apr 2007 22:48

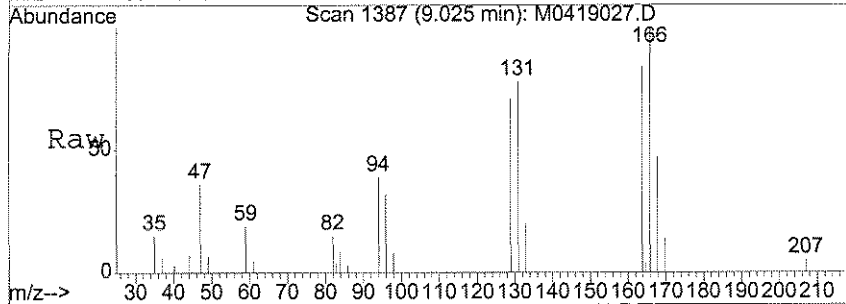
Tgt Ion	Resp	Lower	Upper
92	16061		
92	100		
91	170.2	133.7	200.5



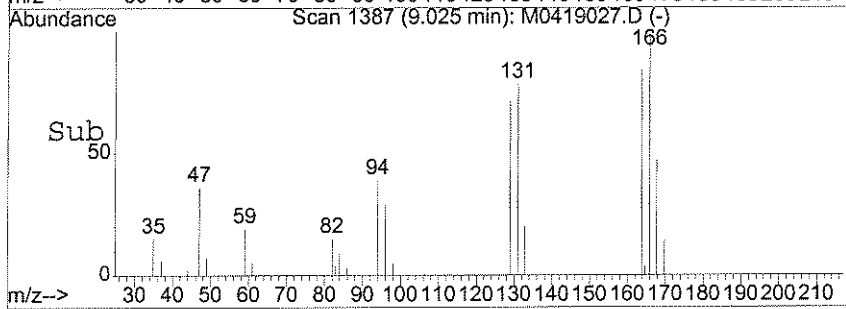
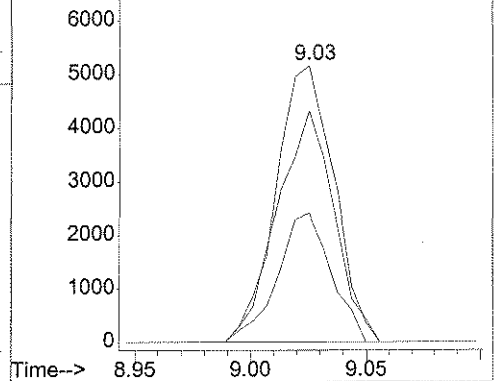


#56  
 Tetrachloroethene  
 Concen: 1.19 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0419027.D  
 Acq: 19 Apr 2007 22:48

Tgt Ion	Ratio	Lower	Upper
166	100		
164	81.9	63.3	94.9
168	43.2	39.6	59.4



Abundance Ion 165.95 (165.65 to 166.65): M04190  
 7000 Ion 163.95 (163.65 to 164.65): M04190  
 Ion 167.95 (167.65 to 168.65): M04190



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-14-4/13/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-014  
 Lab File ID: M0419016.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 18:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-14-4/13/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL33-014

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419016.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/13/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 18:24

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-14-4/13/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL33  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-014  
 Lab File ID: M0419016.D  
 Date Collected: 04/13/2007  
 Date/Time Analyzed: 04/19/2007 18:24  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

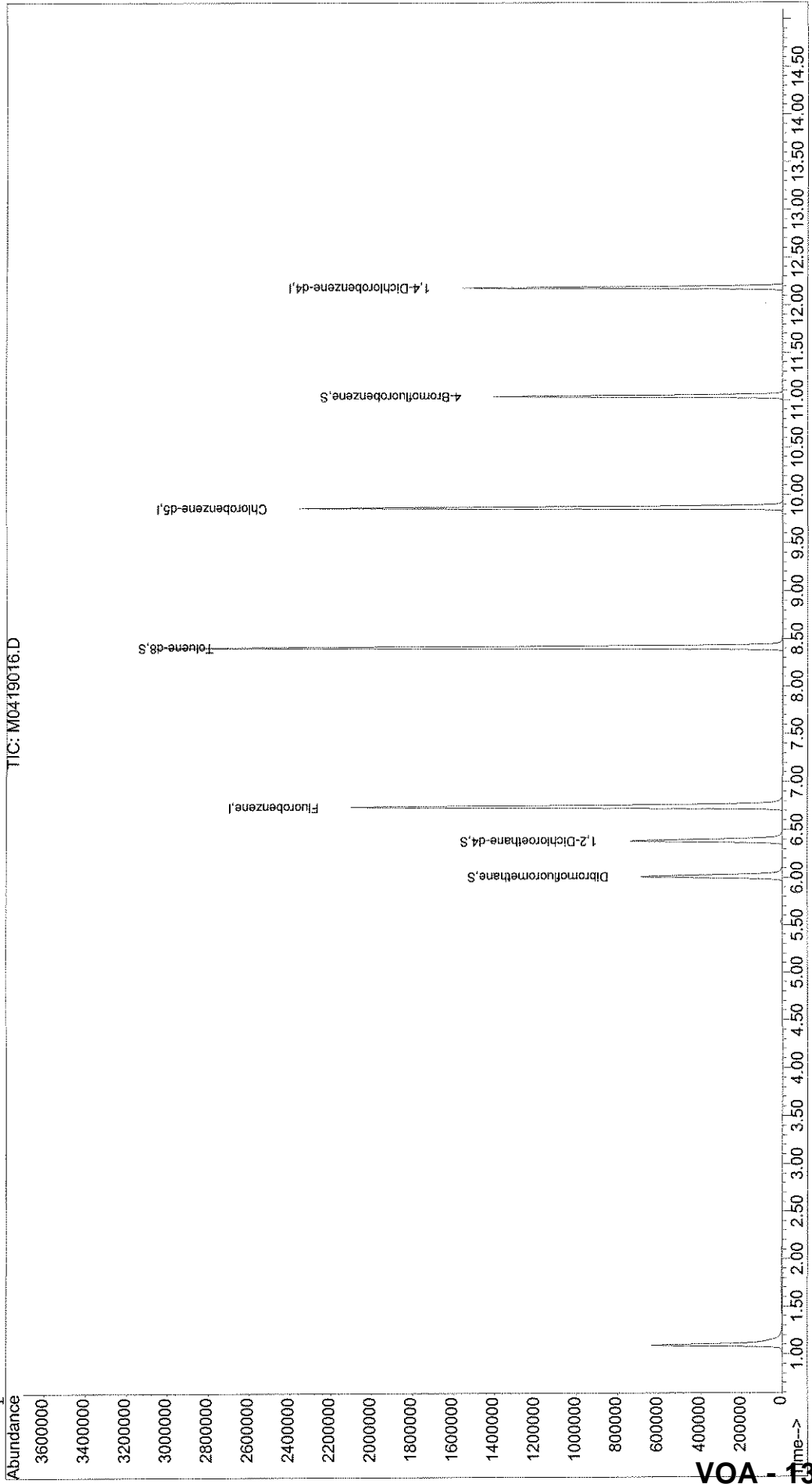
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419016.D Vial: 59  
Acq On : 19 Apr 2007 18:24 Operator: LH  
Sample : JPL33-014 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:06 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260 - 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration





Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419016.D  
 Acq On : 19 Apr 2007 18:24  
 Sample : JPL33-014 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:06 2007

Vial: 59  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	Rcv(Ar )
1) Fluorobenzene	6.74	96	2029196	50.00	ug/l	0.00	103.96%
50) Chlorobenzene-d5	9.87	82	692365	50.00	ug/l	0.00	98.45%
70) 1,4-Dichlorobenzene-d4	12.18	152	370936	50.00	ug/l	0.00	95.30%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	484910	50.65	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.30%	
37) 1,2-Dichloroethane-d4	6.39	65	548069	50.01	ug/l	0.00	
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.02%	
51) Toluene-d8	8.41	98	2002228	52.36	ug/l	0.00	
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.72%	
72) 4-Bromofluorobenzene	11.04	95	420672	54.98	ug/l	0.00	
Spiked Amount	50.000	Range	75 - 120	Recovery	=	109.96%	

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	1.41	50	61	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	1.81	96	71	N.D.		
6) Chloroethane	1.97	64	70	N.D.		
7) Trichlorofluoromethane	2.20	101	57	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	0.00	96	0	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	2.98	43	211	N.D.		
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.04	76	454	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	3.30	43	58	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	0.00	96	0	N.D.		
20) Acrylonitrile	0.00	53	0	N.D.		
21) Methyl tert-butyl ether	3.89	73	62	N.D.		
22) 1,1-Dichloroethane	4.56	63	107	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419016.D  
 Acq On : 19 Apr 2007 18:24  
 Sample : JPL33-014 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:06 2007

Vial: 59  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	0.00	43	0		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.35	43	78		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	0.00	41	0		N.D.	
31) Chloroform	0.00	83	0		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	6.03	56	60		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.20	75	57		N.D.	
38) Benzene	6.40	78	884		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	0.00	130	0		N.D.	
42) Methylcyclohexane	7.29	83	59		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	7.49	41	58		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	831		N.D.	
53) trans-1,3-Dichloropropene	8.66	75	74		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	0.00	97	0		N.D.	
56) Tetrachloroethene	9.01	166	58		N.D.	
57) 1,3-Dichloropropane	9.09	76	68		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.90	112	346		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	260		N.D.	
65) m,p-Xylene	10.11	106	373		N.D.	
66) o-xylene	10.51	106	64		N.D.	
67) Styrene	10.51	104	76		N.D.	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419016.D  
 Acq On : 19 Apr 2007 18:24  
 Sample : JPL33-014 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:06 2007

Vial: 59  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.74	173	111		N.D.	
69) Isopropylbenzene	10.87	105	125		N.D.	
71) trans-1,4-Dichloro-2-buten	0.00	53	0		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	281		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	148		N.D.	
76) n-Propylbenzene	11.26	120	60		N.D.	
77) 2-Chlorotoluene	11.47	91	195		N.D.	
78) 4-Chlorotoluene	11.47	91	195		N.D.	
79) 1,3,5-Trimethylbenzene	11.45	105	160		N.D.	
80) tert-Butylbenzene	11.76	119	220		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	292		N.D.	
82) sec-butylbenzene	11.98	105	327		N.D.	
83) 1,3-Dichlorobenzene	12.21	146	263		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.21	146	263		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	146		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	d
90) Hexachlorobutadiene	14.33	225	146		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	

**TIC FORMS**

SDG JPL33

VOLATILES ANALYSIS

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041907MVOWM1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: B041907MVOWM1

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419012.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419012.D Vial: 55  
Acq On : 19 Apr 2007 16:47 Operator: LH  
Sample : B041907MVOWM1 Inst : MOBY  
Misc : 5ml PFW+IS/SS (MV8-38-11) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419012.D 8260B.M Fri Apr 20 10:06:43 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-26-2

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-001

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419018.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/11/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419018.D Vial: 61  
Acq On : 19 Apr 2007 19:12 Operator: LH  
Sample : JPL33-001 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419018.D 8260B.M Fri Apr 20 10:09:56 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-26-1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-002

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419019.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/11/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419019.D Vial: 62  
Acq On : 19 Apr 2007 19:36 Operator: LH  
Sample : JPL33-002 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419019.D 8260B.M Fri Apr 20 10:11:06 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

EB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-003

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419020.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/11/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419020.D Vial: 63  
Acq On : 19 Apr 2007 20:00 Operator: LH  
Sample : JPL33-003 Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419020.D 8260B.M Fri Apr 20 10:13:37 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-11-4/10/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419013.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/11/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419013.D Vial: 56  
Acq On : 19 Apr 2007 17:11 Operator: LH  
Sample : JPL33-004 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419013.D 8260B.M Fri Apr 20 10:00:33 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-13

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-005  
 Lab File ID: M0419021.D  
 Date Collected: 04/12/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419021.D Vial: 64  
Acq On : 19 Apr 2007 20:24 Operator: LH  
Sample : JPL33-005 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419021.D 8260B.M Fri Apr 20 10:15:27 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-16
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Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-006  
 Lab File ID: M0419022.D  
 Date Collected: 04/12/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
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Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419022.D Vial: 65  
Acq On : 19 Apr 2007 20:48 Operator: LH  
Sample : JPL33-006 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419022.D 8260B.M Fri Apr 20 10:17:01 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-12-4/11/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-007

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419014.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/12/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
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30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419014.D Vial: 57  
Acq On : 19 Apr 2007 17:35 Operator: LH  
Sample : JPL33-007 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419014.D 8260B.M Fri Apr 20 10:03:12 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

DUPE-7-1Q07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-008  
 Lab File ID: M0419023.D  
 Date Collected: 04/12/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
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21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419023.D Vial: 66  
Acq On : 19 Apr 2007 21:12 Operator: LH  
Sample : JPL33-008 Inst : MOBY  
Misc : 5ml +IS/SS #4 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419023.D 8260B.M Fri Apr 20 10:18:40 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-7

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-009  
 Lab File ID: M0419024.D  
 Date Collected: 04/13/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
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24					
25					
26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419024.D Vial: 67  
Acq On : 19 Apr 2007 21:36 Operator: LH  
Sample : JPL33-009 Inst : MOBY  
Misc : 5ml +IS/SS #4 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419024.D 8260B.M Fri Apr 20 10:20:15 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-8

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-010  
 Lab File ID: M0419025.D  
 Date Collected: 04/13/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
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24					
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26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419025.D Vial: 68  
Acq On : 19 Apr 2007 22:00 Operator: LH  
Sample : JPL33-010 Inst : MOBY  
Misc : 5ml +IS/SS #4 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419025.D 8260B.M Fri Apr 20 10:21:44 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-13-4/12/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL33

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL33-011

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419015.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/13/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
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11					
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28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419015.D Vial: 58  
Acq On : 19 Apr 2007 18:00 Operator: LH  
Sample : JPL33-011 TB Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419015.D 8260B.M Fri Apr 20 10:04:55 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-5

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-012  
 Lab File ID: M0419026.D  
 Date Collected: 04/13/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
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28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419026.D Vial: 69  
Acq On : 19 Apr 2007 22:24 Operator: LH  
Sample : JPL33-012 Inst : MOBY  
Misc : 5ml +IS/SS #2 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419026.D 8260B.M Fri Apr 20 10:23:04 2007

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-10

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-013  
 Lab File ID: M0419027.D  
 Date Collected: 04/13/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
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26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419027.D Vial: 70  
Acq On : 19 Apr 2007 22:48 Operator: LH  
Sample : JPL33-013 Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419027.D 8260B.M Fri Apr 20 10:24:53 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-14-4/13/07

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL33  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL33-014  
 Lab File ID: M0419016.D  
 Date Collected: 04/13/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
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26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419016.D Vial: 59  
Acq On : 19 Apr 2007 18:24 Operator: LH  
Sample : JPL33-014 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419016.D 8260B.M Fri Apr 20 10:06:27 2007

**Metals Data**

**JPL33**

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-26-2	JPL33-001
MW-26-1	JPL33-002
EB-11-4/10/07	JPL33-003
MW-13	JPL33-005
MW-16	JPL33-006
DUPE-7-1Q07	JPL33-008
MW-7	JPL33-009
MW-8	JPL33-010
MW-5	JPL33-012
MW-10	JPL33-013
MW-10MS	JPL33-013MS
MW-10MSD	JPL33-013MSD

Were ICP interelement corrections applied? Yes/No YES

Were ICP background corrections applied? Yes/No NO

If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:

\_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Cheronne Ozeio

Date: 05/03/2007

Title: Metals Lead

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-26-2

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-001

Level (low/med): LOW

Date Received: 04/11/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	10.0			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-26-1

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-002

Level (low/med): LOW

Date Received: 04/11/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	9.66			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

EB-11-4/10/07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-003

Level (low/med): LOW

Date Received: 04/11/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	3.40			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-13

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-005

Level (low/med): LOW

Date Received: 04/12/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	70.3			M	R017204

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-16

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-006

Level (low/med): LOW

Date Received: 04/12/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	11.3			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-7-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-008

Level (low/med): LOW

Date Received: 04/12/2007

\* Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	10.5			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-7

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-009

Level (low/med): LOW

Date Received: 04/13/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	10.6			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-8

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-010

Level (low/med): LOW

Date Received: 04/13/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	12.7			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-5

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-012

Level (low/med): LOW

Date Received: 04/13/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.58			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-10

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL33

Matrix (soil/water): Water

Lab Sample ID: JPL33-013

Level (low/med): LOW

Date Received: 04/13/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	47.5			M	R017204

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL33**



## COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL33

Client Identification	Lab Sample Work Order Number
MW-26-2	JPL33-001DL
MW-26-1	JPL33-002DL
EB-11-4/10/07	JPL33-003
MW-13	JPL33-005
MW-16	JPL33-006
DUPE-7-1Q07	JPL33-008
DUPE-7-1Q07MS	JPL33-008MS
DUPE-7-1Q07MSD	JPL33-008MSD
MW-7	JPL33-009
MW-8	JPL33-010
MW-8MS	JPL33-010MS
MW-8MSD	JPL33-010MSD
MW-5	JPL33-012DL
MW-10	JPL33-013DL
MW-10MS	JPL33-013MS
MW-10MSD	JPL33-013MSD

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Bill Ambacher

Name: Bill Ambacher

Date: 5/4/07

Title: Inorganics/ Metals Manager

## **Inorganic Analysis Data Sheets**

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** MW-26-2 **Date/Time Collected:** 04/10/2007 07:57  
**Lab Sample ID:** JPL33-001 **Date/Time Received:** 04/11/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	2.0	U	2.0	0.28	05/02/2007	05/03/2007	R017239

Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL33  
Sample Number: MW-26-1 Date/Time Collected: 04/10/2007 08:40  
Lab Sample ID: JPL33-002 Date/Time Received: 04/11/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	05/02/2007	05/03/2007	R017239

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** EB-11-4/10/07 **Date/Time Collected:** 04/10/2007 08:24  
**Lab Sample ID:** JPL33-003 **Date/Time Received:** 04/11/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	1	1.0	U	1.0	0.14	05/02/2007	05/03/2007	R017239



**Laucks Testing Laboratories, Inc.**

**Final Results**

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** MW-16 **Date/Time Collected:** 04/11/2007 11:40  
**Lab Sample ID:** JPL33-006 **Date/Time Received:** 04/12/2007 08:30  
**Method:** E300.0 **Unit:** mg/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Nitrate - N	14797-55-8	10	4.4		2.0	0.55	04/12/2007	04/12/2007	R016780
Nitrite - N	14797-65-0	1	0.10	U	0.10	0.017	04/12/2007	04/12/2007	R016780
Sulfate as SO <sub>4</sub>	14808-79-8	10	26		10	1.7	04/12/2007	04/12/2007	R016780
Chloride	16887-00-6	10	28		10	0.76	04/12/2007	04/12/2007	R016780
Orthophosphate	7723-14-0	1	1.0	U	1.0	0.33	04/12/2007	04/12/2007	R016780

**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	50	1500		50	7.0	05/02/2007	05/03/2007	R017239

**Laucks Testing Laboratories, Inc.**

**Final Results**

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** DUPE-7-1Q07 **Date/Time Collected:** 04/11/2007 00:00  
**Lab Sample ID:** JPL33-008 **Date/Time Received:** 04/12/2007 08:30  
**Method:** E300.0 **Unit:** mg/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Nitrate - N	14797-55-8	10	4.8		2.0	0.55	04/12/2007	04/12/2007	R016780
Nitrite - N	14797-65-0	1	0.10	U	0.10	0.017	04/12/2007	04/12/2007	R016780
Sulfate as SO4	14808-79-8	10	26		10	1.7	04/12/2007	04/12/2007	R016780
Chloride	16887-00-6	10	31		10	0.76	04/12/2007	04/12/2007	R016780
Orthophosphate	7723-14-0	1	1.0	U	1.0	0.33	04/12/2007	04/12/2007	R016780

**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	50	1500		50	7.0	05/02/2007	05/03/2007	R017239







Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** MW-5 **Date/Time Collected:** 04/13/2007 09:24  
**Lab Sample ID:** JPL33-012 **Date/Time Received:** 04/16/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	2	16		2.0	0.28	05/02/2007	05/03/2007	R017239

Laucks Testing Laboratories, Inc.

Final Results

**Client:** Battelle **Project:** JPL Groundwater Monitoring  
**SDG Number:** JPL33  
**Sample Number:** MW-10 **Date/Time Collected:** 04/13/2007 12:00  
**Lab Sample ID:** JPL33-013 **Date/Time Received:** 04/16/2007 08:30  
**Method:** E314.0 **Unit:** ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	05/02/2007	05/03/2007	R017239

**SAMPLE DATA**

SDG JPL34

VOLATILES ANALYSIS

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-6

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL34  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL34-001  
 Lab File ID: M0419028.D  
 Date Collected: 04/16/2007  
 Date/Time Analyzed: 04/19/2007 23:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.84	
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.40	J
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.52	
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-6

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL34  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL34-001  
 Lab File ID: M0419028.D  
 Date Collected: 04/16/2007  
 Date/Time Analyzed: 04/19/2007 23:12  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	1.6	
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

MW-6

Lab Name: Laucks Testing Laboratories, Inc. Contract: JPL Groundwater Monitorin  
 SDG No.: JPL34 Run Sequence: R017025  
 Matrix: (SOIL/SED/WATER) Water Lab Sample ID: JPL34-001  
 Sample wt/vol: 5.00 (g/mL) mL Lab File ID: M0419028.D  
 Level: (LOW/MED) \_\_\_\_\_ Date Collected: 04/16/2007  
 % Moisture: not dec. \_\_\_\_\_ Date/Time Analyzed: 04/19/2007 23:12  
 GC Column: ZB-624 20m ID: 0.18 (mm) Dilution Factor: 1.0  
 Soil Extract Volume: \_\_\_\_\_ (uL) Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

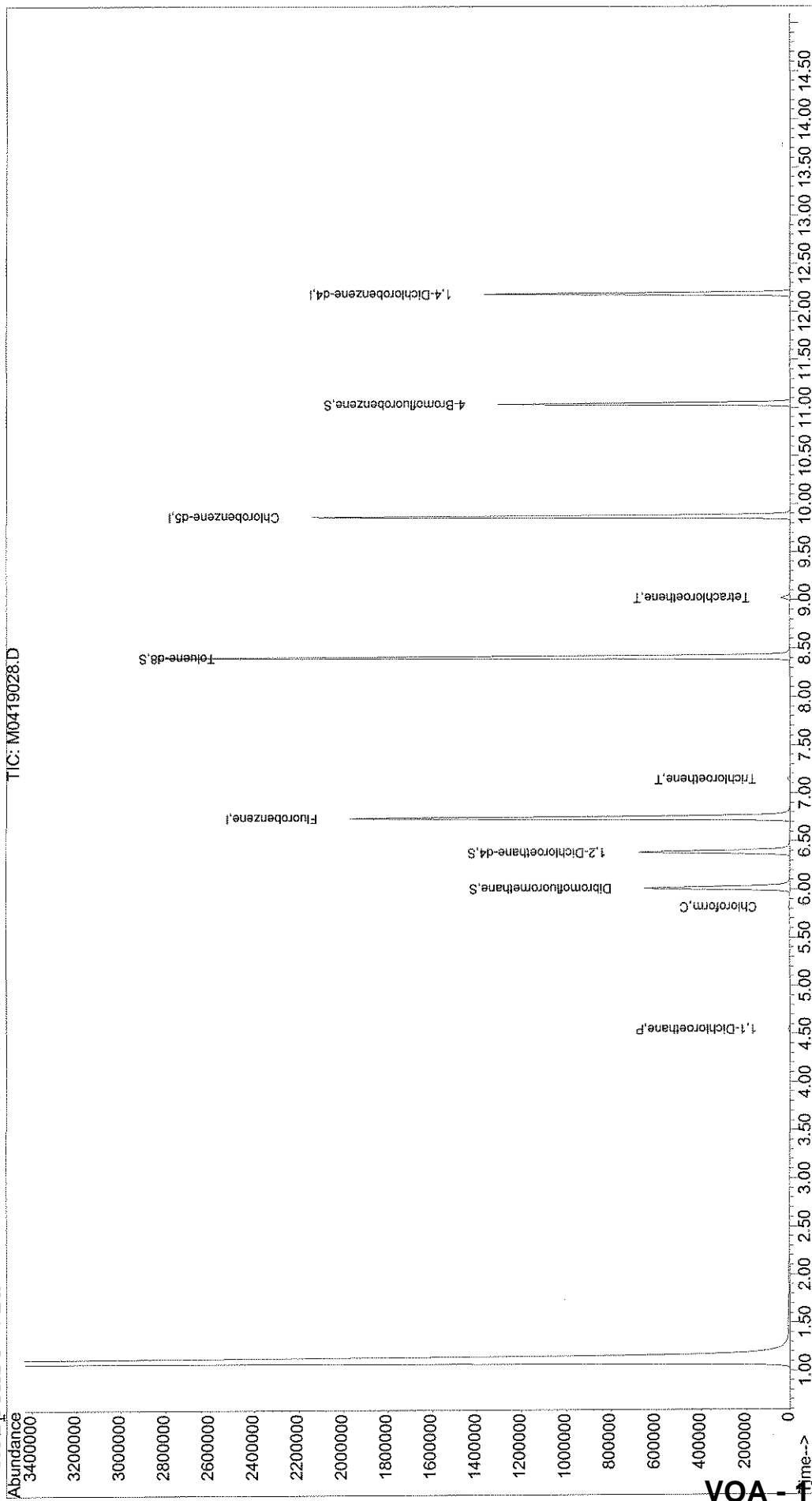
Comments:



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419028.D Vial: 71  
Acq On : 19 Apr 2007 23:12 Operator: LH  
Sample : JPL34-001 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:25 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



VOA - 17

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419028.D  
 Acq On : 19 Apr 2007 23:12  
 Sample : JPL34-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:25 2007

Vial: 71  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1  
 IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	1874949	50.00	ug/l	0.00 96.06%
50) Chlorobenzene-d5	9.87	82	637011	50.00	ug/l	0.00 90.57%
70) 1,4-Dichlorobenzene-d4	12.18	152	337964	50.00	ug/l	0.00 86.83%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	447867	50.63	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.26%
37) 1,2-Dichloroethane-d4	6.39	65	507705	50.14	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.28%
51) Toluene-d8	8.41	98	1843649	52.40	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.80%
72) 4-Bromofluorobenzene	11.04	95	386379	55.42	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.84%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0		N.D.	
3) Chloromethane	0.00	50	0		N.D.	
4) Vinyl Chloride	0.00	62	0		N.D.	
5) Bromomethane	1.79	96	55		N.D.	
6) Chloroethane	0.00	64	0		N.D.	
7) Trichlorofluoromethane	2.18	101	1243		N.D.	
8) Acrolein	0.00	56	0		N.D.	
9) 1,1-Dichloroethene	2.77	96	819		N.D.	
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0		N.D.	
11) Acetone	2.97	43	121		N.D.	
12) Iodomethane	0.00	142	0		N.D.	
13) Bromoethane	0.00	108	0		N.D.	
14) Carbon Disulfide	3.02	76	407		N.D.	
15) Allyl chloride	0.00	76	0		N.D.	
16) Acetonitrile	0.00	40	0		N.D.	d
17) Methyl Acetate	0.00	43	0		N.D.	
18) Methylene Chloride	0.00	84	0		N.D.	d
19) trans-1,2-Dichloroethene	0.00	96	0		N.D.	
20) Acrylonitrile	0.00	53	0		N.D.	
21) Methyl tert-butyl ether	3.93	73	57		N.D.	
22) 1,1-Dichloroethane	4.54	63	12099	0.84	ug/l	

(#) = qualifier out of range (m) = manual integration  
 M0419028.D 8260B.M Fri Apr 20 10:26:02 2007

96  
 W 4/20/07  
 VOA-18  
 Page 1

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419028.D  
 Acq On : 19 Apr 2007 23:12  
 Sample : JPL34-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:25 2007

Vial: 71  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.60	43	57		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	0.00	77	0		N.D.	
26) cis-1,2-Dichloroethene	0.00	96	0		N.D.	
27) 2-Butanone	5.45	43	67		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.75	41	58		N.D.	
31) Chloroform	5.81	83	6031	0.40	ug/l	84
32) 1,1,1-Trichloroethane	5.97	97	55		N.D.	
34) Cyclohexane	6.00	56	68		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	0.00	75	0		N.D.	
38) Benzene	6.40	78	1036		N.D.	
39) 1,2-Dichloroethane	0.00	62	0		N.D.	
40) Isobutanol	0.00	43	0		N.D. d	
41) Trichloroethene	7.15	130	4377	0.52	ug/l #	53
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	7.52	63	68		N.D.	
44) Dibromomethane	7.48	93	86		N.D.	
45) Methyl methacrylate	7.38	41	74		N.D.	
46) Bromodichloromethane	7.71	83	62		N.D.	
47) 2-Chloroethyl vinyl ether	0.00	63	0		N.D.	
48) cis-1,3-Dichloropropene	8.26	75	55		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D. d	
52) Toluene	8.48	92	76		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	0.00	69	0		N.D.	
55) 1,1,2-Trichloroethane	8.86	97	74		N.D.	
56) Tetrachloroethene	9.03	166	12069	1.58	ug/l	98
57) 1,3-Dichloropropane	0.00	76	0		N.D.	
58) 2-Hexanone	0.00	43	0		N.D. d	
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.91	112	127		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D. d	
63) 1,1,1,2-Tetrachloroethane	0.00	131	0		N.D.	
64) Ethylbenzene	9.99	91	137		N.D.	
65) m,p-Xylene	10.11	106	60		N.D.	
66) o-xylene	10.49	106	57		N.D.	
67) Styrene	0.00	104	0		N.D.	

*LH 4/20/07*

(#) = qualifier out of range (m) = manual integration  
 M0419028.D 8260B.M Fri Apr 20 10:26:03 2007

Quantitation Report

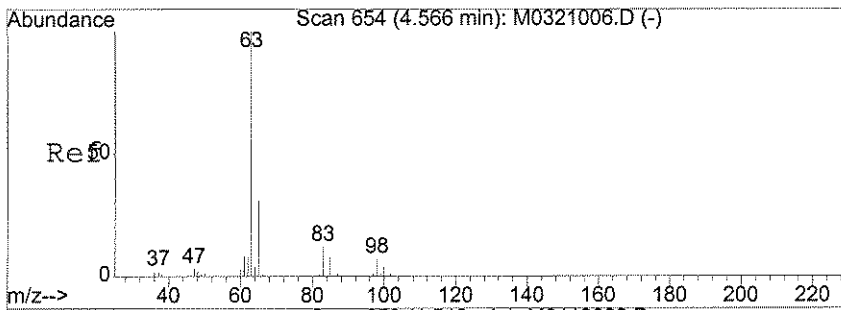
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 Acq On : 19 Apr 2007 23:12  
 Sample : JPL34-001  
 Misc : 5ml +IS/SS #3  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:25 2007

Vial: 71  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

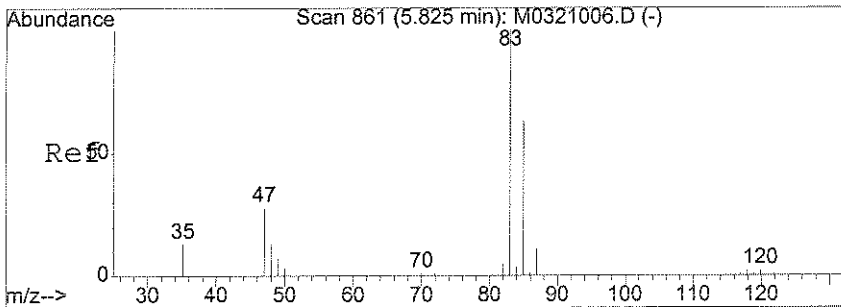
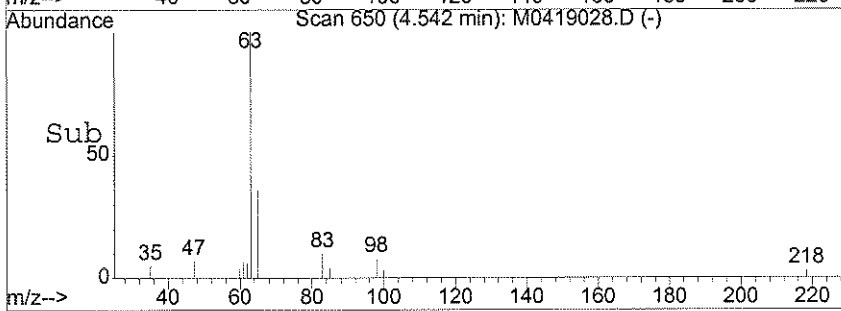
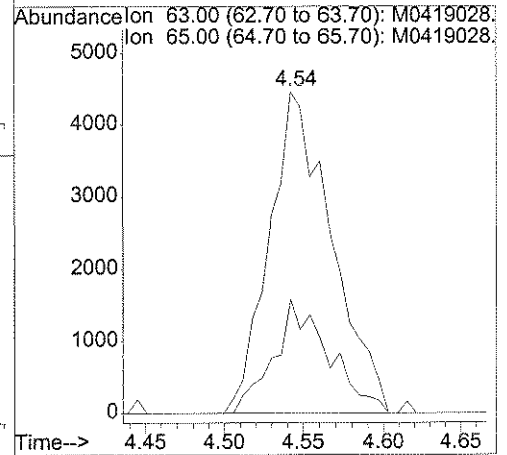
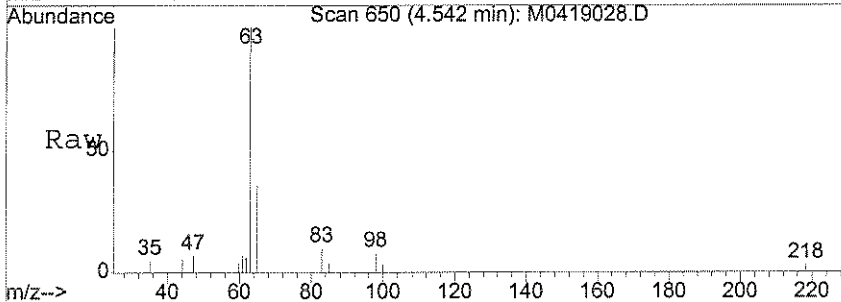
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1)

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.75	173	58		N.D.	
69) Isopropylbenzene	11.04	105	1096		N.D.	
71) trans-1,4-Dichloro-2-buten	11.12	53	55		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	131		N.D.	
75) 1,2,3-Trichloropropane	11.05	110	126		N.D.	
76) n-Propylbenzene	11.42	120	60		N.D.	
77) 2-Chlorotoluene	11.26	91	58		N.D.	
78) 4-Chlorotoluene	11.50	91	60		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	0.00	119	0		N.D.	
81) 1,2,4-Trimethylbenzene	11.83	105	61		N.D.	
82) sec-butylbenzene	11.99	105	243		N.D.	
83) 1,3-Dichlorobenzene	12.24	146	63		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.24	146	63		N.D.	
86) 1,2-Dichlorobenzene	12.58	146	66		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.14	225	75		N.D.	
91) Naphthalene	0.00	128	0		N.D.	
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	



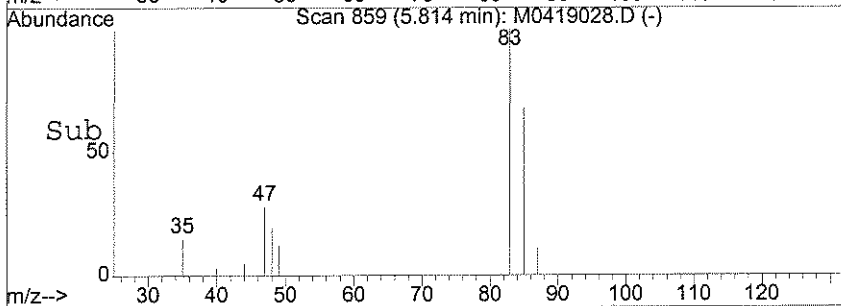
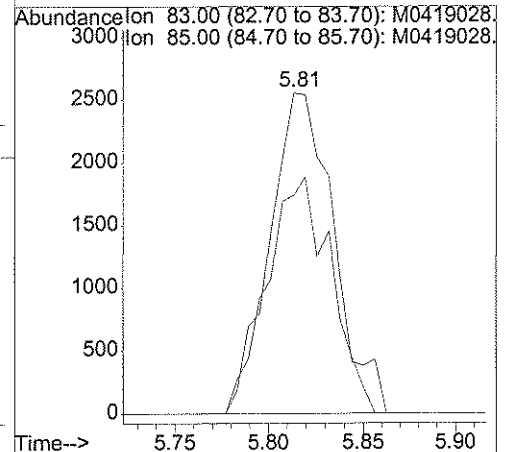
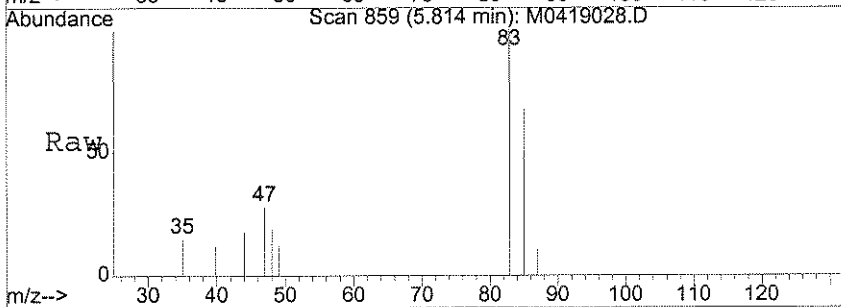
#22  
 1,1-Dichloroethane  
 Concen: 0.84 ug/l  
 RT: 4.54 min Scan# 650  
 Delta R.T. 0.00 min  
 Lab File: M0419028.D  
 Acq: 19 Apr 2007 23:12

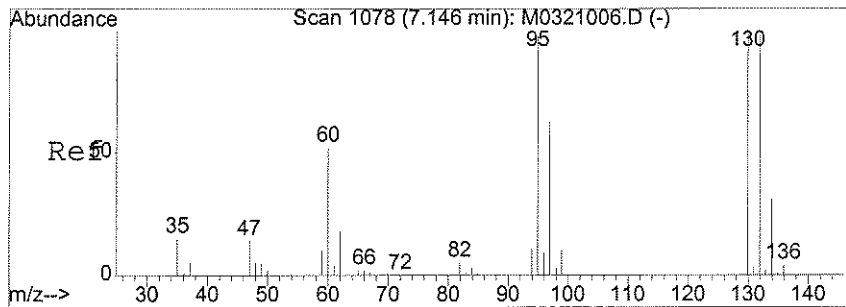
Tgt Ion: 63 Resp: 12099  
 Ion Ratio Lower Upper  
 63 100  
 65 31.3 13.7 53.7



#31  
 Chloroform  
 Concen: 0.40 ug/l  
 RT: 5.81 min Scan# 859  
 Delta R.T. 0.00 min  
 Lab File: M0419028.D  
 Acq: 19 Apr 2007 23:12

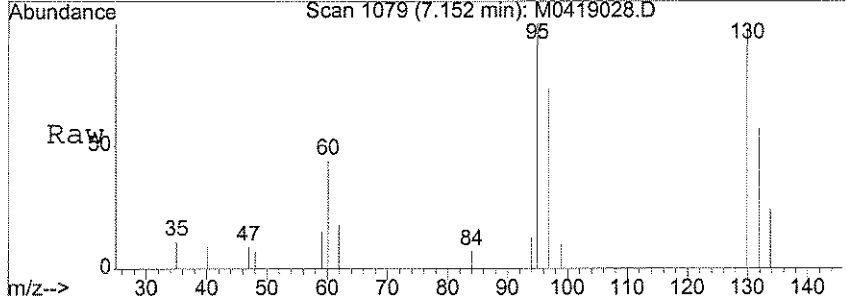
Tgt Ion: 83 Resp: 6031  
 Ion Ratio Lower Upper  
 83 100  
 85 73.4 41.2 81.2



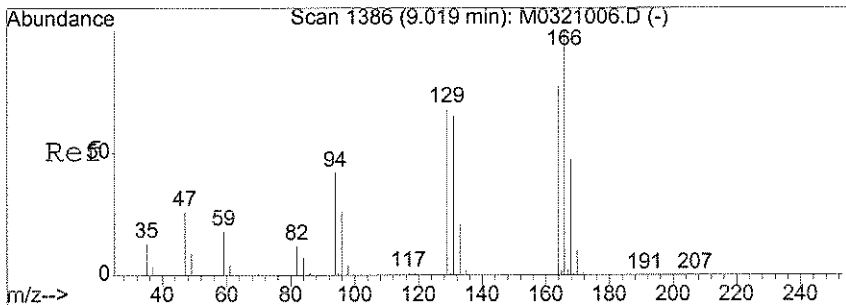
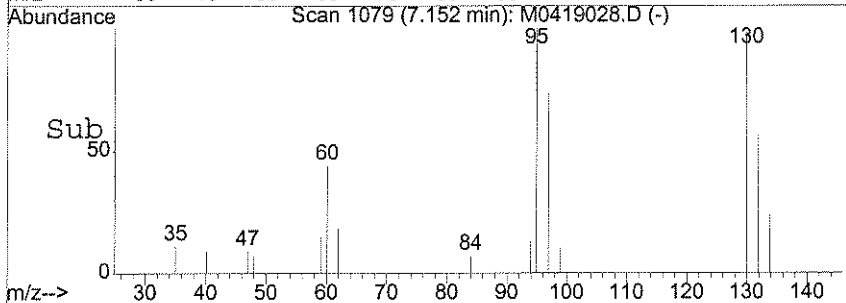
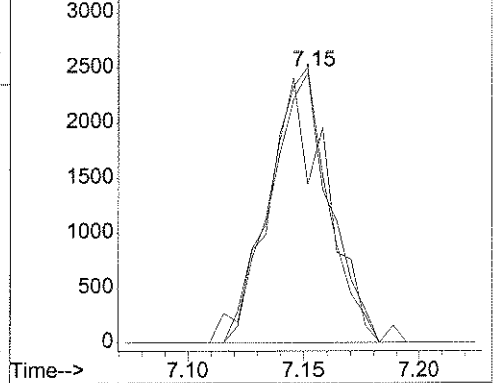


#41  
 Trichloroethene  
 Concen: 0.52 ug/l  
 RT: 7.15 min Scan# 1079  
 Delta R.T. 0.01 min  
 Lab File: M0419028.D  
 Acq: 19 Apr 2007 23:12

Tgt Ion	Resp	Lower	Upper
130	4377		
130	100		
132	96.6	75.0	115.0
95	0.0	69.4	109.4#

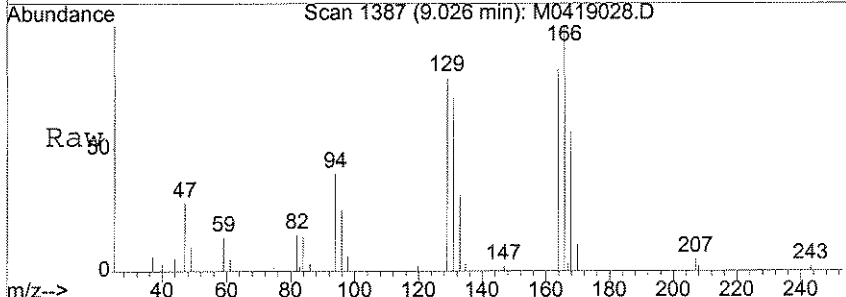


Abundance  
 Ion 130.00 (129.70 to 130.70): M04190  
 Ion 132.00 (131.70 to 132.70): M04190  
 Ion 95.00 (94.70 to 95.70): M0419028

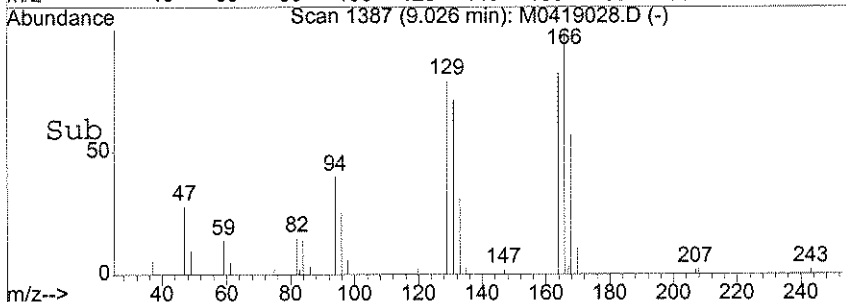
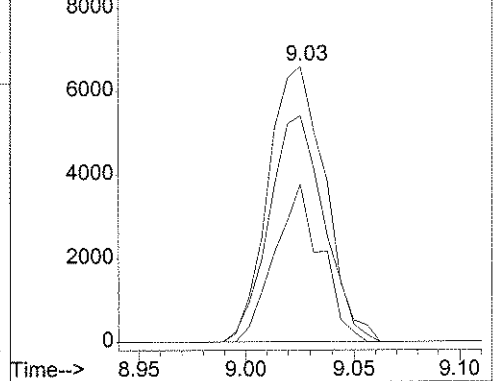


#56  
 Tetrachloroethene  
 Concen: 1.58 ug/l  
 RT: 9.03 min Scan# 1387  
 Delta R.T. 0.00 min  
 Lab File: M0419028.D  
 Acq: 19 Apr 2007 23:12

Tgt Ion	Resp	Lower	Upper
166	12069		
166	100		
164	79.6	63.3	94.9
168	46.8	39.6	59.4



Abundance  
 Ion 165.95 (165.65 to 166.65): M04190  
 Ion 163.95 (163.65 to 164.65): M04190  
 Ion 167.95 (167.65 to 168.65): M04190



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-15-4/16/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL34

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL34-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/16/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 18:48

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
75-71-8	Dichlorodifluoromethane	0.50	U
74-87-3	Chloromethane	0.50	U
75-01-4	Vinyl chloride	0.50	U
74-83-9	Bromomethane	0.50	U
75-00-3	Chloroethane	0.50	U
75-69-4	Trichlorofluoromethane	0.50	U
75-35-4	1,1-Dichloroethene	0.50	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.50	U
75-09-2	Methylene chloride	0.50	U
1634-04-4	Methyl tert-butyl ether	0.50	U
156-60-5	trans-1,2-Dichloroethene	0.50	U
75-34-3	1,1-Dichloroethane	0.50	U
594-20-7	2,2-Dichloropropane	0.50	U
156-59-2	cis-1,2-Dichloroethene	0.50	U
78-93-3	2-Butanone	5.0	U
74-97-5	Bromochloromethane	0.50	U
67-66-3	Chloroform	0.50	U
71-55-6	1,1,1-Trichloroethane	0.50	U
56-23-5	Carbon tetrachloride	0.50	U
563-58-6	1,1-Dichloropropene	0.50	U
71-43-2	Benzene	0.50	U
107-06-2	1,2-Dichloroethane	0.50	U
79-01-6	Trichloroethene	0.50	U
78-87-5	1,2-Dichloropropane	0.50	U
74-95-3	Dibromomethane	0.50	U
75-27-4	Bromodichloromethane	0.50	U
10061-01-	cis-1,3-Dichloropropene	0.50	U
108-10-1	4-Methyl-2-pentanone	5.0	U
108-88-3	Toluene	0.50	U

1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-15-4/16/07

Lab Name: Laucks Testing Laboratories, Inc.  
 SDG No.: JPL34  
 Matrix: (SOIL/SED/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_(uL)  
 Heated Purge: (Y/N) N

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL34-004  
 Lab File ID: M0419017.D  
 Date Collected: 04/16/2007  
 Date/Time Analyzed: 04/19/2007 18:48  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_(uL)

CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
10061-02-	trans-1,3-Dichloropropene	0.50	U
79-00-5	1,1,2-Trichloroethane	0.50	U
127-18-4	Tetrachloroethene	0.50	U
142-28-9	1,3-Dichloropropane	0.50	U
124-48-1	Dibromochloromethane	0.50	U
106-93-4	1,2-Dibromoethane	0.50	U
108-90-7	Chlorobenzene	0.50	U
100-41-4	Ethylbenzene	0.50	U
630-20-6	1,1,1,2-Tetrachloroethane	0.50	U
179601-23	m,p-Xylene	1.0	U
95-47-6	o-Xylene	0.50	U
100-42-5	Styrene	0.50	U
75-25-2	Bromoform	0.50	U
98-82-8	Isopropylbenzene	0.50	U
79-34-5	1,1,2,2-Tetrachloroethane	0.50	U
103-65-1	n-Propylbenzene	0.50	U
108-86-1	Bromobenzene	0.50	U
96-18-4	1,2,3-Trichloropropane	0.50	U
95-49-8	2-Chlorotoluene	0.50	U
108-67-8	1,3,5-Trimethylbenzene	0.50	U
106-43-4	4-Chlorotoluene	0.50	U
98-06-6	tert-Butylbenzene	0.50	U
95-63-6	1,2,4-Trimethylbenzene	0.50	U
135-98-8	sec-Butylbenzene	0.50	U
99-87-6	4-Isopropyltoluene	0.50	U
541-73-1	1,3-Dichlorobenzene	0.50	U
106-46-7	1,4-Dichlorobenzene	0.50	U
104-51-8	n-Butylbenzene	0.50	U
95-50-1	1,2-Dichlorobenzene	0.50	U



1  
VOLATILE ORGANICS ANALYSIS DATA SHEET

CLIENT SAMPLE NO.

TB-15-4/16/07

Lab Name: Laucks Testing Laboratories, Inc.

Contract: JPL Groundwater Monitorin

SDG No.: JPL34

Run Sequence: R017025

Matrix: (SOIL/SED/WATER) Water

Lab Sample ID: JPL34-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/16/2007

% Moisture: not dec. \_\_\_\_\_

Date/Time Analyzed: 04/19/2007 18:48

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Heated Purge: (Y/N) N

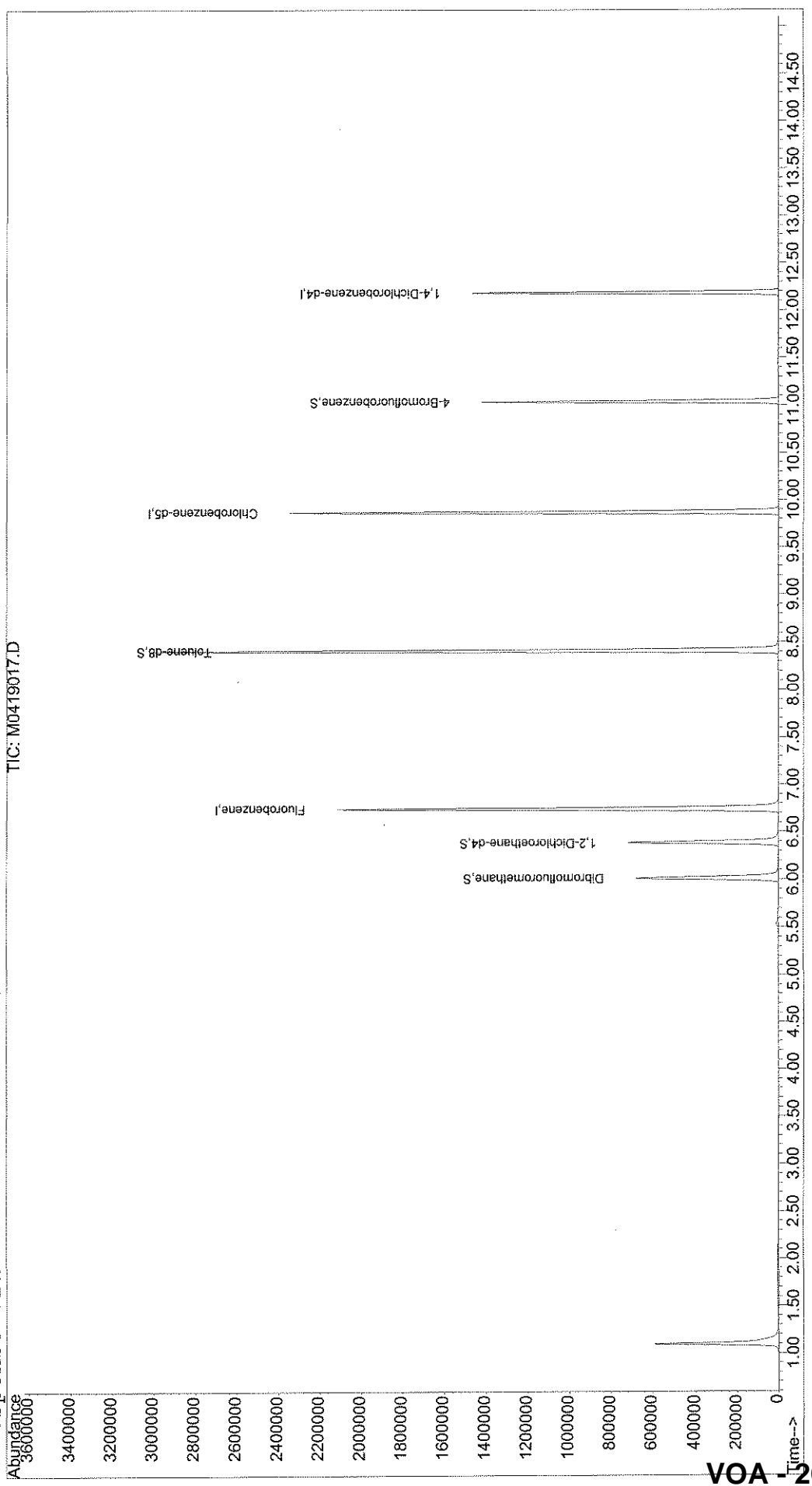
CAS NO.	COMPOUND	CONCENTRATION UNITS: (ug/L or ug/kg) <u>ug/L</u>	Q
96-12-8	1,2-Dibromo-3-chloropropane	0.50	U
120-82-1	1,2,4-Trichlorobenzene	0.50	U
87-68-3	Hexachlorobutadiene	0.50	U
91-20-3	Naphthalene	0.50	U
87-61-6	1,2,3-Trichlorobenzene	0.50	U

Comments:

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419017.D Vial: 60  
Acq On : 19 Apr 2007 18:48 Operator: LH  
Sample : JPL34-004 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: rteint.p  
Quant Time: Apr 20 10:07 2007 Quant Results File: 8260B.RES

Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Last Update : Tue Apr 17 16:33:44 2007  
Response via : Initial Calibration



Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419017.D  
 Acq On : 19 Apr 2007 18:48  
 Sample : JPL34-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:07 2007

Vial: 60  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

IS QA File : X:\MSVOA\MOBY\041607\M0416012.D (16 Apr 2007 12:11)

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min) Rcv(Ar )
1) Fluorobenzene	6.74	96	2020158	50.00	ug/l	0.00 103.50%
50) Chlorobenzene-d5	9.87	82	684983	50.00	ug/l	0.00 97.40%
70) 1,4-Dichlorobenzene-d4	12.19	152	366955	50.00	ug/l	0.00 94.28%

System Monitoring Compounds

33) Dibromofluoromethane	6.01	111	481368	50.51	ug/l	0.00
Spiked Amount	50.000	Range	85 - 115	Recovery	=	101.02%
37) 1,2-Dichloroethane-d4	6.38	65	545264	49.98	ug/l	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	99.96%
51) Toluene-d8	8.41	98	1983966	52.44	ug/l	0.00
Spiked Amount	50.000	Range	85 - 120	Recovery	=	104.88%
72) 4-Bromofluorobenzene	11.04	95	417528	55.16	ug/l	0.00
Spiked Amount	50.000	Range	75 - 120	Recovery	=	110.32%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) Dichlorodifluoromethane	0.00	85	0	N.D.		
3) Chloromethane	0.00	50	0	N.D.		
4) Vinyl Chloride	0.00	62	0	N.D.		
5) Bromomethane	0.00	96	0	N.D.		
6) Chloroethane	0.00	64	0	N.D.		
7) Trichlorofluoromethane	0.00	101	0	N.D.		
8) Acrolein	0.00	56	0	N.D.		
9) 1,1-Dichloroethene	2.84	96	57	N.D.		
10) 1,1,2-Trichloro-1,2,2-trif	0.00	101	0	N.D.		
11) Acetone	0.00	43	0	N.D.	d	
12) Iodomethane	0.00	142	0	N.D.		
13) Bromoethane	0.00	108	0	N.D.		
14) Carbon Disulfide	3.03	76	270	N.D.		
15) Allyl chloride	0.00	76	0	N.D.		
16) Acetonitrile	0.00	40	0	N.D.	d	
17) Methyl Acetate	0.00	43	0	N.D.		
18) Methylene Chloride	0.00	84	0	N.D.	d	
19) trans-1,2-Dichloroethene	3.95	96	65	N.D.		
20) Acrylonitrile	3.86	53	83	N.D.		
21) Methyl tert-butyl ether	0.00	73	0	N.D.		
22) 1,1-Dichloroethane	0.00	63	0	N.D.		

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419017.D  
 Acq On : 19 Apr 2007 18:48  
 Sample : JPL34-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:07 2007

Vial: 60  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
23) Vinyl acetate	4.71	43	55		N.D.	
24) Chloroprene	0.00	53	0		N.D.	
25) 2,2-Dichloropropane	5.40	77	56		N.D.	
26) cis-1,2-Dichloroethene	5.37	96	62		N.D.	
27) 2-Butanone	5.53	43	525		N.D.	
28) Propionitrile	0.00	54	0		N.D.	
29) Bromochloromethane	0.00	128	0		N.D.	
30) Methacrylonitrile	5.78	41	69		N.D.	
31) Chloroform	5.82	83	57		N.D.	
32) 1,1,1-Trichloroethane	0.00	97	0		N.D.	
34) Cyclohexane	6.01	56	131		N.D.	
35) Carbon Tetrachloride	0.00	117	0		N.D.	
36) 1,1-Dichloropropene	6.15	75	67		N.D.	
38) Benzene	6.40	78	769		N.D.	
39) 1,2-Dichloroethane	6.40	62	482		N.D.	
40) Isobutanol	0.00	43	0		N.D.	d
41) Trichloroethene	7.15	130	78		N.D.	
42) Methylcyclohexane	0.00	83	0		N.D.	
43) 1,2-Dichloropropane	0.00	63	0		N.D.	
44) Dibromomethane	0.00	93	0		N.D.	
45) Methyl methacrylate	0.00	41	0		N.D.	
46) Bromodichloromethane	0.00	83	0		N.D.	
47) 2-Chloroethyl vinyl ether	7.98	63	55		N.D.	
48) cis-1,3-Dichloropropene	0.00	75	0		N.D.	
49) 4-Methyl-2-pentanone	0.00	43	0		N.D.	d
52) Toluene	8.48	92	756		N.D.	
53) trans-1,3-Dichloropropene	0.00	75	0		N.D.	
54) Ethyl methacrylate	8.63	69	71		N.D.	
55) 1,1,2-Trichloroethane	8.98	97	102		N.D.	
56) Tetrachloroethene	0.00	166	0		N.D.	
57) 1,3-Dichloropropane	9.19	76	58		N.D.	
58) 2-Hexanone	0.00	43	0		N.D.	d
59) Dibromochloromethane	0.00	129	0		N.D.	
60) 1,2-Dibromoethane	0.00	107	0		N.D.	
61) Chlorobenzene	9.89	112	223		N.D.	
62) 1-Chlorohexane	0.00	91	0		N.D.	d
63) 1,1,1,2-Tetrachloroethane	9.86	131	72		N.D.	
64) Ethylbenzene	9.99	91	372		N.D.	
65) m,p-Xylene	10.11	106	431		N.D.	
66) o-xylene	0.00	106	0		N.D.	
67) Styrene	10.51	104	66		N.D.	

(#) = qualifier out of range (m) = manual integration  
 M0419017.D 8260B.M Fri Apr 20 10:08:00 2007

Quantitation Report

Data File : X:\MSVOA\MOBY\041907\M0419017.D  
 Acq On : 19 Apr 2007 18:48  
 Sample : JPL34-004 TB  
 Misc : 5ml +IS/SS #1  
 MS Integration Params: rteint.p  
 Quant Time: Apr 20 10:07 2007

Vial: 60  
 Operator: LH  
 Inst : MOBY  
 Multiplr: 1.00

Quant Results File: 8260B.RES

Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
 Title : VOA 8260- 5ML Calibration 5973M  
 Last Update : Tue Apr 17 16:33:44 2007  
 Response via : Initial Calibration  
 DataAcq Meth : 8260(0~1

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
68) Bromoform	10.73	173	127		N.D.	
69) Isopropylbenzene	10.86	105	286		N.D.	
71) trans-1,4-Dichloro-2-buten	11.04	53	73		N.D.	
73) Bromobenzene	0.00	156	0		N.D.	
74) 1,1,2,2-Tetrachloroethane	11.04	83	346		N.D.	
75) 1,2,3-Trichloropropane	11.04	110	353		N.D.	
76) n-Propylbenzene	11.28	120	96		N.D.	
77) 2-Chlorotoluene	11.37	91	71		N.D.	
78) 4-Chlorotoluene	11.48	91	316		N.D.	
79) 1,3,5-Trimethylbenzene	0.00	105	0		N.D.	
80) tert-Butylbenzene	11.77	119	61		N.D.	
81) 1,2,4-Trimethylbenzene	11.82	105	59		N.D.	
82) sec-butylbenzene	12.18	105	333		N.D.	
83) 1,3-Dichlorobenzene	12.09	146	56		N.D.	
84) 4-Isopropyltoluene	0.00	119	0		N.D.	d
85) 1,4-Dichlorobenzene	12.22	146	140		N.D.	
86) 1,2-Dichlorobenzene	0.00	146	0		N.D.	
87) n-Butylbenzene	0.00	91	0		N.D.	d
88) 1,2-Dibromo-3-chloropropan	0.00	157	0		N.D.	
89) 1,2,4-Trichlorobenzene	0.00	180	0		N.D.	
90) Hexachlorobutadiene	14.01	225	59		N.D.	
91) Naphthalene	0.00	128	0		N.D.	d
92) 1,2,3-Trichlorobenzene	0.00	180	0		N.D.	d

**TIC FORMS**

SDG JPL34

VOLATILES ANALYSIS

1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

B041907MVOWM1

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL34

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: B041907MVOWM1

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419012.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: \_\_\_\_\_

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
(ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419012.D Vial: 55  
Acq On : 19 Apr 2007 16:47 Operator: LH  
Sample : B041907MVOWM1 Inst : MOBY  
Misc : 5ml PFW+IS/SS (MV8-38-11) Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419012.D 8260B.M Mon Apr 30 16:42:27 2007



1 TIC  
VOLATILE ORGANICS ANALYSIS DATA SHEET  
TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

MW-6

Lab Name: Laucks Testing Laboratories, Inc  
 SDG No.: JPL34  
 Matrix: (SOIL/WATER) Water  
 Sample wt/vol: 5.00 (g/mL) mL  
 Level: (LOW/MED) \_\_\_\_\_  
 % Moisture: not dec. \_\_\_\_\_  
 GC Column: ZB-624 20m ID: 0.18 (mm)  
 Soil Extract Volume: \_\_\_\_\_ (uL)  
 Number TICs Found: 0

Contract: JPL Groundwater Monitorin  
 Run Sequence: R017025  
 Lab Sample ID: JPL34-001  
 Lab File ID: M0419028.D  
 Date Collected: 04/17/2007  
 Date Analyzed: 04/19/2007  
 Dilution Factor: 1.0  
 Soil Aliquot Volume: \_\_\_\_\_ (uL)  
 CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01				
02				
03				
04				
05				
06				
07				
08				
09				
10				
11				
12				
13				
14				
15				
16				
17				
18				
19				
20				
21				
22				
23				
24				
25				
26				
27				
28				
29				
30				

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419028.D Vial: 71  
Acq On : 19 Apr 2007 23:12 Operator: LH  
Sample : JPL34-001 Inst : MOBY  
Misc : 5ml +IS/SS #3 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419028.D 8260B.M Fri Apr 20 10:26:09 2007

1 TIC  
 VOLATILE ORGANICS ANALYSIS DATA SHEET  
 TENTATIVELY IDENTIFIED COMPOUNDS

CLIENT SAMPLE NO.

TB-15-4/16/07

Lab Name: Laucks Testing Laboratories, Inc

Contract: JPL Groundwater Monitorin

SDG No.: JPL34

Run Sequence: R017025

Matrix: (SOIL/WATER) Water

Lab Sample ID: JPL34-004

Sample wt/vol: 5.00 (g/mL) mL

Lab File ID: M0419017.D

Level: (LOW/MED) \_\_\_\_\_

Date Collected: 04/17/2007

% Moisture: not dec. \_\_\_\_\_

Date Analyzed: 04/19/2007

GC Column: ZB-624 20m ID: 0.18 (mm)

Dilution Factor: 1.0

Soil Extract Volume: \_\_\_\_\_ (uL)

Soil Aliquot Volume: \_\_\_\_\_ (uL)

Number TICs Found: 0

CONCENTRATION UNITS:  
 (ug/L or ug/kg) ug/L

	CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
01					
02					
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

Comments:

Library Search Compound Report

Data File : X:\MSVOA\MOBY\041907\M0419017.D Vial: 60  
Acq On : 19 Apr 2007 18:48 Operator: LH  
Sample : JPL34-004 TB Inst : MOBY  
Misc : 5ml +IS/SS #1 Multiplr: 1.00  
MS Integration Params: LSCINT.P  
Quant Method : X:\MSVOA\MOBY\QUANT\8260B.M (RTE Integrator)  
Title : VOA 8260- 5ML Calibration 5973M  
Library : D:\DATABASE\NIST129K.L

No Library Search Compounds Detected

\*\*\*\*\*  
M0419017.D 8260B.M Fri Apr 20 10:08:05 2007

**Metals Data**

**JPL34**

COVER PAGE-INORGANIC ANALYSES DATA PACKAGE

Lab Name: Laucks Laboratories Contract: JPL Groundwater Monitorin  
 Lab Code: LAUCKS SDG No.: JPL34  
 SOW No.: \_\_\_\_\_

Sample No.	Lab Sample ID
MW-6	JPL34-001
MW-6MS	JPL34-001MS
MW-6MSD	JPL34-001MSD
MW-15	JPL34-002
DUPE-8-1Q07	JPL34-003
DUPE-8-1Q07MS	JPL34-003MS
DUPE-8-1Q07MSD	JPL34-003MSD

Were ICP interelement corrections applied? Yes/No YES  
 Were ICP background corrections applied? Yes/No NO  
 If yes-was raw data generated before application of background corrections? Yes/No NO

Comments:  
 \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

I certify that this data package is technically complete, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: *Bill Ambacher* Name: Bill Ambacher  
 Date: 5/7/07 Title: Inorganics/Metals Manager

## **Metals Analysis Data Sheets**

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-6

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL34

Matrix (soil/water): Water

Lab Sample ID: JPL34-001

Level (low/med): LOW

Date Received: 04/17/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	10.1			M	R017204

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_



INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

MW-15

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL34

Matrix (soil/water): Water

Lab Sample ID: JPL34-002

Level (low/med): LOW

Date Received: 04/17/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.45			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_

INORGANIC ANALYSES DATA SHEET

SAMPLE NO.

DUPE-8-1Q07

Lab Name: Laucks Laboratories

Contract: JPL Groundwater Monitorin

Lab Code: LAUCKS

SDG No.: JPL34

Matrix (soil/water): Water

Lab Sample ID: JPL34-003

Level (low/med): LOW

Date Received: 04/17/2007

% Solids: \_\_\_\_\_

Concentration Units : ug/L

CAS No.	Analyte	Concentration	C	Q	M	Run Seq.
7440-47-3	Chromium	8.05			M	R017069

Color Before: \_\_\_\_\_ Clarity Before: \_\_\_\_\_ Texture: \_\_\_\_\_

Color After: \_\_\_\_\_ Clarity After: \_\_\_\_\_ Artifacts: No

Comment \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

**Miscellaneous Inorganic Data**

**JPL34**

COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

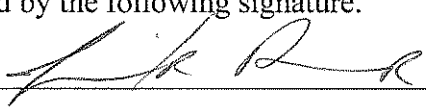
Lab Name: LAUCKS TESTING LABS, INC.

Lab Code: LAUCKS

SDG No.: JPL34

Client Identification	Lab Sample Work Order Number
MW-6	JPL34-001DL
MW-6MS	JPL34-001MS
MW-6MSD	JPL34-001MSD

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: 

Name: Jennifer Penner

Date: 5-7-07

Title: Inorganics Lead

## **Inorganic Analysis Data Sheets**

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Laucks Testing Laboratories, Inc.

Final Results

Client: Battelle Project: JPL Groundwater Monitoring  
SDG Number: JPL34  
Sample Number: MW-6 Date/Time Collected: 04/16/2007 10:04  
Lab Sample ID: JPL34-001 Date/Time Received: 04/17/2007 08:30  
Method: E314.0 Unit: ug/L

Analyte	CAS	DF	Result	Q	PQL	MDL	Prepared	Analyzed	Run Seq.
Perchlorate	14797-73-0	4	4.0	U	4.0	0.56	05/02/2007	05/03/2007	R017239

**DIVIDER SHEET**

**ANALYTICAL DATA**  
**FOR**

**Hexavalent Chromium**

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**ANALYSIS**

**COLUMBIA ANALYTICAL SERVICES, INC.**

Analytical Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

**Service Request :** P0700303  
**Date Collected :** 03/27-30/07  
**Date Received :** 03/27-30/07

Chromium, Hexavalent

**Prep Method :** None  
**Analysis Method :** 7196A  
**Test Notes :**

**Units :** mg/L (ppm)  
**Basis :** NA

Sample Name	Lab Code	PQL	MDL	Dilution Factor	Date Extracted	Date/Time Analyzed	Result	Result Notes
MW-21-5	P0700303-001	0.01	0.004	1	NA	03/27/07 15:10	ND	
MW-21-4	P0700303-002	0.01	0.004	1	NA	03/27/07 15:10	ND	
MW-21-3	P0700303-003	0.01	0.004	1	NA	03/27/07 15:10	ND	
MW-21-2	P0700303-004	0.01	0.004	1	NA	03/27/07 15:10	ND	
MW-21-1	P0700303-005	0.01	0.004	1	NA	03/27/07 15:10	ND	
EB-1-3/27/07	P0700303-006	0.01	0.004	1	NA	03/27/07 15:10	ND	
MW-14-3	P0700303-007	0.01	0.004	1	NA	03/28/07 15:05	ND	
MW-14-2	P0700303-008	0.01	0.004	1	NA	03/28/07 15:05	ND	
MW-14-1	P0700303-009	0.01	0.004	1	NA	03/28/07 15:05	ND	
EB-2-3/28/07	P0700303-010	0.01	0.004	1	NA	03/28/07 15:05	ND	
MW-17-4	P0700303-011	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-17-3	P0700303-012	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-17-2	P0700303-013	0.01	0.004	1	NA	03/29/07 15:50	ND	
DUPE-1-1Q07	P0700303-014	0.01	0.004	1	NA	03/29/07 15:50	ND	
EB-3-3/29/07	P0700303-015	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-18-4	P0700303-016	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-18-3	P0700303-017	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-18-2	P0700303-018	0.01	0.004	1	NA	03/29/07 15:50	ND	
DUPE-2-1Q07	P0700303-019	0.01	0.004	1	NA	03/29/07 15:50	ND	
MW-20-5	P0700303-020	0.01	0.004	1	NA	03/30/07 14:50	ND	
MW-20-4	P0700303-021	0.01	0.004	1	NA	03/30/07 14:50	ND	
MW-20-3	P0700303-022	0.01	0.004	1	NA	03/30/07 14:50	ND	
MW-20-2	P0700303-023	0.01	0.004	1	NA	03/30/07 14:50	ND	
MW-20-1	P0700303-024	0.01	0.004	1	NA	03/30/07 14:50	ND	
DUPE-3-1Q07	P0700303-025	0.01	0.004	1	NA	03/30/07 14:50	ND	
EB-4-3/30/07	P0700303-026	0.01	0.004	1	NA	03/30/07 14:50	ND	
Method Blank	P0700303-MB	0.01	0.004	1	NA	03/27/07 15:10	ND	
Method Blank	P0700303-MB	0.01	0.004	1	NA	03/28/07 15:05	ND	
Method Blank	P0700303-MB	0.01	0.004	1	NA	03/29/07 15:50	ND	
Method Blank	P0700303-MB	0.01	0.004	1	NA	03/30/07 14:50	ND	

Approved By

  
 18

Date :

4/4/07



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/27/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By: \_\_\_\_\_  
ICCBMDL/120594



Date: 4/4/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

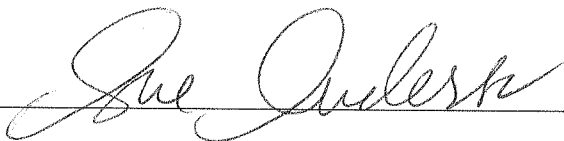
**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/27/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0387	0.0404	104	90-110
CCV1	0.0387	0.0379	98	90-110

Approved By:



Date:

4/4/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

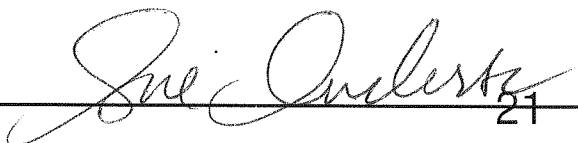
**Service Request :** P0700303  
**Date Collected :** NA  
**Date Received :** NA  
**Date Extracted :** NA  
**Date Analyzed :** 03/27/07

Laboratory Control Sample Summary  
 Inorganic Parameters

**Sample Name :** Laboratory Control Sample  
**Lab Code :** P0700303-LCS  
**Test Notes :**

**Units :** mg/L (ppm)  
**Basis :** NA

Analyte	Prep Method	Analysis Method	True Value	Result	Percent Recovery	CAS	Result Notes
						Percent Recovery Acceptance Limits	
Chromium, Hexavalent	None	7196A	0.0400	0.0387	97	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0402	101	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0385	96	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0389	97	90-110	

Approved By  Date : 4/4/07

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

**Service Request :** P0700303  
**Date Collected :** 03/27/07  
**Date Received :** 03/27/07  
**Date Extracted :** NA  
**Date Analyzed :** 03/27/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-21-1 Units : mg/L (ppm)  
 Lab Code : P0700303-005MS P0700303-005DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0488	0.0522	98	104	85-115	7	

Approved By

  
 \_\_\_\_\_  
 22

Date :

4/4/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/28/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/4/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/28/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0387	0.0385	99	90-110
CCV1	0.0387	0.0376	97	90-110

Approved By:  
CCV1A/120594



Date:

4/4/07

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report


**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

**Service Request :** P0700303  
**Date Collected :** 03/28/07  
**Date Received :** 03/28/07  
**Date Extracted :** NA  
**Date Analyzed :** 03/28/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-14-3 Units : mg/L (ppm)  
 Lab Code : P0700303-007MS P0700303-007DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0550	0.0498	110	100	85-115	10	

Approved By  25 Date : 4/4/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report


**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/29/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND
CCB2	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/4/07

ICCBMDL/120594



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/29/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0387	0.0402	104	90-110
CCV1	0.0387	0.0385	99	90-110
CCV2	0.0387	0.0394	102	90-110

Approved By:



Date:

4/4/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report


**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

**Service Request :** P0700303  
**Date Collected :** 03/29/07  
**Date Received :** 03/29/07  
**Date Extracted :** NA  
**Date Analyzed :** 03/29/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-17-4 Units : mg/L (ppm)  
 Lab Code : P0700303-011MS P0700303-011DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0504	0.0487	101	97	85-115	3	

Approved By  Date : 4/4/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/30/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND
CCB2	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/4/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07

**Service Request:** P0700303  
**Date Analyzed:** 03/30/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0387	0.0380	98	90-110
CCV1	0.0387	0.0380	98	90-110
CCV2	0.0387	0.0380	98	90-110

Approved By:



Date:

4/4/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report


**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** NA  
**Sample Matrix :** WATER

**Service Request :** P0700303  
**Date Collected :** 03/30/07  
**Date Received :** 03/30/07  
**Date Extracted :** NA  
**Date Analyzed :** 03/30/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-20-3 Units : mg/L (ppm)  
 Lab Code : P0700303-022MS P0700303-022DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0496	0.0487	99	97	85-115	2	

Approved By  Date : 4/4/07

**DIVIDER SHEET**

**ANALYTICAL DATA**

**FOR**

**Hexavalent Chromium**

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**ANALYSIS**

**COLUMBIA ANALYTICAL SERVICES, INC.**

Analytical Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700321  
**Date Collected :** 04/02-06/07  
**Date Received :** 04/02-06/07


Chromium, Hexavalent

**Prep Method :** None  
**Analysis Method :** 7196A  
**Test Notes :**

**Units :** mg/L (ppm)  
**Basis :** NA

Sample Name	Lab Code	MRL	MDL	Dilution Factor	Date Extracted	Date/Time Analyzed	Result	Result Notes
MW-3-4	P0700321-001	0.01	0.004	1	NA	04/02/07 17:10	ND	
MW-3-3	P0700321-002	0.01	0.004	1	NA	04/02/07 17:10	ND	
MW-3-2	P0700321-003	0.01	0.004	1	NA	04/02/07 17:10	ND	
EB-5-4/2/07	P0700321-004	0.01	0.004	1	NA	04/02/07 17:10	ND	
MW-4-3	P0700321-005	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-4-2	P0700321-006	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-4-1	P0700321-007	0.01	0.004	1	NA	04/03/07 16:25	ND	
EB-6-4/3/07	P0700321-008	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-11-3	P0700321-009	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-11-2	P0700321-010	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-11-1	P0700321-011	0.01	0.004	1	NA	04/03/07 16:25	ND	
MW-23-4	P0700321-012	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-23-3	P0700321-013	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-23-2	P0700321-014	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-23-1	P0700321-015	0.01	0.004	1	NA	04/04/07 16:50	ND	
EB-7-4/4/07	P0700321-016	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-12-3	P0700321-017	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-12-2	P0700321-018	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-12-1	P0700321-019	0.01	0.004	1	NA	04/04/07 16:50	ND	
MW-24-4	P0700321-020	0.01	0.004	1	NA	04/05/07 17:55	0.006	J
MW-24-3	P0700321-021	0.01	0.004	1	NA	04/05/07 17:55	ND	
MW-24-2	P0700321-022	0.01	0.004	1	NA	04/05/07 17:55	ND	
MW-24-1	P0700321-023	0.01	0.004	1	NA	04/05/07 17:55	0.015	
DUPE-5-1Q07	P0700321-024	0.01	0.004	1	NA	04/05/07 17:55	ND	
EB-8-4/5/07	P0700321-025	0.01	0.004	1	NA	04/05/07 17:55	ND	
MW-22-3	P0700321-026	0.01	0.004	1	NA	04/06/07 13:45	ND	
MW-22-2	P0700321-027	0.01	0.004	1	NA	04/06/07 13:45	ND	
MW-22-1	P0700321-028	0.01	0.004	1	NA	04/06/07 13:45	ND	
DUPE-6-1Q07	P0700321-029	0.01	0.004	1	NA	04/06/07 13:45	ND	
EB-9-4/6/07	P0700321-030	0.01	0.004	1	NA	04/06/07 13:45	ND	

J Estimated concentration. The result is less than the PQL but greater than the MDL.

Approved By  **21** Date : 4/16/07

COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Report

Client : Battelle  
Project Name : JPL Groundwater Monitoring 1Q07  
Project Number : G486090  
Sample Matrix : WATER


Service Request : P0700321  
Date Collected : 04/02-06/07  
Date Received : 04/02-06/07

Chromium, Hexavalent

Prep Method : None  
Analysis Method : 7196A  
Test Notes :

Units : mg/L (ppm)  
Basis : NA

Sample Name	Lab Code	MRL	MDL	Dilution Factor	Date Extracted	Date/Time Analyzed	Result	Result Notes
Method Blank	P0700321-MB	0.01	0.004	1	NA	04/02/07 17:10	ND	
Method Blank	P0700321-MB	0.01	0.004	1	NA	04/03/07 16:25	ND	
Method Blank	P0700321-MB	0.01	0.004	1	NA	04/04/07 16:50	ND	
Method Blank	P0700321-MB	0.01	0.004	1	NA	04/05/07 17:55	ND	
Method Blank	P0700321-MB	0.01	0.004	1	NA	04/06/07 13:45	ND	

Approved By  **22** Date : 4/16/07



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/02/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/02/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0292	97	90-110
CCV1	0.0301	0.0292	97	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

CCV1A/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client : Battelle  
 Project Name : JPL Groundwater Monitoring 1Q07  
 Project Number : G486090  
 Sample Matrix : WATER

Service Request : P0700321  
 Date Collected : NA  
 Date Received : NA  
 Date Extracted : NA  
 Date Analyzed : 04/02/07

Laboratory Control Sample Summary  
 Inorganic Parameters

Sample Name : Laboratory Control Sample  
 Lab Code : P0700321-LCS  
 Test Notes :

Units : mg/L (ppm)  
 Basis : NA

Analyte	Prep Method	Analysis Method	True Value	Result	Percent Recovery	CAS Percent Recovery Acceptance Limits	Result Notes
Chromium, Hexavalent	None	7196A	0.0400	0.0378	95	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0405	101	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0400	100	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0389	97	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0425	106	90-110	

Approved By



25

Date :

4/16/07

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700321  
**Date Collected :** 04/02/07  
**Date Received :** 04/02/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/02/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-3-3 Units : mg/L (ppm)  
 Lab Code : P0700321-002MS P0700321-002DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0465	0.0465	93	93	85-115	<1	

Approved By \_\_\_\_\_

*Sue Janderts* **26**

Date : \_\_\_\_\_

*4/16/07*

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/03/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND
CCB2	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/03/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0294	98	90-110
CCV1	0.0301	0.0284	94	90-110
CCV2	0.0301	0.0294	98	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700321  
**Date Collected :** 04/03/07  
**Date Received :** 04/03/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/03/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-4-3 Units : mg/L (ppm)  
 Lab Code : P0700321-005MS P0700321-005DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0456	0.0476	91	95	85-115	4	

Approved By \_\_\_\_\_

*Jane Anderson* **29** Date : 4/16/07

Report By:RWONG

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700321  
**Date Collected :** 04/03/07  
**Date Received :** 04/03/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/03/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-11-3 Units : mg/L (ppm)  
 Lab Code : P0700321-009MS P0700321-009DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0486	0.0496	97	99	85-115	2	

Approved By \_\_\_\_\_

*[Signature]* **30**

Date : \_\_\_\_\_

*4/16/07*



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/04/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND
CCB2	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

ICCBMDL120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/04/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0293	97	90-110
CCV1	0.0301	0.0293	97	90-110
CCV2	0.0301	0.0293	97	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700321  
**Date Collected :** 04/04/07  
**Date Received :** 04/04/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/04/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-23-3 Units : mg/L (ppm)  
 Lab Code : P0700321-013MS P0700321-013DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0507	0.0507	101	101	85-115	<1	

Approved By



**33**

Date :

4/16/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/05/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND
CCB2	0.01	0.004	ND
CCB3	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

ICCBMDL120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/05/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0317	105	90-110
CCV1	0.0301	0.0308	102	90-110
CCV2	0.0301	0.0272	90	90-110
CCV3	0.0301	0.0272	90	90-110

Approved By:



Date:

4/16/07

CCV1A/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client : Battelle  
Project Name : JPL Groundwater Monitoring 1Q07  
Project Number : G486090  
Sample Matrix : WATER

Service Request : P0700321  
Date Collected : 04/05/07  
Date Received : 04/05/07  
Date Extracted : NA  
Date Analyzed : 04/05/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-24-4 Units : mg/L (ppm)  
Lab Code : P0700321-020MS P0700321-020DMS Basis : NA  
Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	0.00553	0.0479	0.0470	85	83	85-115	2	E1

E1 See case narrative.

Approved By  **36** Date : 4/16/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/06/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By: \_\_\_\_\_  
ICCBMDL/120594



Date: \_\_\_\_\_

4/16/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report


**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700321  
**Date Analyzed:** 04/06/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0324	108	90-110
CCV1	0.0301	0.0315	105	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/16/07

CCV1A/120594



QA/QC Report


Client : Battelle  
 Project Name : JPL Groundwater Monitoring 1Q07  
 Project Number : G486090  
 Sample Matrix : WATER

Service Request : P0700321  
 Date Collected : 04/06/07  
 Date Received : 04/06/07  
 Date Extracted : NA  
 Date Analyzed : 04/06/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-22-3 Units : mg/L (ppm)  
 Lab Code : P0700321-026MS P0700321-026DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	MRL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0462	0.0490	92	98	85-115	6	

Approved By  **39** Date : 4/16/07

**DIVIDER SHEET**

**ANALYTICAL DATA**  
**FOR**

**Hexavalent Chromium**

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**ANALYSIS**

**COLUMBIA ANALYTICAL SERVICES, INC.**

Analytical Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** 04/09-16/07  
**Date Received :** 04/09-16/07

Chromium, Hexavalent

Prep Method : None  
 Analysis Method : 7196A  
 Test Notes :

Units : mg/L (ppm)  
 Basis : NA

Sample Name	Lab Code	PQL	MDL	Dilution Factor	Date Extracted	Date/Time Analyzed	Result	Result Notes
MW-25-5	P0700347-001	0.01	0.004	1	NA	04/09/07 14:25	ND	
MW-25-4	P0700347-002	0.01	0.004	1	NA	04/09/07 14:25	ND	
MW-25-3	P0700347-003	0.01	0.004	1	NA	04/09/07 14:25	ND	
MW-25-2	P0700347-004	0.01	0.004	1	NA	04/09/07 14:25	ND	
MW-25-1	P0700347-005	0.01	0.004	1	NA	04/09/07 14:25	ND	
EB-10-4/9/07	P0700347-006	0.01	0.004	1	NA	04/09/07 14:25	ND	
MW-26-2	P0700347-007	0.01	0.004	1	NA	04/10/07 14:10	ND	
MW-26-1	P0700347-008	0.01	0.004	1	NA	04/10/07 14:10	ND	
EB-11-4/10/07	P0700347-009	0.01	0.004	1	NA	04/10/07 14:10	ND	
MW-13	P0700347-010	0.01	0.004	1	NA	04/11/07 17:45	0.041	
MW-16	P0700347-011	0.01	0.004	1	NA	04/11/07 17:45	ND	
DUPE-7-1Q07	P0700347-012	0.01	0.004	1	NA	04/11/07 17:45	ND	
MW-7	P0700347-013	0.01	0.004	1	NA	04/12/07 14:45	0.005	J
MW-8	P0700347-014	0.01	0.004	1	NA	04/12/07 14:45	ND	
MW-5	P0700347-015	0.01	0.004	1	NA	04/13/07 15:30	ND	
MW-10	P0700347-016	0.01	0.004	1	NA	04/13/07 15:30	ND	
MW-6	P0700347-017	0.01	0.004	1	NA	04/16/07 15:40	ND	
MW-15	P0700347-018	0.01	0.004	1	NA	04/16/07 15:40	ND	
DUPE-8-1Q07	P0700347-019	0.01	0.004	1	NA	04/16/07 15:40	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/09/07 14:25	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/10/07 14:10	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/11/07 17:45	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/12/07 14:45	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/13/07 15:30	ND	
Method Blank	P0700347-MB	0.01	0.004	1	NA	04/16/07 15:40	ND	

J Estimated concentration. The result is less than the PQL but greater than the MDL.

Approved By  21

Date : 4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/09/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By:



Date:

4/27/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/09/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0291	97	90-110
CCV1	0.0301	0.0291	97	90-110

Approved By:



Date:

4/27/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** NA  
**Date Received :** NA  
**Date Extracted :** NA  
**Date Analyzed :** 04/09/07


Laboratory Control Sample Summary  
 Inorganic Parameters

**Sample Name :** Laboratory Control Sample  
**Lab Code :** P0700347-LCS  
**Test Notes :**

**Units :** mg/L (ppm)  
**Basis :** NA

Analyte	Prep Method	Analysis Method	True Value	Result	Percent Recovery	CAS	Result Notes
						Percent Recovery Acceptance Limits	
Chromium, Hexavalent	None	7196A	0.0400	0.0374	94	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0390	98	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0419	105	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0393	98	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0369	92	90-110	
Chromium, Hexavalent	None	7196A	0.0400	0.0377	94	90-110	

J Estimated concentration. The result is less than the PQL but greater than the MDL.

Approved By  24 Date: 4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report


Client : Battelle  
 Project Name : JPL Groundwater Monitoring 1Q07  
 Project Number : G486090  
 Sample Matrix : WATER

Service Request : P0700347  
 Date Collected : 04/09/07  
 Date Received : 04/09/07  
 Date Extracted : NA  
 Date Analyzed : 04/09/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-25-5 Units : mg/L (ppm)  
 Lab Code : P0700347-001MS P0700347-001DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0436	0.0436	87	87	85-115	<1	

Approved By  Date : 4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/10/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By:  Date: 4/27/07

ICCBMDL/120594



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/10/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0293	97	90-110
CCV1	0.0301	0.0284	94	90-110

Approved By:  
CCV1A/120594



Date:

4/27/07

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

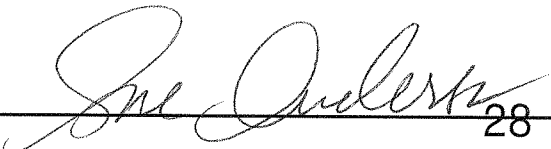
**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** 04/10/07  
**Date Received :** 04/10/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/10/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-26-2 Units : mg/L (ppm)  
 Lab Code : P0700347-007MS P0700347-007DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0468	0.0487	94	97	85-115	4	

Approved By  28 Date : 4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/11/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By:



Date:

4/27/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report


**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/11/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0299	99	90-110
CCV1	0.0301	0.0316	105	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/27/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** 04/11/07  
**Date Received :** 04/11/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/11/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-13 Units : mg/L (ppm)  
 Lab Code : P0700347-010MS P0700347-010DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	0.0411	0.0868	0.0877	91	93	85-115	1	

Approved By \_\_\_\_\_

*Sue Anderson*  
 31

Date : \_\_\_\_\_

4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/12/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By:



Date:

4/27/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/12/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0295	98	90-110
CCV1	0.0301	0.0286	95	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/27/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report

**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** 04/12/07  
**Date Received :** 04/12/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/12/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-8 Units : mg/L (ppm)  
 Lab Code : P0700347-014MS P0700347-014DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0465	0.0483	93	97	85-115	4	

Approved By

  
 34

Date :

4/27/07



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/13/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By:



Date:

4/27/07

ICCBMDL120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/13/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0294	98	90-110
CCV1	0.0301	0.0304	101	90-110

Approved By:



Date:

4/27/07

CCV1A/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report


Client : Battelle  
Project Name : JPL Groundwater Monitoring 1Q07  
Project Number : G486090  
Sample Matrix : WATER

Service Request : P0700347  
Date Collected : 04/13/07  
Date Received : 04/13/07  
Date Extracted : NA  
Date Analyzed : 04/13/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-10 Units : mg/L (ppm)  
Lab Code : P0700347-016MS P0700347-016DMS Basis : NA  
Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0529	0.0520	106	104	85-115	2	

Approved By  37 Date : 4/27/07

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/16/07

**Title:** Initial and Continuing Calibration Blank (ICB and CCB) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	PQL	MDL	Result
ICB	0.01	0.004	ND
CCB1	0.01	0.004	ND

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/27/07

ICCBMDL/120594

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

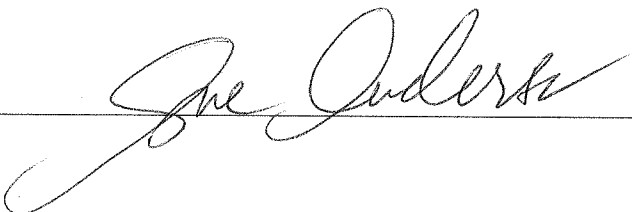
**Client:** Battelle  
**Project:** JPL Groundwater Monitoring 1Q07/G486090

**Service Request:** P0700347  
**Date Analyzed:** 04/16/07

**Title:** Initial and Continuing Calibration Verification (ICV and CCV) Summary  
**Analyte:** Chromium, Hexavalent  
**Method:** 7196A  
**Units:** mg/L (ppm)

Sample Name	True Value	Result	Percent Recovery	Acceptance Criteria
ICV	0.0301	0.0287	95	90-110
CCV1	0.0301	0.0297	99	90-110

Approved By: \_\_\_\_\_



Date: \_\_\_\_\_

4/27/07

CCV1A/120594

**COLUMBIA ANALYTICAL SERVICES, INC.**

QA/QC Report


**Client :** Battelle  
**Project Name :** JPL Groundwater Monitoring 1Q07  
**Project Number :** G486090  
**Sample Matrix :** WATER

**Service Request :** P0700347  
**Date Collected :** 04/16/07  
**Date Received :** 04/16/07  
**Date Extracted :** NA  
**Date Analyzed :** 04/16/07

Matrix Spike/Duplicate Matrix Spike Summary

Sample Name : MW-6 Units : mg/L (ppm)  
 Lab Code : P0700347-017MS P0700347-017DMS Basis : NA  
 Test Notes :

Analyte	Prep Method	Analysis Method	PQL	Spike Level		Sample Result	Spike Result		Spike Recovery		CAS Acceptance Limits	Relative Percent Difference	Result Notes
				MS	DMS		MS	DMS	MS	DMS			
Chromium, Hexavalent	None	7196A	0.01	0.0500	0.0500	ND	0.0468	0.0478	94	96	85-115	2	

Approved By  40 Date : 4/27/07