

Level C Data Package Deliverables

# Metals



Applied P & Ch Laboratory

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.

Project ID: JPL

Sample ID: 03M1458-MB-01

Sample Type: Method Blank

Project No: 04-4428.10

Service ID: 33082

Lab Sample ID: 03M1458-MB-01

Sample Matrix: Water

Collection Date: 05/13/2003

Collected by:

Received Date: 05/13/2003

Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	< 200	U	P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	6.5	B	P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	22.5	B	P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	39.1	B	P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	400	B	P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor

C Qualifier: U - Not Detected or less than IDL

B - Less than RL (PQL, EQL or CRDL), but greater than IDL.

Q Qualifier: N - Spike recovery out of control

\* - Duplicate analysis out of control

W - Post digestion spike for GFAA out of control

E - Serial dilution difference out of control

M Qualifier: P - ICP

A - FLAA

F - GFAA

CV - Cold Vapor

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.

Project ID: JPL

Sample ID: **EB-10-5/6/03**

Sample Type: Field Sample

Project No: 04-4428.10

Service ID: 33082

Lab Sample ID: 03-3082-1

Sample Matrix: Water

Collection Date: 05/06/2003

Collected by:

Received Date: 05/06/2003

Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	< 200	U	P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	31.4	B	P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	56.5	B	P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	59.3	B	P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	1200	B	P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor

C Qualifier: U - Not Detected or less than IDL

B - Less than RL (PQL, EQL or CRDL), but greater than IDL.

Q Qualifier: N - Spike recovery out of control

\* - Duplicate analysis out of control

W - Post digestion spike for GFAA out of control

E - Serial dilution difference out of control

M Qualifier: P - ICP

A - FLAA

F - GFAA

CV - Cold Vapor

## Applied P & Ch Laboratory Metal Analysis Results

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/06/2003
Project ID: JPL	Service ID: 33082	Collected by:
Sample ID: MW-11-1	Lab Sample ID: 03-3082-2	Received Date: 05/06/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	μg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	μg/L	200	56200		P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	μg/L	50	92.6		P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	μg/L	100	18700		P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	μg/L	400	3130		P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	μg/L	2000	24700		P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL    D-Date: Digestion Date;    A-Date: Analysis Date;    DF: Dilution Factor

C Qualifier: U - Not Detected or less than IDL	B - Less than RL (PQL, EQL or CRDL), but greater than IDL.
Q Qualifier: N - Spike recovery out of control	* - Duplicate analysis out of control
W - Post digestion spike for GFAA out of control	E - Serial dilution difference out of control
M Qualifier: P - ICP                    A - FLAA                    F - GFAA	CV - Cold Vapor

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.  
 Project ID: JPL  
 Sample ID: MW-11-2  
 Sample Type: Field Sample

Project No: 04-4428.10  
 Service ID: 33082  
 Lab Sample ID: 03-3082-3  
 Sample Matrix: Water

Collection Date: 05/06/2003  
 Collected by:  
 Received Date: 05/06/2003  
 Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	39200		P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	216		P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	15900		P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	2770		P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	21200		P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor  
 C Qualifier: U - Not Detected or less than IDL B - Less than RL (PQL, EQL or CRDL), but greater than IDL.  
 Q Qualifier: N - Spike recovery out of control \* - Duplicate analysis out of control  
 W - Post digestion spike for GFAA out of control E - Serial dilution difference out of control  
 M Qualifier: P - ICP A - FLAA F - GFAA CV - Cold Vapor

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.  
 Project ID: JPL  
 Sample ID: MW-11-3  
 Sample Type: Field Sample

Project No: 04-4428.10  
 Service ID: 33082  
 Lab Sample ID: 03-3082-4  
 Sample Matrix: Water

Collection Date: 05/06/2003  
 Collected by:  
 Received Date: 05/06/2003  
 Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	38600		P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	2310		P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	11900		P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	2050		P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	24700		P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor  
 C Qualifier: U - Not Detected or less than IDL B - Less than RL (PQL, EQL or CRDL), but greater than IDL.  
 Q Qualifier: N - Spike recovery out of control \* - Duplicate analysis out of control  
 W - Post digestion spike for GFAA out of control E - Serial dilution difference out of control  
 M Qualifier: P - ICP A - FLAA F - GFAA CV - Cold Vapor

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.  
 Project ID: JPL  
 Sample ID: MW-11-4  
 Sample Type: Field Sample

Project No: 04-4428.10  
 Service ID: 33082  
 Lab Sample ID: 03-3082-5  
 Sample Matrix: Water

Collection Date: 05/06/2003  
 Collected by:  
 Received Date: 05/06/2003  
 Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	9050		P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	61.5		P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	7550		P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	1820		P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	25100		P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor  
 C Qualifier: U - Not Detected or less than IDL B - Less than RL (PQL, EQL or CRDL), but greater than IDL.  
 Q Qualifier: N - Spike recovery out of control \* - Duplicate analysis out of control  
 W - Post digestion spike for GFAA out of control E - Serial dilution difference out of control  
 M Qualifier: P - ICP A - FLAA F - GFAA CV - Cold Vapor

Applied P & Ch Laboratory  
**Metal Analysis Results**

Client Name: GEOFON, Inc.  
 Project ID: JPL  
 Sample ID: MW-11-5  
 Sample Type: Field Sample

Project No: 04-4428.10  
 Service ID: 33082  
 Lab Sample ID: 03-3082-6  
 Sample Matrix: Water

Collection Date: 05/06/2003  
 Collected by:  
 Received Date: 05/06/2003  
 Moisture %: -

Element Name	CAS No	Unit	RL	Result	C	M	Q	Batch	D-Date	A-Date	DF	Method
ARSENIC	7440-38-2	µg/L	5	< 5	U	F		03M1458E	05/13/03	05/13/03	1	200.9
CALCIUM	7440-70-2	µg/L	200	13300		P		03M1442L	05/08/03	05/08/03	1	200.7
IRON	7439-89-6	µg/L	50	250		P		03M1442L	05/08/03	05/08/03	1	200.7
MAGNESIUM	7439-95-4	µg/L	100	1610		P		03M1442L	05/08/03	05/08/03	1	200.7
POTASSIUM	7440-09-7	µg/L	400	1170		P		03M1442L	05/08/03	05/08/03	1	200.7
SODIUM	7440-23-5	µg/L	2000	46800		P		03M1442L	05/08/03	05/08/03	1	200.7

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: RL: PQL (EQL) or CRDL D-Date: Digestion Date; A-Date: Analysis Date; DF: Dilution Factor

C Qualifier: U - Not Detected or less than IDL

B - Less than RL (PQL, EQL or CRDL), but greater than IDL.

Q Qualifier: N - Spike recovery out of control

\* - Duplicate analysis out of control

W - Post digestion spike for GFAA out of control

E - Serial dilution difference out of control

M Qualifier: P - ICP

A - FLAA

F - GFAA

CV - Cold Vapor



FORM-2A Metal  
Applied P & Ch Laboratory  
**Initial and Continuing Calibration Verification**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1442L
Batch No.(s): 03M1442	Instrument: ICP -L	Method: 200.7

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	ICV 12:47			CCV 13:06			CCV 13:33			CCV 14:01		
		True	Result	%R	True	Result	%R	True	Result	%R	True	Result	%R
1	Calcium	100000.0	99157.68	99.2	50000.0	50933.48	101.9	50000.0	49353.50	98.7	50000.0	50677.68	101.4
2	Iron	10000.0	9950.21	99.5	5000.0	4980.48	99.6	5000.0	4891.39	97.8	5000.0	5003.66	100.1
3	Magnesium	50000.0	51051.63	102.1	25000.0	25016.86	100.1	25000.0	24834.93	99.3	25000.0	25378.85	101.5
4	Potassium	30000.0	29959.65	99.9	15000.0	15169.41	101.1	15000.0	15006.94	100.0	15000.0	15035.59	100.2
5	Sodium	200000.0	197263.13	98.6	100000.0	104546.71	104.5	100000.0	102475.75	102.5	100000.0	105263.09	105.3

- (a) ICV Control Limit 95-105%; For Hg, 90-110%.
- (b) CCV Control Limit 90-110%; For Hg, 80-120%.

FORM-2A Metal  
Applied P & Ch Laboratory  
**Initial and Continuing Calibration Verification**

Client Name: GEOFON, Inc.

Project No: 04-4428.10

Lab Code: APCL

Project Name: JPL

Service ID: 033082

Sequence No.: 03M1442L

Instrument: ICP -L

Method: 200.7

Batch No.(s): 03M1442

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	CCV 14:24			CCV 14:52			CCV 15:18			CCV 15:36		
		True	Result	%R	True	Result	%R	True	Result	%R	True	Result	%R
1	Calcium	50000.0	50736.71	101.5	50000.0	50446.97	100.9	50000.0	50846.14	101.7	50000.0	50191.61	100.4
2	Iron	5000.0	5071.85	101.4	5000.0	4977.75	99.6	5000.0	4958.01	99.2	5000.0	5057.88	101.2
3	Magnesium	25000.0	25642.50	102.6	25000.0	25159.45	100.6	25000.0	25248.49	101.0	25000.0	25320.68	101.3
4	Potassium	15000.0	15551.38	103.7	15000.0	15194.11	101.3	15000.0	15141.38	100.9	15000.0	15156.78	101.0
5	Sodium	100000.0	107049.55	107.0	100000.0	105128.01	105.1	100000.0	104240.20	104.2	100000.0	103058.88	103.1

(a) ICV Control Limit 95-105%; For Hg, 90-110%.

(b) CCV Control Limit 90-110%; For Hg, 80-120%.

FORM-2A Metal  
Applied P & Ch Laboratory  
**Initial and Continuing Calibration Verification**

Client Name: GEOFON, Inc.  
Project Name: JPL  
Batch No.(s): 03M1458

Project No: 04-4428.10      Lab Code: APCL  
Service ID: 033082      Sequence No.: 03M1458E  
Instrument: GFAA-E      Method: 200.9

Analysis Date: 05/13/03

Concentration Units: UG/L

#	Analyte	ICV 11:10			CCV 12:27			CCV 13:42			CCV 14:59		
		True	Result	%R	True	Result	%R	True	Result	%R	True	Result	%R
1	Arsenic	50.0	51.40	102.8	50.0	56.60	113.2	50.0	55.60	111.2	50.0	64.60	129.2

(a) ICV Control Limit 95-105%; For Hg, 90-110%.

(b) CCV Control Limit 90-110%; For Hg, 80-120%.

FORM-2A Metal  
Applied P & Ch Laboratory  
**Initial and Continuing Calibration Verification**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1458E
Batch No.(s): 03M1458	Instrument: GFAA-E	Method: 200.9

Analysis Date: 05/13/03

Concentration Units: UG/L

#	Analyte	CCV 15:15			CCV 16:06			True	Result	%R	True	Result	%R
		True	Result	%R	True	Result	%R						
1	Arsenic	50.0	51.00	102.0	50.0	49.70	99.4						

(a) ICV Control Limit 95-105%; For Hg, 90-110%.

(b) CCV Control Limit 90-110%; For Hg, 80-120%.

FORM-2B Metal  
 Applied P & Ch Laboratory  
**CRDL Standard For AA and ICP**

Client Name: GEOFON, Inc.  
 Project Name: JPL  
 Batch No.(s): 03M1442

Project No: 04-4428.10  
 Service ID: 033082  
 Instrument: ICP -L

Lab Code: APCL  
 Sequence No.: 03M1442L  
 Method: 200.7

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	True	12:58		Time	
			Found	R%	Found	R%
1	Calcium	1000.0	1044.07	104.4		
2	Iron	50.0	46.86	93.7		
3	Magnesium		14.04			
4	Potassium		45.67			
5	Sodium		375.85			

FORM-3 Metal  
Applied P & Ch Laboratory  
**Metal ICB/CCB Summary**

Client Name: GEOFON, Inc.  
Project Name: JPL

Project No: 04-4428.10  
Service ID: 033082  
Instrument: ICP -L

Lab Code: APCL  
Sequence No.: 03M1442L  
Method: 200.7

Batch No.(s): 03M1442

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	ICB 12:51		CCB 13:08		CCB 13:36		CCB 14:04		CCB 14:27	
		Result	C	Result	C	Result	C	Result	C	Result	C
1	Calcium	-550.40		59.00	U	-283.01		-501.76		-295.18	
2	Iron	1.50	U	11.92	B	1.50	U	2.03	B	2.67	B
3	Magnesium	11.37	B	40.30	B	9.70	U	16.76	B	17.63	B
4	Potassium	37.50	B	42.52	B	41.79	B	43.33	B	41.42	B
5	Sodium	309.43	B	463.29	B	303.00	B	537.71	B	239.23	B

FORM-3 Metal  
Applied P & Ch Laboratory  
Metal ICB/CCB Summary

Client Name: GEOFON, Inc.  
Project Name: JPL  
Batch No.(s): 03M1442

Project No: 04-4428.10      Lab Code: APCL  
Service ID: 033082      Sequence No.: 03M1442L  
Instrument: ICP -L      Method: 200.7

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	CCB 14:55		CCB 15:20		CCB 15:38		CCB Time		CCB Time	
		Result	C	Result	C	Result	C	Result	C	Result	C
1	Calcium	-404.48		-190.57	B	-127.93	B				
2	Iron	3.10	B	3.23	B	70.00					
3	Magnesium	20.34	B	50.91	B	163.85					
4	Potassium	49.06	B	46.57	B	46.48	B				
5	Sodium	234.15	B	568.97	B	233.38	B				

FORM-3 Metal  
 Applied P & Ch Laboratory  
**Metal ICB/CCB Summary**

Client Name: GEOFON, Inc.  
 Project Name: JPL  
 Batch No.(s): 03M1458

Project No: 04-4428.10      Lab Code: APCL  
 Service ID: 033082      Sequence No.: 03M1458E  
 Instrument: GFAA-E      Method: 200.9

Analysis Date: 05/13/03

Concentration Units: UG/L

#	Analyte	ICB 11:17		CCB 12:34		CCB 13:49		CCB 15:06		CCB 15:21	
		Result	C	Result	C	Result	C	Result	C	Result	C
1	Arsenic	2.10	U	2.10	U	2.10	U	2.10	U	2.10	U



FORM-3 Metal  
 Applied P & Ch Laboratory  
**Metal ICB/CCB Summary**

Client Name: GEOFON, Inc.  
 Project Name: JPL  
 Batch No.(s): 03M1458

Project No: 04-4428.10      Lab Code: APCL  
 Service ID: 033082      Sequence No.: 03M1458E  
 Instrument: GFAA-E      Method: 200.9

Analysis Date: 05/13/03

Concentration Units: UG/L

#	Analyte	CCB 16:12		CCB Time		CCB Time		CCB Time		CCB Time	
		Result	C	Result	C	Result	C	Result	C	Result	C
1	Arsenic	2.10	U								

FORM-4 Metal  
Applied P & Ch Laboratory  
**ICP Interference Check Sample**

Client Name: GEOFON, Inc.  
Project Name: JPL

Project No: 04-4428.10  
Service ID: 033082  
ICP ID Number: ICP -L

Lab Code: APCL  
Sequence No.: 03M1442L

Batch No.(s): 03M1442

Analysis Date: 05/08/03

Concentration Units: UG/L

#	Analyte	Expected		Initial	Found	%R	Final	Found	%R
		Sol. A	Sol. AB	13:00	13:03		15:31	15:33	
1	Calcium	500000	500000	468054	470353.8	94.1	459157	462841.1	92.6
2	Iron	200000	200000	179857	184181.8	92.1	178418	186233.3	93.1
3	Magnesium	500000	500000	472613	481539.6	96.3	473200	491814.1	98.4
4	Potassium	0	0	42	33.1		50	38.3	
5	Sodium	0	0	35	375.7		121	494.6	

FORM-5A Metal

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 200.9

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03M1458E	
MS Filename: -	Date Analyzed: 051303	Time Analyzed: 12:01
MSD Filename: -	Date Analyzed: 051303	Time Analyzed: 12:08
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
ARSENIC	µg/L	50	0	54.9	110	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
ARSENIC	µg/L	50	55.0	110	0	20	75-125
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-5A Metal

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 200.7

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03M1442L	
MS Filename: -	Date Analyzed: 050803	Time Analyzed: 13:26
MSD Filename: -	Date Analyzed: 050803	Time Analyzed: 13:28
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
CALCIUM	µg/L	20000	38600	59200	103	75-125
IRON	µg/L	1000	2310	3420	111	75-125
MAGNESIUM	µg/L	10000	11900	22700	108	75-125
POTASSIUM	µg/L	5000	2050	7690	113	75-125
SODIUM	µg/L	40000	24700	69200	111	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CALCIUM	µg/L	20000	57800	96	7	20	75-125
IRON	µg/L	1000	3270	96	14	20	75-125
MAGNESIUM	µg/L	10000	21600	97	11	20	75-125
POTASSIUM	µg/L	5000	7850	116	3	20	75-125
SODIUM	µg/L	40000	66000	103	7	20	75-125
# of Out-of-control					0	0	

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-5B Metal  
Applied P & Ch Laboratory  
**Post Digest Spike Sample Recovery**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1442L
	Batch No.: 03M1442	Method: 200.7
Spike Sample No. : 03-3082-04	Matrix: WATER	Instrument: ICP -L
Client Sample No.: MW-11-3	Analysis Date: 05/08/03	

Concentration Units: UG/L

#	Analyte	Spiked Sample Result(SSR)	13:31 C	Sample Result(SR)	13:18 C	Spike Added(SA)	% Rec.	Control Limit	Q
1	Calcium	58408.4219		38602.4375		20000.00	99.0	75-125	
2	Iron	3324.6311		2311.5544		1000.00	101.3	75-125	
3	Magnesium	22020.9512		11939.2461		10000.00	100.8	75-125	
4	Potassium	7691.8008		2051.4998		5000.00	112.8	75-125	
5	Sodium	67122.6719		24669.8223		40000.00	106.1	75-125	

FORM-5B Metal  
Applied P & Ch Laboratory  
**Post Digest Spike Sample Recovery**

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Lab Code:	APCL
Project Name:	JPL	Service ID:	033082	Sequence No.:	03M1458E
		Batch No.:	03M1458	Method:	200.9
Spike Sample No. :	03-3082-04	Matrix:	WATER	Instrument:	GFAA-E
Client Sample No.:	MW-11-3	Analysis Date:	05/13/03		

Concentration Units: UG/L

#	Analyte	Spiked Sample 12:14		Sample 11:42		Spike Added(SA)	% Rec.	Control Limit	Q
		Result(SSR)	C	Result(SR)	C				
1	Arsenic	56.4000		0.9000	U	50.00	112.8	75-125	

FORM-6 Metal  
Applied P & Ch Laboratory  
**Duplicates Verification**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1442L
	Batch No.: 03M1442	Method: 200.7
Spike Sample No. 03-3082-04	Matrix: WATER	Instrument: ICP -L
Client Sample No. MW-11-3	% Solid: 0.00	Analysis Date: 05/08/03

Concentration Unit: UG/L

#	Analyte	13:18		13:21		RPD(%)	Q
		Sample(s)	C	Duplicate	C		
1	Calcium	38602.4375		39149.0078		1.4	
2	Iron	2311.5544		2342.1619		1.3	
3	Magnesium	11939.2461		12161.7412		1.8	
4	Potassium	2051.4998		2012.4011		1.9	
5	Sodium	24669.8223		25210.3086		2.2	

FORM-6 Metal  
Applied P & Ch Laboratory  
**Duplicates Verification**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1458E
Spike Sample No. 03-3082-04	Batch No.: 03M1458	Method: 200.9
Client Sample No. MW-11-3	Matrix: WATER	Instrument: GFAA-E
	% Solid: 0.00	Analysis Date: 05/13/03

Concentration Unit: UG/L

#	Analyte	11:42 Sample(s)	C	11:48 Duplicate	C	RPD(%)	Q
1	Arsenic	0.9000	U	0.5000	U		



FORM-7 Metal

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 200.9

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03M1458E	
LCS Filename: -	Date Analyzed: 051303	Time Analyzed: 11:29
LCSD Filename: -	Date Analyzed: 051303	Time Analyzed: 11:36

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
ARSENIC	µg/L	50	0	56.8	114	80-120
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
ARSENIC	µg/L	50	57.1	114	0	20	80-120
# of Out-of-control					0	0	

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits      D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-7 Metal

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 200.7

Client Name: GEOFON, Inc.

Contract No:

Lab Code: APCL

Case No:

SAS No:

Service ID: 33082

Project ID: JPL

Project No: 04-4428.10

Sample Matrix: Water

Batch No: 03M1442L

LCS Filename: -

Date Analyzed: 050803

Time Analyzed: 13:13

LCSD Filename: -

Date Analyzed: 050803

Time Analyzed: 13:16

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
CALCIUM	µg/L	20000	0	20400	102	80-120
IRON	µg/L	1000	0	1030	103	80-120
MAGNESIUM	µg/L	10000	0	9390	94	80-120
POTASSIUM	µg/L	5000	0	4890	98	80-120
SODIUM	µg/L	40000	0	41800	105	80-120
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CALCIUM	µg/L	20000	20500	103	1	20	80-120
IRON	µg/L	1000	1040	104	1	20	80-120
MAGNESIUM	µg/L	10000	9430	94	0	20	80-120
POTASSIUM	µg/L	5000	4900	98	0	20	80-120
SODIUM	µg/L	40000	41300	103	2	20	80-120
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-9 Metal  
Applied P & Ch Laboratory  
Serial Dilution

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Lab Code:	APCL
Project Name:	JPL	Service ID:	033082	Sequence No.:	03M1442L
		Batch No.:	03M1442	Method:	200.7
Dilution Sample No.:	03-3082-04	Matrix:	WATER	Instrument:	ICP -L
Client Sample No.:	MW-11-3	Analysis Date:	05/08/03		

Concentration Units: UG/L

#	Analyte	Initial Sample Results(I)	13:18 C	Serial Dilut Results(S)	13:23 C	% Diff.	Q
1	Calcium	38602.44		38426.93		0.5	
2	Iron	2311.55		2332.19		0.9	
3	Magnesium	11939.25		11456.36		4.0	
4	Potassium	2051.50		1838.65	B	10.4	
5	Sodium	24669.82		23330.49		5.4	

FORM-9 Metal  
Applied P & Ch Laboratory  
**Serial Dilution**

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Lab Code:	APCL
Project Name:	JPL	Service ID:	033082	Sequence No.:	03M1458E
		Batch No.:	03M1458	Method:	200.9
Dilution Sample No.:	03-3082-04	Matrix:	WATER	Instrument:	GFAA-E
Client Sample No.:	MW-11-3	Analysis Date:	05/13/03		

Concentration Units: **UG/L**

#	Analyte	Initial Sample Results(I)	11:42 C	Serial Dilut Results(S)	11:55 C	% Diff.	Q
1	Arsenic	0.90	U	-0.50	U		

FORM-13 Metal  
Applied P & Ch Laboratory  
**Preparation Log**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1442L
	Batch No.: 03M1442	Method: 200.7
Preparation Matrix: WATER	Instrument: ICP -L	

#	Client Sample No.	APCL Sample No.	Preparation Date	Weight (gram)	Volume (ml)
1	MW-11-3	03-3082-04DM	05/08/03		50.0
2	EB-10-5/6/03	03-3082-01	05/08/03		50.0
3	MW-11-1	03-3082-02	05/08/03		50.0
4	MW-11-2	03-3082-03	05/08/03		50.0
5	MW-11-4	03-3082-05	05/08/03		50.0
6	MW-11-5	03-3082-06	05/08/03		50.0
7	DUPE-6-2Q03	03-3102-01	05/08/03		50.0
8	EB-11-5/7/03	03-3102-02	05/08/03		50.0
9	MW-12-1	03-3102-03	05/08/03		50.0
10	MW-12-2	03-3102-04	05/08/03		50.0
11	MW-12-3	03-3102-05	05/08/03		50.0
12	MW-12-4	03-3102-06	05/08/03		50.0
13	MW-12-5	03-3102-07	05/08/03		50.0
14	A	03-3105-01	05/08/03		50.0
15	B	03-3105-02	05/08/03		50.0
16	C	03-3105-03	05/08/03		50.0
17	D	03-3105-04	05/08/03		50.0
18		03M1442MB	05/08/03		50.0
19		03M1442LCS	05/08/03		50.0
20		03M1442LCSD	05/08/03		50.0
21	MW-11-3 Dup.	03M1442MD	05/08/03		50.0
22	MW-11-3 MS	03M1442MS	05/08/03		50.0
23	MW-11-3 MSD	03M1442MSD	05/08/03		50.0

FORM-13 Metal  
Applied P & Ch Laboratory  
Preparation Log

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Lab Code: APCL
Project Name: JPL	Service ID: 033082	Sequence No.: 03M1458E
	Batch No.: 03M1458	Method: 200.9
Preparation Matrix: WATER	Instrument: GFAA-E	

#	Client Sample No.	APCL Sample No.	Preparation Date	Weight (gram)	Volume (ml)
1	MW-11-3	03-3082-04DM	05/13/03		50.0
2	EB-10-5/6/03	03-3082-01	05/13/03		50.0
3	MW-11-1	03-3082-02	05/13/03		50.0
4	MW-11-2	03-3082-03	05/13/03		50.0
5	MW-11-4	03-3082-05	05/13/03		50.0
6	MW-11-5	03-3082-06	05/13/03		50.0
7	DUPE-6-2Q03	03-3102-01	05/13/03		50.0
8	EB-11-5/7/03	03-3102-02	05/13/03		50.0
9	MW-12-1	03-3102-03	05/13/03		50.0
10	MW-12-2	03-3102-04	05/13/03		50.0
11	MW-12-3	03-3102-05	05/13/03		50.0
12	MW-12-4	03-3102-06	05/13/03		50.0
13	MW-12-5	03-3102-07	05/13/03		50.0
14	EB-12-5/8/03	03-3130-01	05/13/03		50.0
15	MW-22-1	03-3130-02	05/13/03		50.0
16	MW-22-2	03-3130-03	05/13/03		50.0
17	MW-22-3	03-3130-04	05/13/03		50.0
18	MW-22-4	03-3130-05	05/13/03		50.0
19	MW-22-5	03-3130-06	05/13/03		50.0
20		03M1458MB	05/13/03		50.0
21		03M1458LCS	05/13/03		50.0
22		03M1458LCSD	05/13/03		50.0
23	MW-11-3 Dup.	03M1458MD	05/13/03		50.0
24	MW-11-3 MS	03M1458MS	05/13/03		50.0
25	MW-11-3 MSD	03M1458MSD	05/13/03		50.0

FORM-14 Metal  
Applied P & Ch Laboratory  
Analysis Run Log

Client Name: GEOFON, Inc.  
Project Name: JPL

Project No: 04-4428.10  
Service ID: 033082  
Instrument: ICP -L  
Start Date: 05/08/03

Lab Code: APCL  
Sequence No.: 03M1442L  
Method: 200.7  
End Date: 05/08/03

Batch No.(s): 03M1442

#	APCL Sample No.	D/F	Time	Al	Sb	As	Ba	Be	Cd	Ca	Cr	Co	Cu	Fe	Pb	Mg	Mn	Hg	Ni	K	Se	Ag	Na	Tl	V	Zn	Mo	Sr	Ti	Sn	Li	B	Si
1	Calib Blank	1.00	12:37							✓			✓	✓						✓		✓											
2	STD1 1423A	1.00	12:40							✓			✓	✓						✓		✓											
3	STD2 1423B	1.00	12:42							✓			✓	✓						✓		✓											
4	STD3 1423C	1.00	12:45							✓			✓	✓						✓		✓											
5	ICV 1447A	1.00	12:47							✓			✓	✓						✓		✓											
6	ICB	1.00	12:51							✓			✓	✓						✓		✓											
7	CRI A1432	1.00	12:58							✓			✓	✓						✓		✓											
8	ICSA 1441	1.00	13:00							✓			✓	✓						✓		✓											
9	ICSAB 1443	1.00	13:03							✓			✓	✓						✓		✓											
10	CCV 1447B	1.00	13:06							✓			✓	✓						✓		✓											
11	CCB	1.00	13:08							✓			✓	✓						✓		✓											
12	M-BL 03M1442 W	1.00	13:11							✓			✓	✓						✓		✓											
13	LCS-03M1442	1.00	13:13							✓			✓	✓						✓		✓											
14	LCSD-03M1442	1.00	13:16							✓			✓	✓						✓		✓											
15	3082-4 S F=1	1.00	13:18							✓			✓	✓						✓		✓											
16	3082-4 D F=1	1.00	13:21							✓			✓	✓						✓		✓											
17	3082-4 1/5 F=5	5.00	13:23							✓			✓	✓						✓		✓											
18	3082-4 MS F=1	1.00	13:26							✓			✓	✓						✓		✓											
19	3082-4 MSD F=1	1.00	13:28							✓			✓	✓						✓		✓											
20	3082-4 PS F=1	1.00	13:31							✓			✓	✓						✓		✓											
21	CCV 1447B	1.00	13:33							✓			✓	✓						✓		✓											
22	CCB	1.00	13:36							✓			✓	✓						✓		✓											
23	3082-1 F=1	1.00	13:38							✓			✓	✓						✓		✓											
24	3082-2 F=1	1.00	13:41							✓			✓	✓						✓		✓											
25	3082-3 F=1	1.00	13:44							✓			✓	✓						✓		✓											
26	3082-5 F=1	1.00	13:46							✓			✓	✓						✓		✓											
27	3082-6 F=1	1.00	13:49							✓			✓	✓						✓		✓											
28	3102-1 F=1	1.00	13:51							✓			✓	✓						✓		✓											
29	3102-2 F=1	1.00	13:54							✓			✓	✓						✓		✓											
30	3102-3 F=1	1.00	13:56							✓			✓	✓						✓		✓											
31	3102-4 F=1	1.00	13:59							✓			✓	✓						✓		✓											
32	CCV 1447B	1.00	14:01							✓			✓	✓						✓		✓											
33	CCB	1.00	14:04							✓			✓	✓						✓		✓											
34	3102-5 F=1	1.00	14:06							✓			✓	✓						✓		✓											
35	3102-6 F=1	1.00	14:09							✓			✓	✓						✓		✓											
36	3102-7 F=1	1.00	14:12							✓			✓	✓						✓		✓											
37	3105-1 F=1	1.00	14:14							✓			✓	✓						✓		✓											
38	3105-2 F=1	1.00	14:17							✓			✓	✓						✓		✓											
39	3105-3 F=1	1.00	14:19							✓			✓	✓						✓		✓											
40	3105-4 F=1	1.00	14:22							✓			✓	✓						✓		✓											

FORM-14 Metal  
Applied P & Ch Laboratory  
Analysis Run Log

Client Name: GEOFON, Inc.  
Project Name: JPL

Project No: 04-4428.10  
Service ID: 033082  
Instrument: ICP -L  
Start Date: 05/08/03

Lab Code: APCL  
Sequence No.: 03M1442L  
Method: 200.7  
End Date: 05/08/03

Batch No.(s): 03M1442

#	APCL Sample No.	D/F	Time	Al	Sb	As	Ba	Be	Cd	Ca	Cr	Co	Cu	Fe	Pb	Mg	Mn	Hg	Ni	K	Se	Ag	Na	Tl	V	Zn	Mo	Sr	Ti	Sn	Li	B	Si	
41	CCV 1447B	1.00	14:24							✓			✓	✓					✓			✓												
42	CCB	1.00	14:27							✓			✓	✓					✓			✓												
43	M-BL 03M1443 W	1.00	14:29							✓			✓	✓					✓			✓												
44	LCS-03M1443	1.00	14:32							✓			✓	✓					✓			✓												
45	LCSD-03M1443	1.00	14:34							✓			✓	✓					✓			✓												
46	3100-1 S F=1	1.00	14:37							✓			✓	✓					✓			✓												
47	3100-1 D F=1	1.00	14:39							✓			✓	✓					✓			✓												
48	3100-1 1/5 F=5	5.00	14:42							✓			✓	✓					✓			✓												
49	3100-1 MS F=1	1.00	14:44							✓			✓	✓					✓			✓												
50	3100-1 MSD F=1	1.00	14:47							✓			✓	✓					✓			✓												
51	3100-1 PS F=1	1.00	14:50							✓			✓	✓					✓			✓												
52	CCV 1447B	1.00	14:52							✓			✓	✓					✓			✓												
53	CCB	1.00	14:55							✓			✓	✓					✓			✓												
54	3100-3 F=1	1.00	14:57							✓			✓	✓					✓			✓												
55	3100-4 F=1	1.00	15:00							✓			✓	✓					✓			✓												
56	3100-5 F=1	1.00	15:02							✓			✓	✓					✓			✓												
57	3100-6 F=1	1.00	15:05							✓			✓	✓					✓			✓												
58	3100-7 F=1	1.00	15:08							✓			✓	✓					✓			✓												
59	3100-8 F=1	1.00	15:10							✓			✓	✓					✓			✓												
60	3100-9 F=1	1.00	15:13							✓			✓	✓					✓			✓												
61	3100-10 F=1	1.00	15:15							✓			✓	✓					✓			✓												
62	CCV 1447B	1.00	15:18							✓			✓	✓					✓			✓												
63	CCB	1.00	15:20							✓			✓	✓					✓			✓												
64	3100-11 F=1	1.00	15:23							✓			✓	✓					✓			✓												
65	3100-12 F=1	1.00	15:26							✓			✓	✓					✓			✓												
66	3100-13 F=1	1.00	15:28							✓			✓	✓					✓			✓												
67	ICSA 1441	1.00	15:31							✓			✓	✓					✓			✓												
68	ICSAB 1443	1.00	15:33							✓			✓	✓					✓			✓												
69	CCV 1447B	1.00	15:36							✓			✓	✓					✓			✓												
70	CCB	1.00	15:38							✓			✓	✓					✓			✓												



FORM-14 Metal  
Applied P & Ch Laboratory  
Analysis Run Log

Client Name: GEOFON, Inc.  
Project Name: JPL

Project No: 04-4428.10  
Service ID: 033082  
Instrument: GFAA-E  
Start Date: 05/13/03

Lab Code: APCL  
Sequence No.: 03M1458E  
Method: 200.9  
End Date: 05/13/03

Batch No.(s): 03M1458

#	APCL Sample No.	D/F	Time	Al	Sb	As	Ba	Be	Cd	Ca	Cr	Co	Cu	Fe	Pb	Mg	Mn	Hg	Ni	K	Se	Ag	Na	Tl	V	Zn	Mo	Sr	Ti	Sn	Li	B	Si			
1	AS Position 002	1.00	10:17			✓																														
2	AS Position 001	1.00	10:25			✓																														
3	Calib. Blank	1.00	10:34			✓																														
4	1/2 STD1 1472A	1.00	10:40			✓																														
5	STD1 1472A	1.00	10:46			✓																														
6	STD2 1472B	1.00	10:53			✓																														
7	STD3 1472C	1.00	10:59			✓																														
8	ICV A1474	1.00	11:10			✓																														
9	ICB	1.00	11:17			✓																														
10	M-BL 03M1458	1.00	11:23			✓																														
11	LCS-03M1458	1.00	11:29			✓																														
12	LCSD-03M1458	1.00	11:36			✓																														
13	3082-4 S F=1	1.00	11:42			✓																														
14	3082-4 D F=1	1.00	11:48			✓																														
15	3082-4 1/5 F=5	5.00	11:55			✓																														
16	3082-4 MS F=1	1.00	12:01			✓																														
17	3082-4 MSD F=1	1.00	12:08			✓																														
18	3082-4 PS F=1	1.00	12:14			✓																														
19	3082-1 F=1	1.00	12:21			✓																														
20	CCV A1474	1.00	12:27			✓																														
21	CCB	1.00	12:34			✓																														
22	3082-2 F=1	1.00	12:40			✓																														
23	3082-3 F=1	1.00	12:46			✓																														
24	3082-5 F=1	1.00	12:52			✓																														
25	3082-6 F=1	1.00	12:59			✓																														
26	3130-1 F=1	1.00	13:05			✓																														
27	3130-2 F=1	1.00	13:17			✓																														
28	CCV A1474	1.00	13:42			✓																														
29	CCB	1.00	13:49			✓																														
30	3130-3 F=1	1.00	14:08			✓																														
31	3130-4 F=1	1.00	14:14			✓																														
32	3130-5 F=1	1.00	14:21			✓																														
33	3130-6 F=1	1.00	14:27			✓																														
34	3102-1 F=1	1.00	14:34			✓																														
35	3102-2 F=1	1.00	14:40			✓																														
36	3102-3 F=1	1.00	14:47			✓																														
37	3102-4 F=1	1.00	14:53			✓																														
38	CCV A1474	1.00	14:59			✓																														
39	CCB	1.00	15:06			✓																														
40	CCV A1474	1.00	15:15			✓																														

FORM-14 Metal  
Applied P & Ch Laboratory  
Analysis Run Log

Client Name: GEOFON, Inc.  
Project Name: JPL  
Batch No.(s): 03M1458

Project No: 04-4428.10  
Service ID: 033082  
Instrument: GFAA-E  
Start Date: 05/13/03

Lab Code: APCL  
Sequence No.: 03M1458E  
Method: 200.9  
End Date: 05/13/03

#	APCL Sample No.	D/F	Time	Al	Sb	As	Ba	Be	Cd	Ca	Cr	Co	Cu	Fe	Pb	Mg	Mn	Hg	Ni	K	Se	Ag	Na	Tl	V	Zn	Mo	Sr	Ti	Sn	Li	B	Si
41	CCB	1.00	15:21			✓																											
42	3102-5 F=1	1.00	15:27			✓																											
43	3102-6 F=1	1.00	15:33			✓																											
44	3102-7 F=1	1.00	15:40			✓																											
45	CCV A1474	1.00	16:06			✓																											
46	CCB	1.00	16:12			✓																											

Batch # B3M1442 Matrix: W Method used: 3010 A Date: 5/8/03 Digested by: XI Diluted by: \_\_\_\_\_  
 Lot #: ASTM Type I water RW1410 HNO<sub>3</sub> 1102/20 H<sub>2</sub>SO<sub>4</sub> \_\_\_\_\_ HCl 4102050 H<sub>2</sub>O<sub>2</sub> \_\_\_\_\_

OP #	Type	Samp ID /Lot #	X (g or mL)	$V_{digest}/X = f_1$	$V_1/V_i = f_2$	$V_j/V_i = f_3$	$F = f_1 f_2 f_3$	Note
1495	Method Blank	Bl. Lot: <u>RW1410</u>	<u>50</u>	<u>50/X = 1</u>	<u>/ =</u>	<u>/ =</u>		
1496	LCS1	Bl. Lot: <u>1</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		<u>5 Me</u>
1497	Sample-1	<u>3082-4</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1498	MS1 on S-1	<u>-4</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		<u>T=95°C</u>
1499	MS2 on S-1	<u>-4</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1500	Sample 2	<u>-1</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1501	Sample 3	<u>-2</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1502	Sample 4	<u>-3</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1503	Sample 5	<u>-5</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1504	Sample 6	<u>-6</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1505	Sample 7	<u>3102-1</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1506	Sample 8	<u>-2</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1507	Sample 9	<u>-3</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1508	Sample 10	<u>-4</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1509	LCS2	Bl. Lot: <u>RW1410</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1510	Sample 11	<u>-5</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1511	Sample 12	<u>-6</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1512	Sample 13	<u>-7</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1513	Sample 14	<u>3105-1</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1514	Sample 15	<u>-2</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1515	Sample 16	<u>-3</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1516	Sample 17	<u>-4</u>		<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1517	Sample 18		<u>↓ ↓</u>	<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1518	Sample 19			<u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1519	Sample 20			<u>XI</u> <u>5/8/03</u> <u>/X =</u>	<u>/ =</u>	<u>/ =</u>		
1520	Duplicate	<u>3082-4</u>	<u>50</u>	<u>50/X = 1</u>	<u>/ =</u>	<u>/ =</u>		

Specification of matrix spike and lab control spike

QC Type	Spiked Element *	Spike Stock Solution Lot #	Spike Stock (Rep.) Conc. C <sub>s</sub> , µg/mL	Spike Stock Volum Used V <sub>s</sub> , mL	Spike Level T' = C <sub>s</sub> V <sub>s</sub> / V ppm or mg/L	Sample Spike T, ppm
MS1	/As/Sb/M <sub>20</sub>	AA- /AA- /AA- /AA- <u>143</u>	<u>/ / 125</u>	<u>/ / 12</u>	<u>/ / /</u>	
MS2	/As/Sb/M <sub>20</sub>	AA- /AA- /AA- /AA- <u>11</u>	<u>/ / /</u>	<u>/ / /</u>	<u>/ / /</u>	
LCS1	/As/Sb/M <sub>20</sub>	AA- /AA- /AA- /AA- <u>1492</u>	<u>/ / /</u>	<u>/ / /</u>	<u>/ / /</u>	
LCS2	/As/Sb/M <sub>20</sub>	AA- /AA- /AA- /AA- <u>11</u>	<u>/ / /</u>	<u>/ / /</u>	<u>/ / /</u>	

\* Notation: T - rep. sample spike level. T' - digest solution spike level. T = f T' = C<sub>s</sub> V<sub>s</sub> / X. M20 (or Mj) represents 20 (or i) metals, (see STD logbook).  
 If digest needs dilution for different metals, use dilution worksheet.  
 APCL form 6-116 April, 03, 1996. Ver. 4.0  
 Root-File: [CUST.DOC.AA]DIGEST\_ROOT.TEX File: [CUST.DOC.AA]DIGEST .TEX

6388  
 Supervisor Initial U.W

1780 Magnolia Ave. Chino CA 91710

Metal Digestion (3010/3050) Worksheet

Phone (909) 590-1828 Fax: (909) 590-1498

Batch # 03M1458 Matrix: W Method used: 3020A Date: 5/13/03 Digested by: DL Diluted by: \_\_\_\_\_

Lot #: ASTM Type I water RW11411 HNO<sub>3</sub> 110280 H<sub>2</sub>SO<sub>4</sub> \_\_\_\_\_ HCl \_\_\_\_\_ H<sub>2</sub>O<sub>2</sub> \_\_\_\_\_

OP #	Type	Samp ID /Lot #	X (g or mL)	V <sub>digest</sub> /X = f <sub>1</sub>	V <sub>1</sub> /V <sub>i</sub> = f <sub>2</sub>	V <sub>j</sub> /V <sub>i</sub> = f <sub>3</sub>	F = f <sub>1</sub> f <sub>2</sub> f <sub>3</sub>	Note
1755	Method Blank	Bl. Lot: <u>RW11411</u>	<u>50</u>	<u>50</u> /X= <u>1</u>	<u>1</u> =	<u>1</u> =		<u>GEAA/10</u>
1756	LCS1	Bl. Lot: <u>11</u>		/X=	<u>1</u> =	<u>1</u> =		<u>T=95°C</u>
1757	Sample-1	<u>3082-4</u>		/X=	<u>1</u> =	<u>1</u> =		
1758	MS1 on S-1	<u>4</u>		/X=	<u>1</u> =	<u>1</u> =		
1759	MS2 on S-1	<u>4</u>		/X=	<u>1</u> =	<u>1</u> =		
1760	Sample 2	<u>1</u>		/X=	<u>1</u> =	<u>1</u> =		
1761	Sample 3	<u>2</u>		/X=	<u>1</u> =	<u>1</u> =		
1762	Sample 4	<u>3</u>		/X=	<u>1</u> =	<u>1</u> =		
1763	Sample 5	<u>5</u>		/X=	<u>1</u> =	<u>1</u> =		
1764	Sample 6	<u>6</u>		/X=	<u>1</u> =	<u>1</u> =		
1765	Sample 7	<u>3130-1</u>		/X=	<u>1</u> =	<u>1</u> =		
1766	Sample 8	<u>2</u>		/X=	<u>1</u> =	<u>1</u> =		<u>DUP/TMS</u>
1767	Sample 9	<u>3</u>		/X=	<u>1</u> =	<u>1</u> =		
1768	Sample 10	<u>4</u>		/X=	<u>1</u> =	<u>1</u> =		
1769	LCS2	Bl. Lot: <u>RW11411</u>		/X=	<u>1</u> =	<u>1</u> =		
1770	Sample 11	<u>5</u>		/X=	<u>1</u> =	<u>1</u> =		
1771	Sample 12	<u>6</u>		/X=	<u>1</u> =	<u>1</u> =		
1772	Sample 13	<u>3102-1</u>		/X=	<u>1</u> =	<u>1</u> =		
1773	Sample 14	<u>2</u>		/X=	<u>1</u> =	<u>1</u> =		
1774	Sample 15	<u>3</u>		/X=	<u>1</u> =	<u>1</u> =		
1775	Sample 16	<u>4</u>		/X=	<u>1</u> =	<u>1</u> =		
1776	Sample 17	<u>5</u>		/X=	<u>1</u> =	<u>1</u> =		
1777	Sample 18	<u>6</u>		/X=	<u>1</u> =	<u>1</u> =		
1778	Sample 19	<u>7</u>		/X=	<u>1</u> =	<u>1</u> =		
1779	Sample 20	<u>DL 5/13/03</u>		/X=	<u>1</u> =	<u>1</u> =		
1780	Duplicate	<u>3082-4</u>	<u>50</u>	<u>50</u> /X= <u>1</u>	<u>1</u> =	<u>1</u> =		

Specification of matrix spike and lab control spike

QC Type	Spiked Element *	Spike Stock Solution Lot #	Spike Stock (Rep.) Conc. C <sub>s</sub> , µg/mL	Spike Stock Volum Used V <sub>s</sub> , mL	Spike Level T' = C <sub>s</sub> V <sub>s</sub> /V ppm or mg/L	Sample Spike T, ppm
MS1	<u>As/Sb/M<sub>20</sub></u>	AA- <u>1AA103</u> /AA- /AA-	<u>151</u> /	<u>10.5</u> /	<u>10.5</u> /	
MS2	<u>As/Sb/M<sub>20</sub></u>	AA- <u>1AA-11</u> /AA- /AA-	<u>1</u> / /	<u>1</u> / /	<u>1</u> / /	
LCS1	<u>As/Sb/M<sub>20</sub></u>	AA- <u>1AA101</u> /AA- /AA-	<u>1</u> / /	<u>1</u> / /	<u>1</u> / /	
LCS2	<u>As/Sb/M<sub>20</sub></u>	AA- <u>1AA-11</u> /AA- /AA-	<u>1</u> / /	<u>1</u> / /	<u>1</u> / /	

\* Notation: T - rep. sample spike level. T' - digest solution spike level. T = f T' = C<sub>s</sub>V<sub>s</sub>/X. M20 (or M<sub>j</sub>) represents 20 (or j) metals, (see STD logbook). If digest needs dilution for different metals, use dilution worksheet.

Level C Data Package Deliverables

# Wet Chemistry



Applied P & Ch Laboratory

Applied P & Ch Laboratory  
**Wet Analysis Results for Method SM2320B**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method SM2320B  
 Collected by:

Component Name: Bicarbonate  
 CAS No:

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	< 2	U
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	208	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	171	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	176	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	80.5	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/10/03	03W2817	mg/L	2	114	
03W2817-MB-01	03W2817-MB-01	Water	05/10/03	05/10/03	05/10/03	03W2817	mg/L	2	< 2	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
**Wet Analysis Results for Method SM2320B**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method SM2320B  
 Collected by:

Component Name: Carbonate  
 CAS No:

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	<2	U
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	<2	U
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	2.6	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	<2	U
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	20.4	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	5.2	
03W2817-MB-01	03W2817-MB-01	Water	05/10/03	05/10/03	05/10/03	03W2817	mg-CaCO <sub>3</sub> /L	2	<2	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
Wet Analysis Results for Method 9040B

Client Name: GEOFON, Inc.  
Project ID: JPL

Project No: 04-4428.10  
Service ID: 33082

Anal. Method 9040B  
Collected by:

Component Name: pH  
CAS No: 10-29-7

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	6.95	
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	7.85	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	8.08	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	7.95	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	8.77	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	8.43	
03W2740-MB-01	03W2740-MB-01	Water	05/06/03	05/06/03	05/06/03	03W2740	pH unit	0.01	6.84	

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.



Applied P & Ch Laboratory  
**Wet Analysis Results for Method 160.1**

Client Name: GEOFON, Inc.                      Project No: 04-4428.10                      Anal. Method 160.1  
 Project ID: JPL                                      Service ID: 33082                              Collected by:

Component Name: Solids, Total Dissolved (TDS)  
 CAS No: 10-33-3

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	<10	U
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	323	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	215	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	222	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	120	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/08/03	03W2783	mg/L	10	156	
03W2783-MB-01	03W2783-MB-01	Water	05/08/03	05/08/03	05/08/03	03W2783	mg/L	10	<10	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
**Wet Analysis Results for Method 7196**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method 7196  
 Collected by:

Component Name: Chromium (VI)  
 CAS No: 1333-82-0

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U
03W2737-MB-01	03W2737-MB-01	Water	05/06/03	05/06/03	05/06/03	03W2737	mg/L	0.01	<0.01	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
**Wet Analysis Results for Method 314.0**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method 314.0  
 Collected by:

Component Name: Perchlorate  
 CAS No:

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/09/03	03W2804	µg/L	4	<4	U
03W2804-MB-01	03W2804-MB-01	Water	05/09/03	05/09/03	05/09/03	03W2804	µg/L	4	<4	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
Wet Analysis Results for Method 300.0

Client Name: GEOFON, Inc.                      Project No: 04-4428.10                      Anal. Method 300.0  
Project ID: JPL                                      Service ID: 33082                              Collected by:

Component Name: Chloride Cl<sup>-</sup>  
CAS No: 16887-00-6

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.25	0.33	
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.5	20.0	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.5	15.4	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.5	11.8	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.4	11.6	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.4	12.2	
03W2750-MB-01	03W2750-MB-01	Water	05/07/03	05/07/03	05/07/03	03W2750	mg/L	0.2	<0.2	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
**Wet Analysis Results for Method 300.0**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method 300.0  
 Collected by:

Component Name: Nitrate as N  
 CAS No: 14797-55-8

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.05	0.16	
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.1	0.97	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.1	0.15	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.1	0.15	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.08	0.11	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.08	0.11	
03W2750-MB-01	03W2750-MB-01	Water	05/07/03	05/07/03	05/07/03	03W2750	mg/L	0.04	<0.04	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

Applied P & Ch Laboratory  
**Wet Analysis Results for Method 300.0**

Client Name: GEOFON, Inc.  
 Project ID: JPL

Project No: 04-4428.10  
 Service ID: 33082

Anal. Method 300.0  
 Collected by:

Component Name: Sulfate  $\text{SO}_4^{--}$   
 CAS No: 14808-79-8

Lab ID	Sample ID	Matrix	Coll. Date	Rcv Date	Anal. Date	Batch	Unit	RL	Result	Q
03-3082-1	EB-10-5/6/03	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	0.63	0.82	
03-3082-2	MW-11-1	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	1.3	43.9	
03-3082-3	MW-11-2	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	1.3	33.6	
03-3082-4	MW-11-3	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	1.3	22.5	
03-3082-5	MW-11-4	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	1	2.2	
03-3082-6	MW-11-5	Water	05/06/03	05/06/03	05/07/03	03W2750	mg/L	1	9.7	
03W2750-MB-01	03W2750-MB-01	Water	05/07/03	05/07/03	05/07/03	03W2750	mg/L	0.5	<0.5	U

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Note: Q - Qualifier.

Qualifier: U - Not Detected or less than MDL

B - Less than RL (PQL, EQL or CRDL), but greater than MDL.

## FORM-3

Applied P &amp; Ch Laboratory

## Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 300.0

Client Name: GEOFON, Inc.

Contract No:

Lab Code: APCL

Case No:

SAS No:

Service ID: 33082

Project ID: JPL

Project No: 04-4428.10

Sample Matrix: Water

Batch No: 03W2750

LCS Filename: -

Date Analyzed: 050703

Time Analyzed: 10:16

LCSD Filename: -

Date Analyzed: 050703

Time Analyzed: 10:29

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
CHLORIDE CL <sup>-</sup>	mg/L	4.0	0	4.17	104	80-120
NITRATE AS N	mg/L	1.5	0	1.58	105	80-120
SULFATE SO <sub>4</sub> <sup>2-</sup>	mg/L	15	0	15.4	103	80-120
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHLORIDE CL <sup>-</sup>	mg/L	4.0	4.20	105	1	20	80-120
NITRATE AS N	mg/L	1.5	1.59	106	1	20	80-120
SULFATE SO <sub>4</sub> <sup>2-</sup>	mg/L	15	15.5	103	0	25	80-120
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_

\_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 300.0

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2750	
MS Filename: -	Date Analyzed: 050703	Time Analyzed: 13:39
MSD Filename: -	Date Analyzed: 050703	Time Analyzed: 14:03
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
CHLORIDE CL <sup>-</sup>	mg/L	20.0	11.8	32.5	104	75-125
NITRATE AS N	mg/L	7.50	0.15	7.70	101	75-125
SULFATE SO <sub>4</sub> <sup>-</sup>	mg/L	75.0	22.5	102	106	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHLORIDE CL <sup>-</sup>	mg/L	20.0	32.7	105	1	20	75-125
NITRATE AS N	mg/L	7.50	7.81	102	1	20	75-125
SULFATE SO <sub>4</sub> <sup>-</sup>	mg/L	75.0	102	106	0	25	75-125
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits      D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_



FORM-3

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 314.0

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2804	
LCS Filename: -	Date Analyzed: 050903	Time Analyzed: 12:52
LCSD Filename: -	Date Analyzed: -	Time Analyzed: -

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
PERCHLORATE	µg/L	25	0	24.7	99	80-120
# of Out-of-control					0	

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 314.0

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2804	
MS Filename: -	Date Analyzed: 050903	Time Analyzed: 17:15
MSD Filename: -	Date Analyzed: 050903	Time Analyzed: 17:34
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
PERCHLORATE	µg/L	50	0	51.3	103	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
PERCHLORATE	µg/L	50	51.6	103	0	20	75-125
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 160.1

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2783	
LCS Filename: -	Date Analyzed: 050803	Time Analyzed: 13:54
LCSD Filename: -	Date Analyzed: 050803	Time Analyzed: 13:54

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	0	412	103	88-108
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	390	98	5	20	88-108
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits      D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 160.1

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2783	
MS Filename: -	Date Analyzed: 050803	Time Analyzed: 13:54
MSD Filename: -	Date Analyzed: 050803	Time Analyzed: 13:54
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	222	631	102	80-119
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	619	99	3	20	80-119
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 7196

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2737	
LCS Filename: -	Date Analyzed: 050603	Time Analyzed: 13:25
LCSD Filename: -	Date Analyzed: 050603	Time Analyzed: 13:25

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
CHROMIUM (VI)	mg/L	0.25	0	0.252	101	80-115
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHROMIUM (VI)	mg/L	0.25	0.246	98	3	19	80-115
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits      D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 7196

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33082
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2737	
MS Filename: -	Date Analyzed: 050603	Time Analyzed: 13:25
MSD Filename: -	Date Analyzed: 050603	Time Analyzed: 13:25
MS Sample No: MW-11-3	Sample Lab ID: 03-3082-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
CHROMIUM (VI)	mg/L	0.25	0	0.231	92	78-115
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHROMIUM (VI)	mg/L	0.25	0.227	91	1	19	78-115
# of Out-of-control				0	0		

# Column to be used to flag recovery and RPD values:

\* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: \_\_\_\_\_  
 \_\_\_\_\_

**Wet Chemistry QC Report B**  
**Duplicate Results**

**Matrix: Water**

**APCL Service ID: 03-3082**

Analysis	Batch ID	Analysis Date	Sample Name	Unit	Result	Duplicate Result	RPD %	RPD Control limit
Bicarbonate	03W2817	05/10/2003	MW-11-1	mg/L	208	210	1	20
Carbonate	03W2817	05/10/2003	MW-11-1	mg-CaCO <sub>3</sub> /L	ND	ND	NC	20
pH	03W2740	05/06/2003	03-3090-1	pH unit	7.11	7.15	1	20

Note: N/A = Not applicable; NR: Not requested; NC= Not Calculated; ND: Not detected.

6A  
INITIAL CALIBRATION DATA

Lab Name: Applied P & Ch Lab Contract: \_\_\_\_\_

Analysis: Chromium (VI) Calibration Date: 01/29/2003

Concentration (mg/L)	0.000	0.0125	0.050	0.125	0.250	0.50
Absorbance	0.000	0.006	0.041	0.109	0.214	0.415

**A = 0.000 + 0.836C**

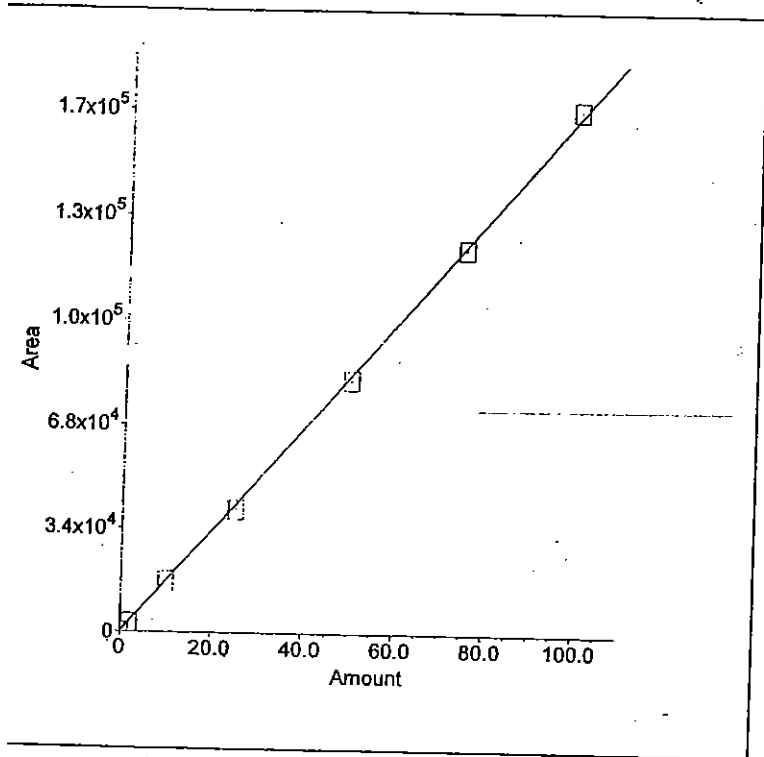
**A=Absorbance**

**C=Concentration (mg/L)**

**r = 0.9997**



1. Component: perchlorate  
Standard: External Fit Type: Linear  
Origin: Force Calibration: Area  
 $r^2=0.999492$   
Amt=0.0005893\*Resp+0



Calibration : 7 points , 0, 2, 10, 25, 50, 75, 100 ppb

Analyst C.W  
Date 03/12/03  
Instrument IC-10

# APCL Perchlorate Analysis Report

Sample Name : ICV 50 ppb w7828a

Data File Name : C:\DATA\E314-011\icv-50pb\_009.DXD

Method File Name : C:\PEAKNET\METHOD\E314-011.met

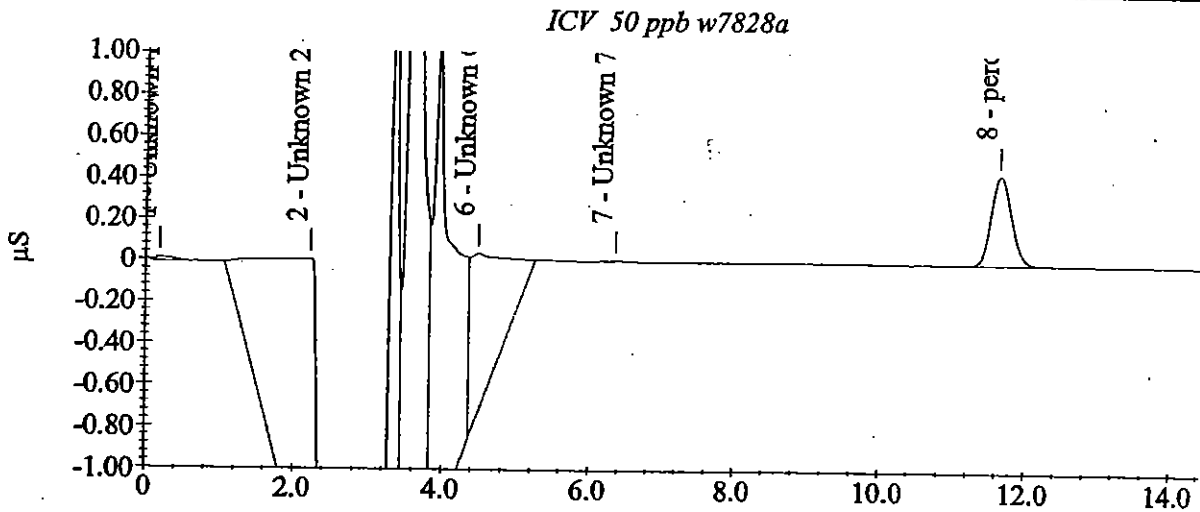
Date Time Collected : 03/12/2003 8:16:15 PM

System Operator : wei wang

Dilution Factor : 1.00

## Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
8	perchlorate	11.65	49.49	83990	4321



# APCL Perchlorate Analysis Report

Sample Name : icb

Data File Name : C:\DATA\E314-011\ICB\_010.DXD

Method File Name : c:\PeakNet\method\E314-011.met

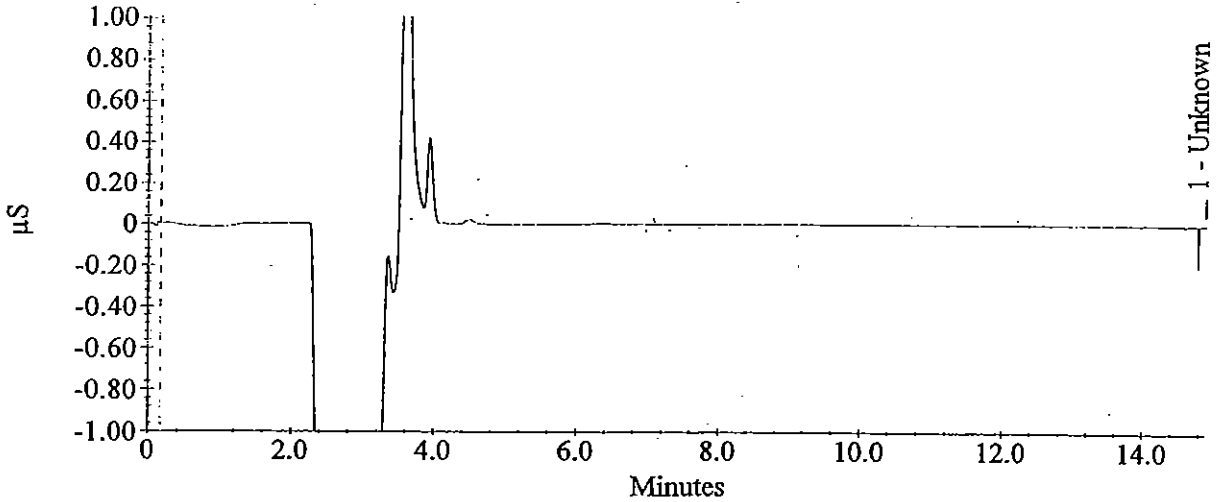
Date Time Collected : 03/12/2003 8:33:51 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
--------	----------------	----------------	--------------	-----------	-------------



# DIONEX METHOD PARAMETERS - E300-063.MET

Method Comment: APCL EPA 300 ANALYSIS DX-100  
 Column ID: Dionex AS4A-SC  
 Analyst ID: David

### System Parameters

System Name: DX-100

Number of Detectors.....	1
Run Time (minutes).....	10.00
Sampling Rate (seconds).....	0.20
Detector 1 Type.....	COND
Detector 1 real time plot scale maximum (uS ).....	30.000
minimum.....	-3.000
Detector 1 Output Equivalent to 1 Volt (in uS ) .....	30.00
Detector 1 ACI Analog Input Connection .....	DET1
Save Data File.....	Yes
Data File Name: C:\DX\DATA\03W1991\W1991Q01.D10	

-- DETECTOR 1 PARAMETERS --

### Report Options

Create ASCII Report File.....	No
Print Report.....	Yes
Print All Components.....	Yes
Print Components Found.....	No
Print Missing Components.....	No
Print All Peaks.....	No
Print Unknown Peaks.....	No
Print Chromatogram.....	Yes
<del>Autoscale Chromatogram Maximum.....</del>	<del>No</del>
Autoscale Chromatogram Minimum.....	No
Fill Peaks with Color .....	No
Draw Grid Lines on Chromatogram.....	No
Show Component Fraction Numbers.....	No
Label with Peak Number.....	No
Label with Retention Times on Chromatogram.....	Yes
Label with Component Name.....	Yes
Format File Name: C:\DX\METHOD\DEFAULT.PRF	

### Integration Parameters

Starting Peak Width (seconds).....	10.0
Peak Threshold .....	2.000
Peak Area Reject.....	1000
Area Reject for Reference Peaks.....	1000

### Data Events

Time	Description
-----	
0.13	Start peak detection
10.00	Stop peak detection

Calibration Parameters

Number Of Levels for Calibration.....	6
Force Calibration Curve Through Origin.....	No
Calibration Fit Type.....	Linear
Replace Or Average Calibrations.....	Replace
External or Internal Calibration.....	External
Calculate Unknowns by Area or Height.....	Area
Default Sample Volume.....	1.0
Default Dilution Factor.....	1.0
Default Response Factor for Unknown Peaks.....	0.0
Calibration Standard Volume .....	1.0
Internal Standard Amount in Samples .....	1.0
Amount Units .....	ppm

Component # 4 Bromide Retention Time 3.45  
 Reference Comp. Nitrate-N Window Size 0.20 min.  
 Amount =  $K0 + K1 \cdot \text{Area}$   
 $K0 = 4.58974E-002$   
 $K1 = 4.83279E-006$

Level	Amount	Area	Height
1	7.50000E-002	13830	1488
2	1.50000E+000	298206	30100
3	3.00000E+000	591234	61776
4	6.00000E+000	1219933	128845
5	7.50000E+000	1559887	166594
6	0.00000E+000	0	0

Component # 5 Nitrate-N Retention Time 3.87  
 Reference Comp. Nitrate-N Window Size 0.25 min.  
 Amount =  $K0 + K1 \cdot \text{Area}$   
 $K0 = 4.24689E-002$   
 $K1 = 8.05553E-007$

Level	Amount	Area	Height
1	3.75000E-002	40157	3802
2	7.50000E-001	849179	77129
3	1.50000E+000	1713421	152776
4	3.00000E+000	3610927	313707
5	3.75000E+000	4688990	396441
6	0.00000E+000	0	0

Component # 6 Phosphate-P Retention Time 6.38  
 Reference Comp. Phosphate-P Window Size 0.60 min.  
 Amount =  $K0 + K1 \cdot \text{Area}$   
 $K0 = 8.68926E-002$   
 $K1 = 2.12227E-006$

Level	Amount	Area	Height
1	7.50000E-002	24783	1450
2	1.50000E+000	642376	38579
3	3.00000E+000	1301126	79971
4	6.00000E+000	2756481	168994
5	7.50000E+000	3546397	217521
6	0.00000E+000	0	0

Component Table -- Last Modified: 17:42 on Fri, 21 Mar 2003

Component # 1 Fluoride Retention Time 1.32  
Reference Comp. Fluoride Window Size 0.15 min.  
Amount = K0 + K1\*Area  
K0 = -9.62851E-004  
K1 = 1.37614E-006

Level	Amount	Area	Height
1	2.50000E-002	28534	2732
2	5.00000E-001	373164	44629
3	1.00000E+000	707646	82595
4	2.00000E+000	1435865	173007
5	2.50000E+000	1837162	220914
6	0.00000E+000	0	0

Component # 2 Chloride Retention Time 1.97  
Reference Comp. Chloride Window Size 0.15 min.  
Amount = K0 + K1\*Area  
K0 = 1.28188E-001  
K1 = 1.95287E-006

Level	Amount	Area	Height
1	1.00000E-001	51206	7044
2	2.00000E+000	909455	126181
3	4.00000E+000	1856586	261681
4	8.00000E+000	3987563	585791
5	1.00000E+001	5142155	754321
6	0.00000E+000	0	0

Component # 3 Nitrite-N Retention Time 2.33  
Reference Comp. Chloride Window Size 0.15 min.  
Amount = K0 + K1\*Area  
K0 = 2.38085E-002  
K1 = 9.76240E-007

Level	Amount	Area	Height
1	3.75000E-002	30884	3582
2	7.50000E-001	734006	79701
3	1.50000E+000	1468106	162005
4	3.00000E+000	3021523	336616
5	3.75000E+000	3856614	429219
6	0.00000E+000	0	0

Component # 7      Sulfate                      Retention Time      7.92  
 Reference Comp.   Sulfate                      Window Size         0.90 min.  
 Amount = K0 + K1\*Area  
 K0     = 5.32283E-001  
 K1     = 2.53252E-006

Level	Amount	Area	Height
1	3.76000E-001	129999	6524
2	7.50000E+000	2598757	138579
3	1.50000E+001	5330209	287851
4	3.00000E+001	11507107	615917
5	3.75000E+001	14859049	776426
6	0.00000E+000	0	0

Timed Events File: C:\DX\METHOD\W761CAL.TE

Step	Time	Description
Init		ACI Autosmp OFF
Init		ACI pump st ON
Init		ACI inject OFF
Init		ACI auto zer OFF
Init		ACI TTL 1 OFF
Init		ACI TTL 2 OFF
Init		ACI TTL 3 OFF
Init		ACI TTL 4 OFF
Init		ACI OFF
Init		ACI OFF
1	0.0	ACI Autosmp ON
1	0.0	ACI auto zer ON
2	2.5	ACI Autosmp OFF
2	2.5	ACI inject ON
2	2.5	ACI TTL 1 ON
2	2.5	Start Sampling



Component: Fluoride

Fit Type: Linear

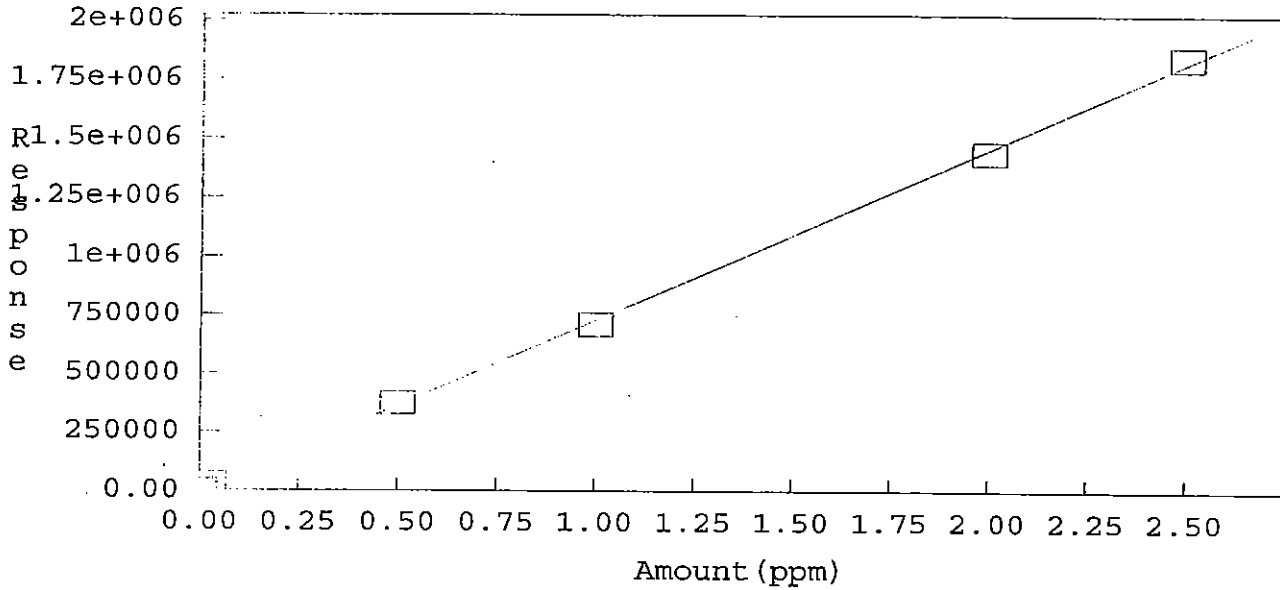
$r^2 = 0.999552$

$\text{Amt} = \text{Resp} * 1.376e-006 + -0.000962$

$\text{Resp} = \text{Amt} * 7.267e+005 + 699.7$

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Chloride

Fit Type: Linear

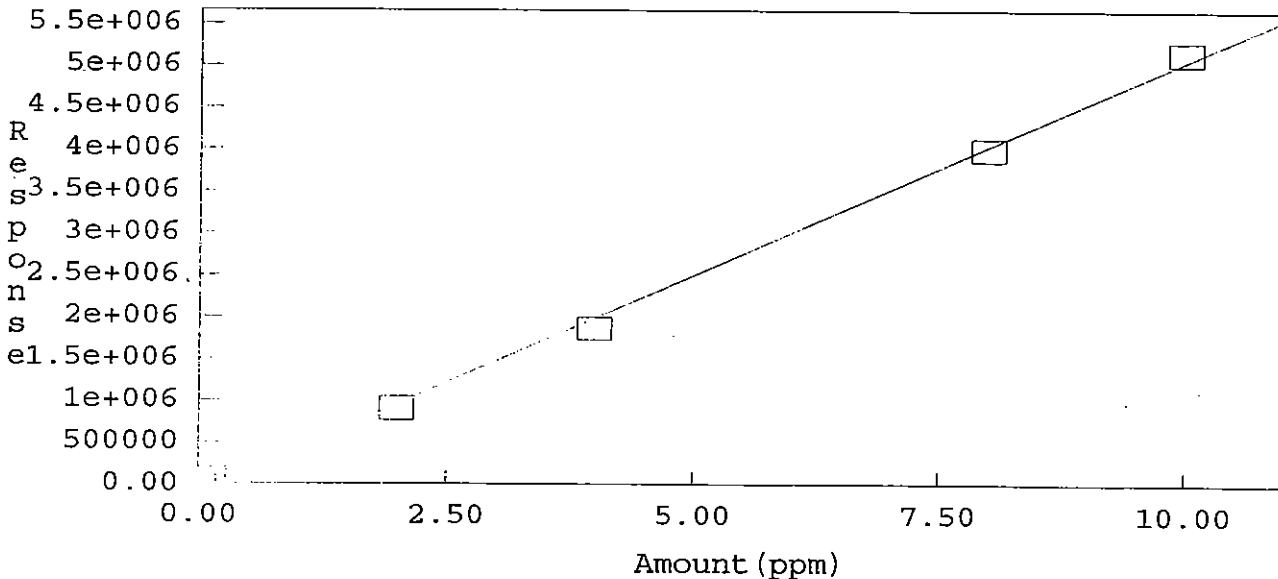
$r^2 = 0.998409$

$\text{Amt} = \text{Resp} * 1.953e-006 + 0.1282$

$\text{Resp} = \text{Amt} * 5.121e+005 + -6.564e+00$

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Nitrite-N

Fit Type: Linear

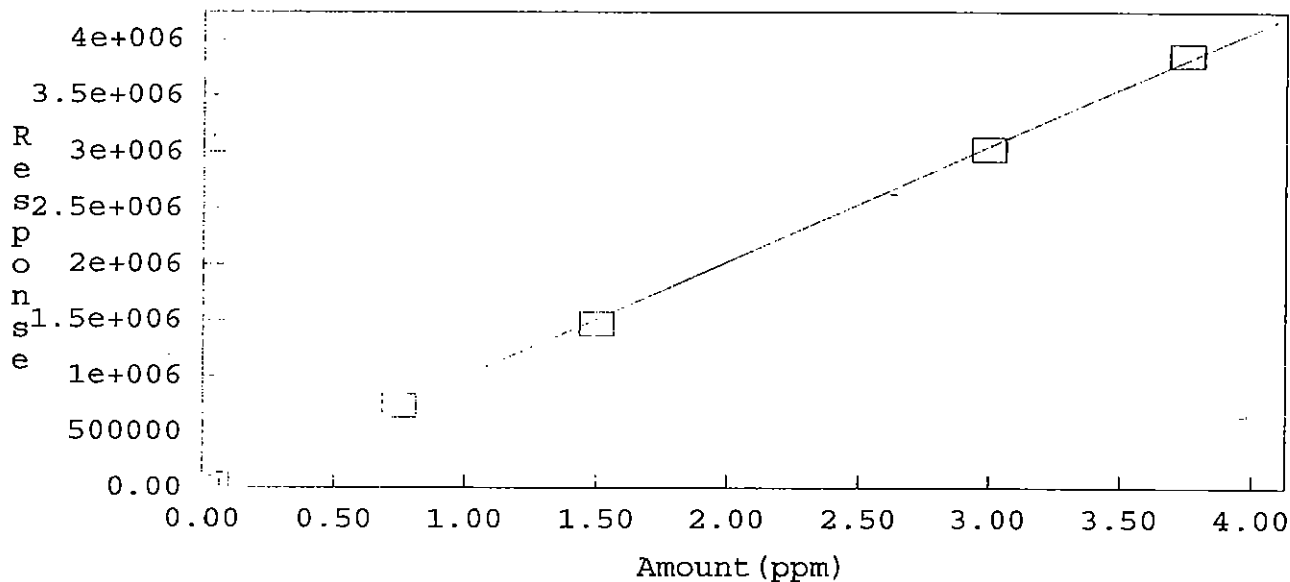
$r^2 = 0.999594$

Amt = Resp \*  $9.762e-007$  + 0.02381

Resp = Amt \*  $1.024e+006$  +  $-2.439e+00$

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Bromide

Fit Type: Linear

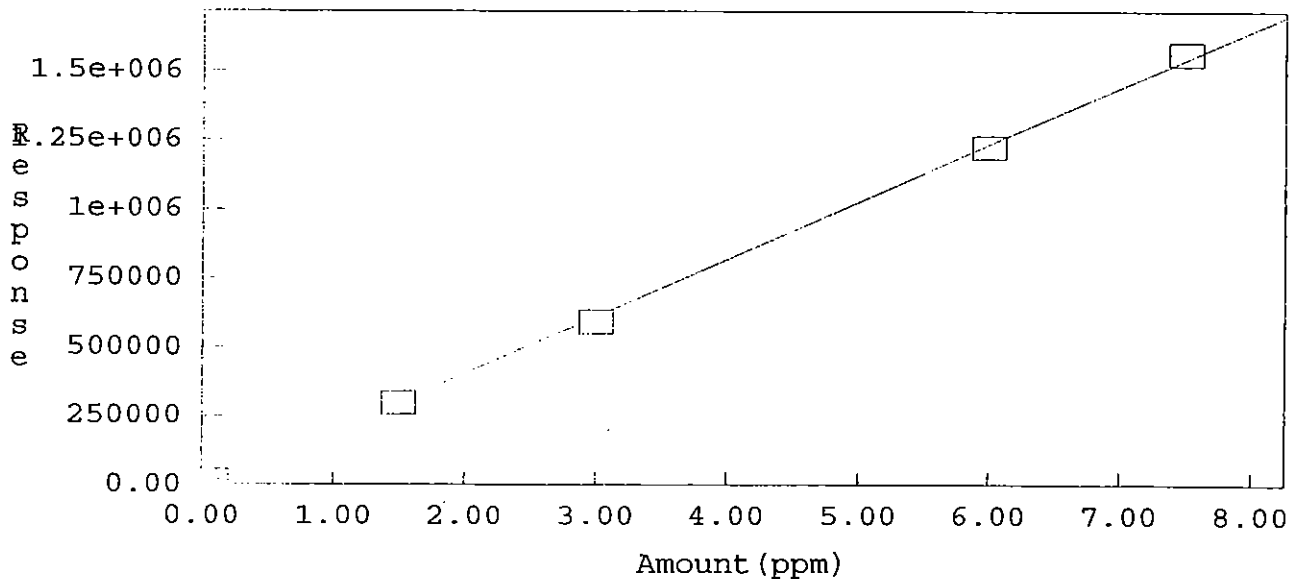
$r^2 = 0.999518$

Amt = Resp \*  $4.833e-006$  + 0.0459

Resp = Amt \*  $2.069e+005$  + -9497

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Nitrate-N

Fit Type: Linear

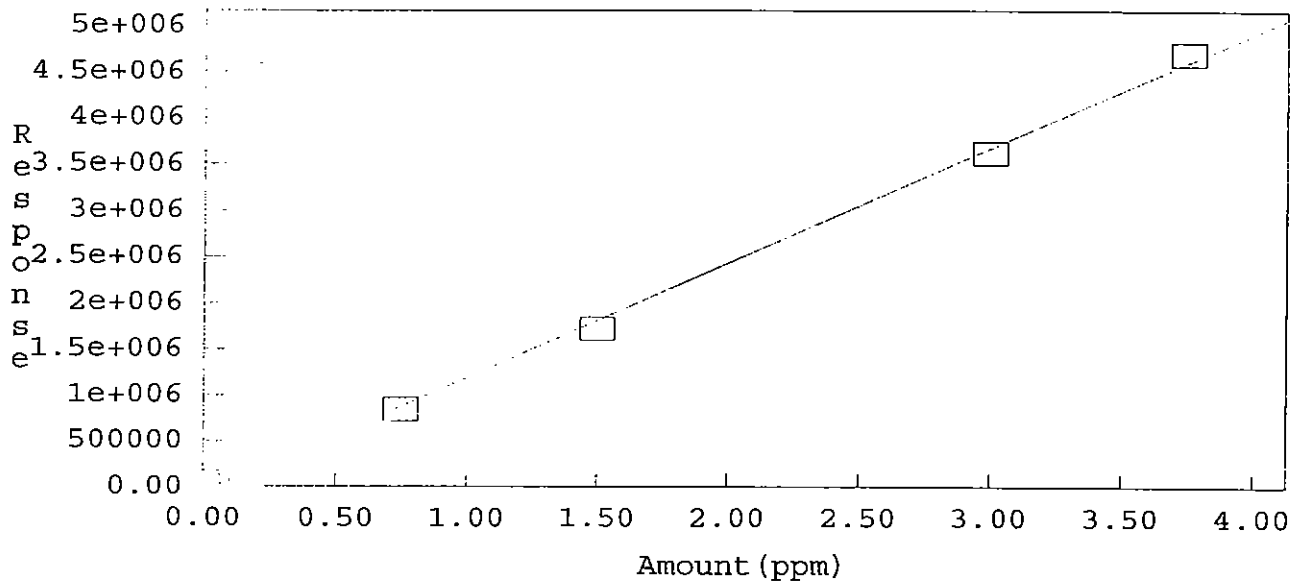
$r^2 = 0.998618$

Amt = Resp \*  $8.056e-007$  + 0.04247

Resp = Amt \*  $1.241e+006$  + -5.272e+00

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Phosphate-P

Fit Type: Linear

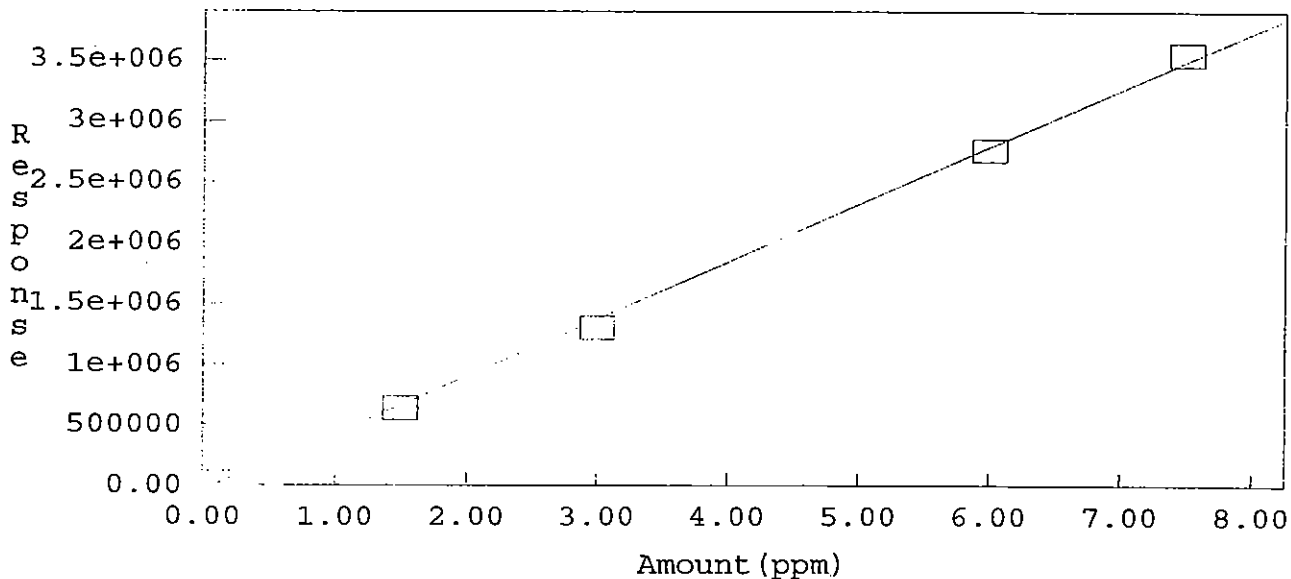
$r^2 = 0.998898$

Amt = Resp \*  $2.122e-006$  + 0.08689

Resp = Amt \*  $4.712e+005$  + -4.094e+00

Standardization: External

Calibration: Area



Method: C:\DX\METHOD\E300-063.MET

Component: Sulfate

Fit Type: Linear

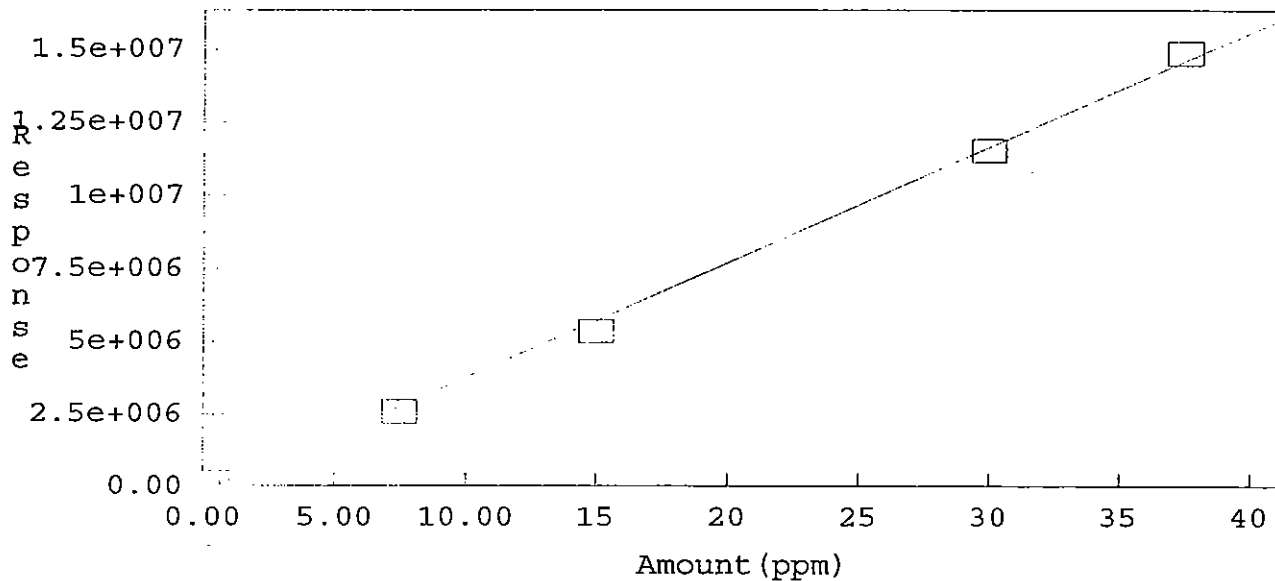
$r^2 = 0.998245$

$\text{Amt} = \text{Resp} * 2.533\text{e-}006 + 0.5323$

$\text{Resp} = \text{Amt} * 3.949\text{e+}005 + -2.102\text{e+}00$

Standardization: External

Calibration: Area



```

=====
Sample Name: ICV-W7768-100X          Date: 03/21/2003 17:56:33
Data File  : C:\DX\DATA\e300-063\W7768Q01.D07
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 7   Detector: COND
Analyst    : David                   Column: Dionex AS4A-SC
=====

```

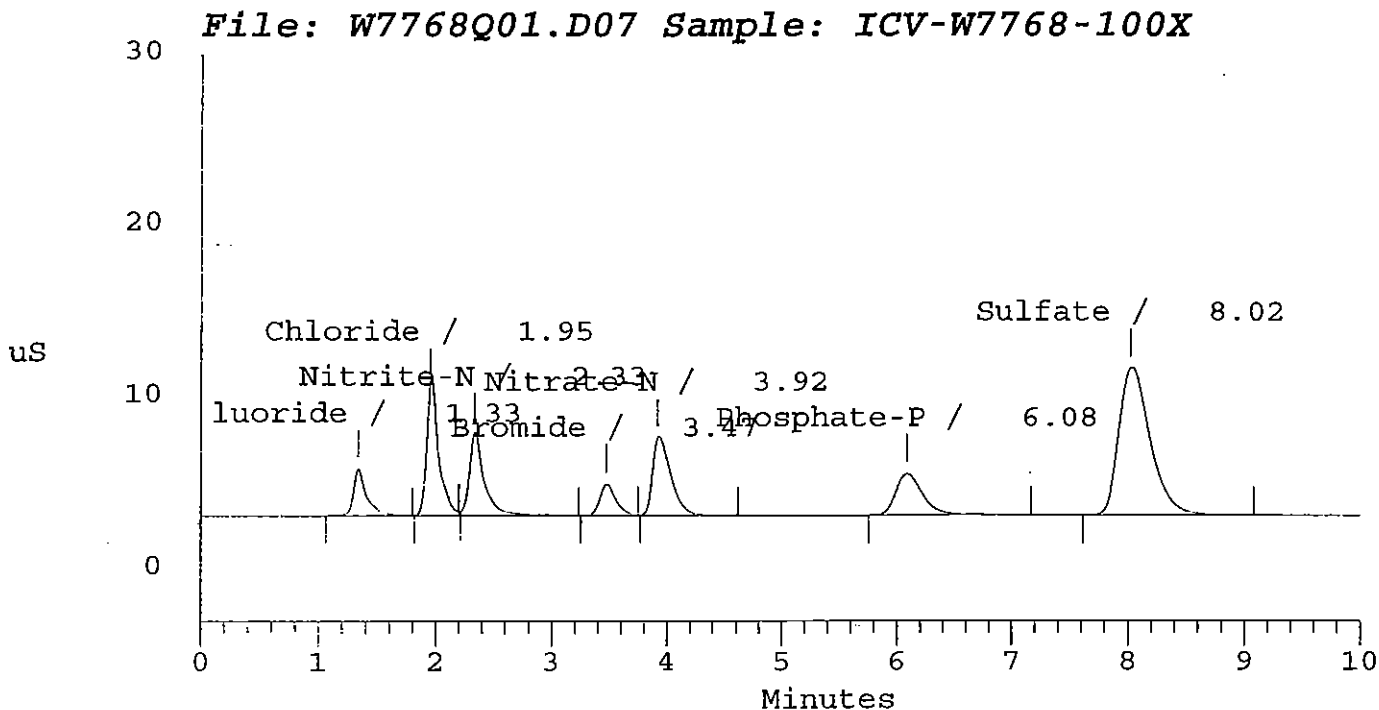
```

-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External           1           1    3000  5Hz   0.00 10.00      1000
-----

```

\*\*\*\*\* Component Report: All Components \*\*\*\*\*

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.33	Fluoride	0.999	89896	726672	2	0.00
2	1.95	Chloride	3.767	248158	1863483	2	0.00
3	2.33	Nitrite-N	1.445	160603	1455681	2	1.17
4	3.47	Bromide	2.881	61007	586729	2	-0.71
5	3.92	Nitrate-N	1.410	149958	1697030	2	0.00
6	6.08	Phosphate-P	2.852	79916	1302965	1	0.00
7	8.02	Sulfate	13.993	285121	5314983	1	0.00
Totals			27.347	1074659	12947545		



```

=====
Sample Name: ICB                               Date: 03/21/2003 18:21:00
Data File  : C:\DX\DATA\E300-063\W7767Q01.D08
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 8           Detector: COND
Analyst    : David                           Column: Dionex AS4A-SC
=====

```

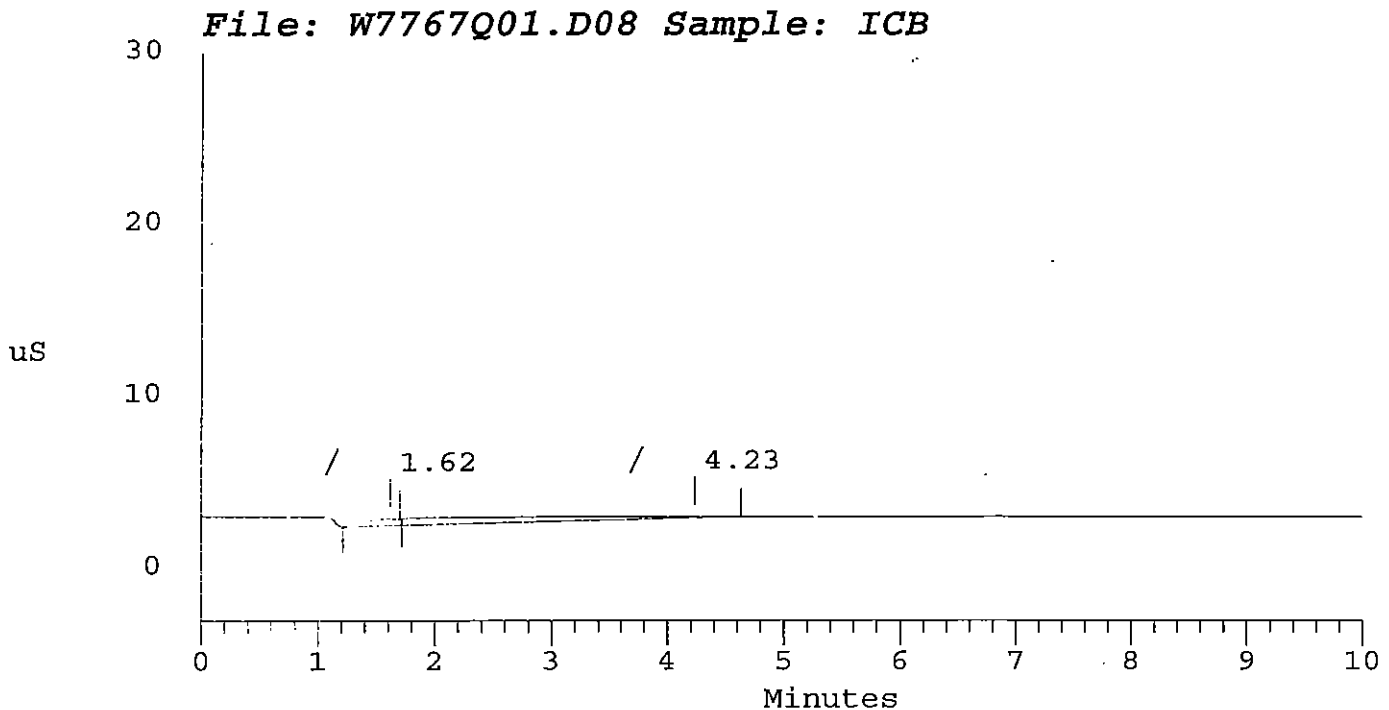
```

-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External          1           1    3000  5Hz   0.00 10.00      1000
-----

```

\*\*\*\*\* Component Report: All Components \*\*\*\*\*

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
0	0.00	Fluoride	0.000	0	0	0	0.00
0	0.00	Chloride	0.000	0	0	0	0.00
0	0.00	Nitrite-N	0.000	0	0	0	0.00
0	0.00	Bromide	0.000	0	0	0	0.00
0	0.00	Nitrate-N	0.000	0	0	0	0.00
0	0.00	Phosphate-P	0.000	0	0	0	0.00
0	0.00	Sulfate	0.000	0	0	0	0.00
Totals			0.000	0	0		



FORM-7  
Applied P & Ch Laboratory  
CCV Recovery for Wet Analysis

Client Name: GEOFON, Inc.  
Case No:  
Project ID: JPL

Contract No.:  
SAS No.:  
Project No.: 04-4428.10

Lab Code: APCL  
Service ID: 33082

#	Component Name	Method	Batch No.	Unit	Expected	Test Result	Rec. %	Dev. %	Flag	Control Limit, %	Test Date
1	Chloride Cl <sup>-</sup>	300.0	03W2750	mg/L	4.0	4.29	107	7	✓	90-110	05/07/2003
	NITRATE as N-NO <sub>3</sub> <sup>-</sup> , BY	300.0	03W2750	mg/L	1.5	1.56	104	4	✓	90-110	05/07/2003
	SULFATE SO <sub>4</sub> <sup>-</sup> , BY I	300.0	03W2750	mg/L	15	15.3	102	2	✓	90-110	05/07/2003
	Chloride Cl <sup>-</sup>	300.0	03W2750	mg/L	4.0	4.20	105	5	✓	90-110	05/07/2003
	NITRATE as N-NO <sub>3</sub> <sup>-</sup> , BY	300.0	03W2750	mg/L	1.5	1.57	104	4	✓	90-110	05/07/2003
	SULFATE SO <sub>4</sub> <sup>-</sup> , BY I	300.0	03W2750	mg/L	15	15.6	104	4	✓	90-110	05/07/2003
	Chloride Cl <sup>-</sup>	300.0	03W2750	mg/L	4.0	4.06	102	2	✓	90-110	05/07/2003
	NITRATE as N-NO <sub>3</sub> <sup>-</sup> , BY	300.0	03W2750	mg/L	1.5	1.58	105	5	✓	90-110	05/07/2003
	SULFATE SO <sub>4</sub> <sup>-</sup> , BY I	300.0	03W2750	mg/L	15	15.5	104	4	✓	90-110	05/07/2003
2	Perchlorate	314.0	03W2804	μg/L	50	49.9	100	0	✓	90-110	05/09/2003
	Perchlorate	314.0	03W2804	μg/L	50	51.1	102	2	✓	90-110	05/09/2003
	Perchlorate	314.0	03W2804	μg/L	50	51.1	102	2	✓	90-110	05/09/2003
	Perchlorate	314.0	03W2804	μg/L	50	52.3	105	5	✓	90-110	05/09/2003
3	Chromium (VI)	7196	03W2737	mg/L	0.25	0.244	98	-2	✓	90-110	05/06/2003
	Chromium (VI)	7196	03W2737	mg/L	0.25	0.255	102	2	✓	90-110	05/06/2003

13760 Magnolia Ave. Chino CA 91710  
 Tel: (909) 590-1828 Fax: (909) 590-1498

pH ( 150.1/9045B ) Worksheet

Temperature compensation must be performed by the instrument automatically.

Analyst W

SOP: G-44

Batch # <u>03W2740</u> Analysis Date: <u>5.6.03</u>	Batch # <u>03W2740</u> Analysis Date: <u>5.6.03</u>
Starting Time: <u>11:37</u> Ending Time: _____	Starting Time: <u>16:19</u> Ending Time: _____
Matrix <input type="checkbox"/> Aqueous <input checked="" type="checkbox"/> Soil	Matrix <input checked="" type="checkbox"/> Aqueous <input type="checkbox"/> Soil

Standard	4.00	7.00	10.00	Standard	4.00	7.00	10.00
Lot #		<u>2120</u>	<u>030619-24</u>	Lot #		<u>2120</u>	<u>030619-24</u>
Temperature °C		<u>24.2</u>	<u>24.2</u>	Temperature °C		<u>24.0</u>	<u>24.0</u>
pH Reading		<u>7.01</u>	<u>10.00</u>	pH Reading		<u>7.02</u>	<u>10.02</u>
T-corrected pH		<u>7.00</u>	<u>10.07</u>	T-corrected pH		<u>7.00</u>	<u>10.07</u>

Control Limit ±0.05 pH unit

#	Sample ID	Pre-treat	pH	Note	#	Sample ID	Pre-treat	pH	Note
MB	<u>0877418</u>		<u>6.86</u>	<u>1:1</u>	MB	<u>T1115</u>		<u>6.84</u>	
1	<u>306-1</u>		<u>7.66</u>	↓	1	<u>3082-1</u>		<u>6.95</u>	
2	<u>-2</u>		<u>7.55</u>		2	<u>-2</u>		<u>7.85</u>	
3	<u>-3</u>		<u>7.95</u>		3	<u>-3</u>		<u>8.08</u>	
4	<u>-4</u>		<u>7.88</u>		4	<u>-4</u>		<u>7.95</u>	
5	<u>2979-1</u>				5	<u>-5</u>		<u>8.77</u>	
6	<u>-2</u>	<u>5/6/03</u>		6	<u>-6</u>		<u>8.43</u>		
7	<u>-3</u>	<u>22</u>		7	<u>3090-1</u>		<u>7.11</u>		
8	<u>-4</u>			8	<u>-7</u>		<u>7.22</u>		
9	<u>-5</u>			9	<u>-3</u>		<u>7.62</u>		
10	<u>-6</u>			10	<u>-4</u>		<u>8.11</u>		
	<u>3000-5</u>			11					
				12					
				13					
				14					
				15					
				16					
				17					
				18					
				19					
				20					
	<u>306-1</u>		<u>7.72</u>	<u>1:1</u>	Dup.	<u>3090-1</u>		<u>7.15</u>	



Alkalinity/OH/CO<sub>3</sub>/HCO<sub>3</sub> (310.1/SM12320B) Worksheet

Batch # 0302287 Matrix: W Titrant H<sub>2</sub>SO<sub>4</sub> Lot # 127900 Concentration (C) 0.025N N: Test Date: 1/6/03 Analyst: AB SOP: G-51

#	Sample ID	Dilution V <sub>1</sub> /V <sub>2</sub> =f <sub>1</sub>	Smp Amnt V, mL	H <sub>2</sub> SO <sub>4</sub> (mL) by Pnhh S <sub>A</sub>	H <sub>2</sub> SO <sub>4</sub> (mL) by MR-BCG S <sub>B</sub>	Pnhh-Alk., P	Toe. Alk., T (in unit of mgCaCO <sub>3</sub> /L)	OH <sup>-</sup>	CO <sub>3</sub> <sup>2-</sup>	HCO <sub>3</sub> <sup>-</sup>	Note & Anomaly
1	MB: T1116	1 =	100	0	0		0	0	0	0	
2	WLS	1 =	100		7.90		100.9				
3	CC57	1 =	100		2.90		102.9				
4	3082-1	1 =	100	0	0		0	0	0		
5	-2	1 =	100	0	8.15		208.2	0	0	208.2	
6	-3	1 =	100	0.05	6.75	1.3	172.7	0	2.6	171.1	
7	-4	1 =	100	0	6.90		176.3	0	0	176.3	
8	-5	1 =	100	0.40	<del>7.5</del> 3.15 4.95	10.2	100.9	0	20.4	80.5	
9	-6	1 =	100	0.10	4.15	2.6	118.8	0	5.2	113.6	
10	3130-1	1 =	100	0	0		0	0	0		
11	-2	1 =	100	0	10.60		270.8	0	0	270.8	
12	-3	1 =	100	0.05	5.85	1.3	170.7	0	2.6	148.1	
13	-4	1 =	100	0.10	5.20	2.6	135.4	0	5.2	130.2	
14	-5	1 =	100	0.05	5.70	1.3	146.9	0	2.6	144.3	
15	-6	1 =	100	1.50	3.65	3.83	126.5	0	76.6	49.9	
16		1 =									
17		1 =									
18		1 =									
19		1 =									
20		1 =									
Dup.	3082-2	1 =	100	0	8.00		209.5	0	0	209.5	

Titration Results	OH <sup>-</sup> (CaCO <sub>3</sub> mg/L)	CO <sub>3</sub> <sup>2-</sup> (CaCO <sub>3</sub> mg/L)	HCO <sub>3</sub> <sup>-</sup> (CaCO <sub>3</sub> mg/L)
P=0	0	0	T
P<T/2	0	2P	T-2P
P=T/2	0	2P	0
P>T/2	2P-T	2(T-P)	0
P=T	T	0	0

Calculation:  
 A=S<sub>A</sub>-E<sub>A</sub>  
 B=S<sub>B</sub>-E<sub>B</sub>  
 P=50,000 f<sub>1</sub> A C / V  
 T=50,000 f<sub>1</sub> (A+B) C / V

APCL Form 5-101, Nov. 29, 1999 Ver 3.2  
 Using blue pen. Correcting by red pen.  
 File: [CUST.DOC.WEFT]ALK.TEX  
 Root-File: [CUST.DOC.WEFT]ALK.ROOT.TEX  
 1-Page-File: [CUST.DOC.WEFT]ALK1.TEX

13760 Magnolia Ave. Chino CA 91710  
Tel: (909) 590-1828 Fax: (909) 590-1498

# Chromium (VI) ( 7196 ) Worksheet

Batch # 03W 2737 Matrix: W

[ Holding Time: 24 hours!! ]

Test Date: 5/6/03 Analyst: hr

Lot #: Reagent Water \_\_\_\_\_ Diphenylcazide solution \_\_\_\_\_

Test Time: 13:25 SOP: G-22

Calibration	STD Lot #	$C_{std} \times V_{std} / V_f = C_i$	$A_i$	$RF_i = A_i / C_i$	Calibration results	Note
STD-1	W-	x / = mg/L			Least Square [RF]=	Cal. Code:
STD-2	W-	x / = mg/L			Average RF=	
STD-3	W-	x / = mg/L			C.C. = 0.999 (> 0.995)	
STD-4	W-	x / = mg/L			RSD= % (< 15%)	
STD-5	W-	x / = mg/L			Ref. page	
STD-6	W-	x / = mg/L			$A = 0.000 + 0.826C$	

Analysis Type	Sample ID or Lot #	Samp. Amnt $X_0$ (g or mL)	Dilu./Ext $X/X_0 = f_1$	Treat. Ratio $V/X = f_2$	- 540 nm A	Concentration $C' = A / RF$	C (Sample) $C = f_1 f_2 C'$	Anomaly Note
CCV	Lot: W- <u>7853</u>	Expected Conc.: x	/	= <u>0.25</u> mg/L	<u>0.204</u>	<u>0.204</u> mg/L	REC. %	90-110 %
Method Blank	Bl. Lot: <u>1146</u>		$X_0 = 1$	95.0/ =	<u>0.000</u>	mg/L	<u>0.000</u> ppm	
LCS1	Bl. Lot: <u>1146</u>		$X_0 = 1$	95.0/ =	<u>0.211</u>	mg/L	<u>0.252</u> ppm	
Sample-1	<u>3082-1</u>		$X_0 = 1$	95.0/ =	<u>0.000</u>	mg/L	<u>0.000</u> ppm	
MS on S-1	<u>4</u>		$X_0 = 1$	95.0/ =	<u>0.193</u>	mg/L	<u>0.231</u> ppm	
MSD on S-1	<u>4</u>		$X_0 = 1$	95.0/ =	<u>0.190</u>	mg/L	<u>0.227</u> ppm	
Sample 2	<u>2</u>		$X_0 = 1$	95.0/ =	<u>0.003</u>	mg/L	<u>0.004</u> ppm	
Sample 3	<u>3</u>		$X_0 = 1$	95.0/ =	<u>0.003</u>	mg/L	<u>0.004</u> ppm	
Sample 4	<u>4</u>		$X_0 = 1$	95.0/ =	<u>0.003</u>	mg/L	<u>0.004</u> ppm	
Sample 5	<u>5</u>		$X_0 = 1$	95.0/ =	<u>0.001</u>	mg/L	<u>0.001</u> ppm	
Sample 6	<u>6</u>		$X_0 = 1$	95.0/ =	<u>0.001</u>	mg/L	<u>0.001</u> ppm	
Sample 7			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 8			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 9			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 10			$X_0 = 1$	95.0/ =		mg/L	ppm	
Blank	Lot:		$X_0 = 1$	95.0/ =		mg/L	ppm	
LCS2	Bl. Lot: <u>1146</u>		$X_0 = 1$	95.0/ =	<u>0.206</u>	mg/L	<u>0.246</u> ppm	
Sample 11			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 12			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 13			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 14			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 15			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 16			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 17			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 18			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 19			$X_0 = 1$	95.0/ =		mg/L	ppm	
Sample 20			$X_0 = 1$	95.0/ =		mg/L	ppm	
MTX Dup.	<u>0.25</u>		$X_0 = 1$	95.0/ =	<u>0.203</u>	<u>0.203</u> mg/L	ppm	

Type	STD Lot #	$C_{STD} (\mu\text{g/mL}) \times V_{STD} (\text{mL}) / X (\text{g or mL}) = T$	Spike Rec.	Ctl Limit (W/S)	PQL/MDL (in ppm)
MS	W- <u>7853</u>	x / = <u>0.25</u> ppm	%	80-120 %/80-120 %	PQL(w) 0.01
MSD	W- <u>1</u>	x / = ppm	%	.. ..	PQL(s) 0.05
LCS	W- <u>7759</u>	x / = ppm	%	80-120 %/80-120 %	MDL(w) 0.005
LCSD	W- <u>1</u>	x / = ppm	%	.. ..	MDL(s) 0.025

Batch # 03W1295 Matrix: W

[ Holding Time: 24 hours!! ]

Test Date: 1/29/03 Analyst: br

Lot #: Reagent Water \_\_\_\_\_ Diphenylcazide solution \_\_\_\_\_ Test Time: \_\_\_\_\_ SOP: G-22

Calibration	STD Lot #	$C_{std} \times V_{std} / V_f = C_i$	$A_i$	$RF_i = A_i / C_i$	Calibration results	Note
STD-1	W-7191	x / = 0.000 mg/L	0.000		Least Square [RF]=	Cal. Code:
STD-2	W-	x / = 0.012 mg/L	0.006		Average RF=	A=0.000+0.836C
STD-3	W-	x / = 0.030 mg/L	0.015		C.C.=0.997 (> 0.995)	
STD-4	W-	x / = 0.115 mg/L	0.109		RSD= % (< 15%)	
STD-5	W-	x / = 0.250 mg/L	0.214		Ref. page	
STD-6	W- ✓	x / = 0.50 mg/L	0.415			A=0.003+0.836C

Analysis Type	Sample ID or Lot #	Samp. Amnt $X_0$ (g or mL)	Dilu./Ext $X/X_0=f_1$	Treat. Ratio $V/X=f_2$	540 nm A	Concentration $C'=A/RF$	C (Sample) $C=f_1 f_2 C'$	Anomaly Note
CCV	Lot: W-7076	Expected Conc.: x	/	= 0.25 mg/L	0.216	0.258 mg/L	REC. %	90-110 %
Method Blank	Bl. Lot: T1115		$1/X_0 = 1$	95.0/ =	0.000	mg/L	0.00 ppm	
LCS1	Bl. Lot: 1		$1/X_0 =$	95.0/ =	0.204	mg/L	0.244 ppm	
Sample-1	1369-1		$1/X_0 =$	95.0/ =	0.000	mg/L	0.00 ppm	
MS on S-1	6		$1/X_0 =$	95.0/ =	0.223	mg/L	0.266 ppm	
MSD on S-1	6		$1/X_0 =$	95.0/ =	0.230	mg/L	0.275 ppm	
Sample 2	2		$1/X_0 =$	95.0/ =	0.004	mg/L	0.005 ppm	
Sample 3	3		$1/X_0 =$	95.0/ =	0.002	mg/L	0.002 ppm	
Sample 4	4		$1/X_0 =$	95.0/ =	0.001	mg/L	0.001 ppm	
Sample 5	5		$1/X_0 =$	95.0/ =	0.002	mg/L	0.002 ppm	
Sample 6	6		$1/X_0 =$ ✓	95.0/ =	0.004	mg/L	0.005 ppm	
Sample 7			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 8			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 9			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 10			$1/X_0 =$	95.0/ =		mg/L	ppm	
Blank	Lot:		$1/X_0 =$	95.0/ =		mg/L	ppm	
LCS2	Bl. Lot: T1115		$1/X_0 = 1$	95.0/ =	0.210	mg/L	0.251 ppm	
Sample 11			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 12			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 13			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 14			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 15			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 16			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 17			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 18			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 19			$1/X_0 =$	95.0/ =		mg/L	ppm	
Sample 20			$1/X_0 =$	95.0/ =		mg/L	ppm	
MTX Dup.	losing 0.25 mg/L		$1/X_0 =$	95.0/ =	0.204	mg/L	0.204 ppm	

Type	STD Lot #	$C_{STD}(\mu\text{g/mL}) \times V_{STD}(\text{mL}) / X(\text{g or mL}) = T$	Spike Rec.	Ctl Limit (W/S)	PQL/MDL (in ppm)
MS	W-7076	x / = 0.25 ppm	%	80-120 %/80-120 %	PQL(w) 0.01
MSD	W- ✓	x / = ppm	%	.. ..	PQL(s) 0.05
LCS	W-7191	x / = ppm	%	80-120 %/80-120 %	MDL(w) 0.005
LCS D	W- ✓	x / = ppm	%	.. ..	MDL(s) 0.025

13760 Magnolia Ave. Chino CA 91710

Solid Analysis ( 160.1, 160.2, 160.3 ) Worksheet

Tel: (909) 590-1828 Fax: (909) 590-1498

Batch # 13W2783 Matrix W Method: 160.1 Balance No. \_\_\_\_\_ Date: 5/8/03 Analyst: DK

EPA 160.1 TDS - Total Dissolved (filterable) Solids - Dry for 1hr. or more at 180 °C

EPA 160.2 TSS - Total Suspended (nonfilterable) Solids - Dry for 1hr. or more at 103-105 °C

EPA 160.3 TS - Total Solids - Dry for 1hr. or more at 103-105 °C

Other method (specify):

Result =  $10^6 \times \Delta W \times f_1 / V$

SOP: G-81

#	Analysis Type	Sample ID (STD Lot #)	Treatment Ratio $V_1/X=f_1$	Volume $V$ , mL	W <sub>1</sub> g	W <sub>2</sub> 1st, g	W <sub>2</sub> 2nd, g	$\Delta W = W_2 - W_1$ , g	Results (ppm)	Note
1	Blank	T116	1 =	100	115.3129	115.3129	115.3130	0.0001	1	38
2	LCS	T116	1 =	100	114.1699	114.2106	114.2104	0.0412	412	41
3	Sample-1	3113-2	1 =	100	108.5429	108.5762	108.5764	0.0335	335	1P
4	MS on S-1	3082-4	1 =	100	115.4120	115.1752	115.1757	0.0631	631	G
5	MSD on S-1	↓ -4	1 =	100	114.7521	114.8142	114.8140	0.0619	619	XX
6	Sample-2	3113-7	1 =	100	112.2714	112.3079	112.3080	0.0364	364	24
7	Sample-3	3105-1	1 =	100	114.7844	114.8381	114.8380	0.0536	536	12
8	Sample-4	↓ -2	1 =	100	111.7980	111.8336	111.8334	0.0354	354	R
9	Sample-5	↓ -3	1 =	100	112.9022	112.9440	112.9439	0.0417	417	木
10	Sample-6	↓ -4	1 =	100	117.8990	117.9218	117.9217	0.0227	227	7
11	Sample-7	3202-1	1 =	100	116.0654	116.0921	116.0920	0.0266	266	31
12	Sample-8	.2	1 =	100	117.6086	117.6089	117.6088	0.0002	2	XS
13	Sample-9	.3	1 =	100	116.9500	116.9869	116.9868	0.0368	368	W
14	Sample-10	.4	1 =	100	99.0897	99.1215	99.1213	0.0316	316	1/2
15	LCSD	T116	1 =	100	103.9474	103.9866	103.9864	0.0390	390	14
16	Sample-11	3102-5	1 =	100	103.5398	103.5647	103.5646	0.0248	248	76
17	Sample-12	.6	1 =	100	114.3109	114.3387	114.3386	0.0277	277	17
18	Sample-13	.7	1 =	100	105.3243	105.3493	105.3490	0.0250	250	12
19	Sample-14	3090-1	3082-1	100	116.6615	116.6617	116.6616	0.0001	1	25
20	Sample-15	3082-2	1 =	100	115.1376	115.1698	115.1699	0.0323	323	10
21	Sample-16	↓ -3	1 =	100	115.2417	115.2625	115.2622	0.0215	215	J
22	Sample-17	↓ -4	1 =	100	107.7129	107.7354	107.7357	0.0222	222	0
23	Sample-18	3086-1	5/8/03/02	100	105.9092	105.9845	105.9848	0.0756	756	19
24	Sample-19	3082-5	1 =	100	121.3345	121.3466	121.3465	0.0120	120	I
25	Sample-20	↓ -6	1 =	100	115.8921	115.9079	115.9077	0.0156	156	CK
	Mix Dup.		1 =							

Std	STD Lot #	$C_{STD}(\mu\text{s/mL}) \times V_{STD}(\text{mL}) / X(\text{g or mL}) = T$	Spike Rec.	Ctl Limit (W/S)	PQL/MDL (in ppm)
W-	7618	x 1 = 400 ppm	%	85-115 %/80-120 %	PQL(w) 10
W-	↓	x 1 = 1 ppm	%	.. ..	PQL(s) 50
W-	7619	x 1 = 1 ppm	%	90-110 %/85-115 %	MDL(w) 4
W-	↓	x 1 = 1 ppm	%	.. ..	MDL(s) 20

# Balance Daily Calibration Worksheet

Weight Set S/N: 12006

Calib. Date	Lab Balance					Digital Balance					Analytical Balance					Calib. by
	Balance #	1 g ±0.05g	10 g ±0.1g	200 g ±0.5g	Note (C)	Balance #	1 g ±0.02g	10 g ±0.05g	200 g ±0.10g	Note (D) (C) (AR)	Balance #	1 g ±0.0002g	10 g ±0.0005g	200 g ±0.0010g	Note (D) (C) (AR)	
5/7/03	A-01	Met	in	Use	✓	B-01	1.00	10.01	200.00	✓✓✓	C-01	1.0000	10.0001	200.0000	✓✓✓	
	<del>A-02</del>					<del>B-05</del>	<del>1.00</del>	<del>10.00</del>	<del>200.01</del>	<del>✓✓✓</del>	<del>C-02</del>	<del>1.0001</del>	<del>10.0000</del>	<del>200.0001</del>	<del>✓✓✓</del>	
	A-03	1.00	9.99	200.00	✓	B-06	1.00	9.99	200.00	✓✓✓	C-					
	A-04					B-07	1.00	9.99	200.00	✓✓✓	C-					
	A-					B-					C-					
5/8/03	A-01	Met	in	Use	✓	B-01	1.00	10.01	200.00	✓✓✓	C-01	1.0000	10.0000	200.0001	✓✓✓	
	<del>A-02</del>					<del>B-05</del>	<del>1.00</del>	<del>10.00</del>	<del>200.00</del>	<del>✓✓✓</del>	<del>C-02</del>	<del>1.0001</del>	<del>10.0001</del>	<del>199.9998</del>	<del>✓✓✓</del>	
	A-03	1.00	9.99	200.00	✓	B-06	1.00	9.99	199.98	✓✓✓	C-					
	A-04					B-07	1.00	10.00	200.00	✓✓✓	C-					
	A-					B-					C-					
5/9/03	A-01	Met	in	Use	✓	B-01	1.00	10.01	200.00	✓✓✓	C-01	1.0000	10.0000	200.0001	✓✓✓	
	<del>A-02</del>					<del>B-05</del>	<del>1.00</del>	<del>10.00</del>	<del>200.00</del>	<del>✓✓✓</del>	<del>C-02</del>	<del>1.0001</del>	<del>10.0001</del>	<del>199.9998</del>	<del>✓✓✓</del>	
	A-03	1.00	9.99	200.00	✓	B-06	1.00	9.99	199.98	✓✓✓	C-					
	A-04					B-07	1.00	10.00	200.00	✓✓✓	C-					
	A-					B-					C-					
5/12/03	A-01	Met	in	Use	✓	B-01	1.00	10.01	200.00	✓✓✓	C-01	1.0000	10.0000	200.0001	✓✓✓	
	<del>A-02</del>					<del>B-05</del>	<del>1.00</del>	<del>10.00</del>	<del>200.00</del>	<del>✓✓✓</del>	<del>C-02</del>	<del>1.0001</del>	<del>10.0001</del>	<del>199.9998</del>	<del>✓✓✓</del>	
	A-03	1.00	9.99	200.00	✓	B-06	1.00	9.99	199.98	✓✓✓	C-					
	A-04					B-07	1.00	10.00	200.00	✓✓✓	C-					
	A-					B-					C-					

Notation: (C) - Cleanliness; (D) - Display; (AR) - Auto Rerzeroing.  
 APCI form 4-213, March 30, 1995. Ver. 4.0 No pencil. Use blue pen for record. Use red pen for correction.  
 File: [CUST.DOC;LAB]BAL-CAL-TEX Root-File: BAL-CAL-ROOT-TEX 1-Page-File: BAL-CAL11-TEX

Line	Sample	Sample Type	Level	Method	Data File	Volume	Dilution
1	Cal blank	Sample		e314-011.met	c:\data\314-01\mb_001.dxd	1	1
2	cal standard 2ppb W7827a	Sample		e314-011.met	c:\data\314-01\std-2pb_002.dxd	1	1
3	cal standard 4ppb W7827b	Sample		e314-011.met	c:\data\314-01\std-4pb_003.dxd	1	1
4	cal standard 10ppb W7827c	Sample		e314-011.met	c:\data\314-01\std-10pb_004.dxd	1	1
5	cal standard 25ppb W7827d	Sample		e314-011.met	c:\data\314-01\std-25pb_005.dxd	1	1
6	cal standard 50ppb W7827e	Sample		e314-011.met	c:\data\314-01\std-50pb_006.dxd	1	1
7	cal standard 75ppb W7827f	Sample		e314-011.met	c:\data\314-01\std-75pb_007.dxd	1	1
8	cal standard 100ppb W7827g	Sample		e314-011.met	c:\data\314-01\std-100pb_008.dxd	1	1
9	ICV 50 ppb w7828a	Sample		e314-011.met	c:\data\314-01\icv-50pb_009.dxd	1	1
10	lcb	Sample		e314-011.met	c:\data\314-01\lcb_010.dxd	1	1
11	anion 100pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-100_011.dxd	1	1
12	anion 200pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-200_012.dxd	1	1
13	anion 300pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-300_013.dxd	1	1
14	anion 400pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-400_014.dxd	1	1
5	anion 500pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-500_015.dxd	1	1
6	anion 600pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-600_016.dxd	1	1
7	anion 800pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-800_017.dxd	1	1
8	anion 1000pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-01\mct-1000_018.dxd	1	1
9	anion 400pm each 2pb	Sample		e314-011.met	c:\data\314-01\ipc-2pb_019.dxd	1	1
0	anion 400pm each 4pb	Sample		e314-011.met	c:\data\314-01\ipc-4pb_020.dxd	1	1
1	anion 400pm each 25pb	Sample		e314-011.met	c:\data\314-01\ipc-25pb_021.dxd	1	1
2	ICV 50 ppb	Sample		e314-011.met	c:\data\314-01\iccv-50pb	1	1
3	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-02_023.dxd	1	1
4	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-03_024.dxd	1	1
5	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-04	1	1
6	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-05	1	1
7	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-06	1	1
8	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-07	1	1
9	MDL 4pb	Sample		e314-011.met	c:\data\314-01\mdl-08	1	1
0	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
1	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
2	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
3	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
4	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
5	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
6	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-01\idap-25pb	1	1
7	MCT anion 800pm each, 25pbCLO4	Sample		e314-011.met	c:\data\314-01\ipo-25pb	1	1
8	MCT anion 800pm each, 25pbCLO4	Sample		e314-011.met	c:\data\314-01\ipc-25pb	1	1
9	MCT anion 800pm each, 4pbCLO4	Sample		e314-011.met	c:\data\314-01\ipc-4pb	1	1
0	MCT anion 800pm each, 4pbCLO4	Sample		e314-011.met	c:\data\314-01\ipc-4pb	1	1
1	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s01	1	5
2	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s02	1	5
3	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s03	1	5
4	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s04	1	5
5	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s05	1	5
6	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s06	1	5
7	MDL 20pb soil	Sample		e314-011.met	c:\data\314-01\mdl-s07	1	5
8	standard 25ppb W7827d	Sample		e314-011.met	c:\data\314-01\std-25pb	1	1
9	anion 100pm each,4pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-100-4pb	1	1
0	anion 200pm each ,4pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-200-4pb	1	1
1	anion 300pm each ,4pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-300-4pb	1	1
2	anion 100pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-100-2pb	1	1
3	anion 200pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-200-2pb	1	1
4	anion 300pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-01\lam-300-2pb	1	1
5	1982-01 B S.C 4450us/cm	Sample		e314-011.met	c:\data\314-01\1982-01	1	1
6	1982-01 B S.C 4450us/cm	Sample		e314-011.met	c:\data\314-01\1982-01	1	2
7	1982-02 f=10	Sample		e314-011.met	c:\data\314-01\1982-02_057.dxd	1	10
8		Sample		aastopcl.met		1	1

ie	Sample	Sample Type	Level	Method	Data File	Volume
	##03w2804kw ipc 25ppb w7759	Sample		e314-011.met	c:\data\03w2804k\w2804k ipc 25ppb_001.dxd	1
	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2804k\w2804k q01	1
	ccb	Sample		e314-011.met	c:\data\03w2804k\w2804k ccb01_003.dxd	1
	lcs 25ppb w7827d	Sample		e314-011.met	c:\data\03w2804k\w2804k l01_004.dxd	1
	LCS 18PPB W7685D	Sample		e314-011.met	c:\data\03w2804k\w2804k j01_005.dxd	1
	ICCS 4ppb w7827b	Sample		e314-011.met	c:\data\03w2804k\w2804k iccs 4ppb_006.dxd	1
	mb	Sample		e314-011.met	c:\data\03w2804k\w2804k k01_007.dxd	1
	3152-01 F=1	Sample		e314-011.met	c:\data\03w2804k\3152-01_008.dxd	1
	3152-02 f=1	Sample		e314-011.met	c:\data\03w2804k\3152-02_009.dxd	1
	3152-03 F=1	Sample		e314-011.met	c:\data\03w2804k\3152-03_010.dxd	1
	3152-04 F=1	Sample		e314-011.met	c:\data\03w2804k\3152-04_011.dxd	1
	3152-05 f=1	Sample		e314-011.met	c:\data\03w2804k\3152-05_012.dxd	1
	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2804k\w2804k q02_013.dxd	1
	ccb	Sample		e314-011.met	c:\data\03w2804k\w2804k k02_014.dxd	1
	3152-06 f=1	Sample		e314-011.met	c:\data\03w2804k\3152-06_015.dxd	1
	3082-04 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-04_016.dxd	1
	3082-01 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-01_017.dxd	1
	3082-04 MS 50PPB F=1	Sample		e314-011.met	c:\data\03w2804k\w2804k m01_018.dxd	1
	3082-04 MSD 50PPB F=1	Sample		e314-011.met	c:\data\03w2804k\w2804k n01_019.dxd	1
	3082-02 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-02_020.dxd	1
	3082-03 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-03_021.dxd	1
	3082-05 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-05_022.dxd	1
	3082-06 F=1	Sample		e314-011.met	c:\data\03w2804k\3082-06_023.dxd	1
	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2804k\w2804k q03_024.dxd	1
	CCB	Sample		e314-011.met	c:\data\03w2804k\w2804k k03_025.dxd	1
	3102-01 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-01_026.dxd	1
	3102-02 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-02_027.dxd	1
	3102-03 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-03_028.dxd	1
	3102-04 f=1	Sample		e314-011.met	c:\data\03w2804k\3102-04_029.dxd	1
	3102-05 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-05_030.dxd	1
	3102-06 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-06_031.dxd	1
	3102-07 F=1	Sample		e314-011.met	c:\data\03w2804k\3102-07_032.dxd	1
	CCV 50PPB W7827E	Sample		e314-011.met	c:\data\03w2804k\w2804k q04_033.dxd	1
		Sample		aastopl.met		1

Analyst Wei Way  
 Date 5/9/03  
 Instrument ZC-1c

DIONEX SCHEDULE - C:\DX\SCHEDULE\E300-063.SCH

Inj#	Sample Name	Method	Data File	Vol.	Dil.	Int.Std.
1	autocal1r	..\E300-063	..\W7767Q01.D01	1	1	1
2	autocal2r	..\E300-063	..\W7767Q01.D02	1	1	1
3	autocal3r	..\E300-063	..\W7767Q01.D03	1	1	1
4	autocal4r	..\E300-063	..\W7767Q01.D04	1	1	1
5	autocal5r	..\E300-063	..\W7767Q01.D05	1	1	1
6	autocal6r	..\E300-063	..\W7767Q01.D06	1	1	1
7	icv-w7768-100X	..\E300-063	..\W7768Q01.D07	1	1	1
8	icb	..\E300-063	..\W7767Q01.D08	1	1	1

Comment :

Analyst       *DN*        
 Date       3/21/03        
 Instrument       J



DIONEX SCHEDULE - C:\DX\SCHEDULE\03W2750.SCH

Inj#	Sample Name	Method	Data File	Vol.	Dil.	Int.Std.
1	##03W2750, W CCVW77	..\E300-063	..\W2750Q01.D01	1	1	1
2	MB RW1410	..\E300-063	..\W2750K01.D02	1	1	1
3	LCS W7768-100X	..\E300-063	..\W2750L01.D03	1	1	1
4	LCSD W7768-100X	..\E300-063	..\W2750J01.D04	1	1	1
5	3082-2 F=2.5	..\E300-063	..\3082-201.D05	1	2.5	1
6	3082-3 F=2.5	..\E300-063	..\3082-301.D06	1	2.5	1
7	3082-4 F=2.5	..\E300-063	..\3082-401.D07	1	2.5	1
8	3082-5 F=2	..\E300-063	..\3082-501.D08	1	2	1
9	3082-6 F=2	..\E300-063	..\3082-601.D09	1	2	1
10	3090-1 F=4	..\E300-063	..\3090-101.D10	1	4	1
11	3082-1 F=1.25	..\E300-063	..\3082-101.D11	1	1.25	1
12	CCV2W7767-100X	..\E300-063	..\W2750Q01.D12	1	1	1
13	MB RW1410	..\E300-063	..\W2750K01.D13	1	1	1
14	\$3082-4 MS F=5	..\E300-063	..\W2750M01.D14	1	5	1
15	\$3082-4 MSD F=5	..\E300-063	..\W2750N01.D15	1	5	1
16	3090-2 F=4	..\E300-063	..\3090-201.D16	1	4	1
17	3090-3 F=5	..\E300-063	..\3090-301.D17	1	5	1
18	3090-4 F=8	..\E300-063	..\3090-401.D18	1	8	1
19	3090-4 MD F=8	..\E300-063	..\W2750D01.D19	1	8	1
20	CCV3W7767-100X	..\E300-063	..\W2750Q11.D20	1	1	1
21		..\STOP.MET		1	1	1

Comment:

LCS/LCSD LOT # W7768

MS/MSD LOT # W7767

ELUENT LOT # W7868

ANALYTICAL METHOD 9056/E300 MATRIX W

Analyst DZ

Date 5/2/03

Instrument J



## Applied Physics & Chemistry Laboratory

13760 Magnolia Ave. Chino CA 91710  
Tel. (909) 590-1828 Fax (909) 590-1498

June 6, 2003

GEOFON, Inc.  
Attention: Leo Williamson  
22632 Golden Spring Dr Ste 270  
Diamond Bar CA 91765

Dear Leo Williamson,

This package contains samples in our Service ID 03-3102 and your project : 04-4428.10 JPL GW  
Mon 2Q03.

Enclosed please find:

- (1) Original analytical report.
- (2) Original Chain of Custody.
- (3) One diskette containing EDD deliverable.
- (4) One original Level C Data Package Deliverable.

If anything is missing or you have any questions, please feel free to contact me.

Respectfully submitted,

A handwritten signature in black ink, appearing to read 'R. Kirakozova', is written over the typed name.

Regina Kirakozova  
Associate QA/QC Director  
Applied P & Ch Laboratory

Applied P & Ch Laboratory

13760 Magnolia Ave. Chino CA 91710

Tel: (909) 590-1828 Fax: (909) 590-1498

Submitted to:  
 GEOFON, Inc.  
 Attention: Leo W. Williamson  
 22632 Golden Spring Dr Ste 270  
 Diamond Bar CA 91765  
 Tel: (909) 396-7662 Fax: (909) 396-1455

# APCL Analytical Report

Service ID #: 801-033102 Received: 05/07/03  
 Collected by: Leo Williamson Extracted: N/A  
 Collected on: 05/07/03 Tested: 05/07-16/03  
 Reported: 05/21/03  
 Sample Description: Water from NW-12 (E. of Bl 302)  
 Project Description: 04-4428.10 JPL GW Mon 2Q03.

## Analysis of Water Samples

Component Analyzed	Method	Unit	PQL	Analysis Result			
				DUPE-6-2Q03 03-03102-1	EB-11-5/7/03 03-03102-2	MW-12-1 03-03102-3	MW-12-2 03-03102-4
BICARBONATE	SM2320B	mg/L	2	185	<2	199	198
CARBONATE	SM2320B	mg-CaCO <sub>3</sub> /L	2	<2	<2	<2	<2
PH	9040B	pH unit	0.01	7.89	7.73	7.56	7.56
SOLIDS, TOTAL DISSOLVED (TDS)	160.1	mg/L	10	266	<10	368	316
CHROMIUM (VI)	7196	mg/L	0.01	<0.01	<0.01	<0.01	<0.01
Dilution Factor				1	1	1	1
PERCHLORATE	314.0	µg/L	4	3.4J	<4	<4	2.0J
Dilution Factor				25	1.25	4	2.5
CHLORIDE CL <sup>-</sup>	300.0	mg/L	0.2	16.8	0.21J	30.9	17.3
NITRATE AS N	300.0	mg/L	0.04	<1	0.098	1.5	1.5
SULFATE SO <sub>4</sub> <sup>-2</sup>	300.0	mg/L	0.5	32.6	0.71	62.4	39.0
Dilution Factor				1	1	1	1
ARSENIC	200.9	µg/L	5	<5	<5	<5	<5
CALCIUM	200.7	µg/L	200	44,800	<200	60,100	56,400
IRON	200.7	µg/L	50	62.8	16.4J	572	204
MAGNESIUM	200.7	µg/L	100	14,400	25.9J	21,600	17,000
POTASSIUM	200.7	µg/L	400	2,790	70.9J	3,430	2,910
SODIUM	200.7	µg/L	2000	23,400	931J	24,500	23,700
VOLATILE ORGANIC COMPOUNDS							
Dilution Factor				1	1	1	1
BENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOCHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMODICHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOFORM	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2-BUTANONE	524.2	µg/L	10	<10	<10	<10	<10
N-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
SEC-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TERT-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CARBON TETRACHLORIDE	524.2	µg/L	0.5	2.6	<0.5	<0.5	<0.5
CHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLORODIBROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLOROFORM	524.2	µg/L	0.5	1.2	<0.5	<0.5	<0.5
CHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2-CHLOROTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
4-CHLOROTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5

# APCL Analytical Report

Component Analyzed	Method	Unit	PQL	Analysis Result			
				DUPE-6-2Q03	EB-11-5/7/03	MW-12-1	MW-12-2
				03-03102-1	03-03102-2	03-03102-3	03-03102-4
1,2-DIBROMO-3-CHLOROPROPANE	524.2	µg/L	1.1 (a)	<1.1	<1.1	<1.1	<1.1
1,2-DIBROMOETHANE (EDB)	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
DIBROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,3-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,4-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
DICHLORODIFLUOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1-DICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CIS-1,2-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRANS-1,2-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROETHENE (TOTAL)	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,3-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2,2-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CIS-1,3-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRANS-1,3-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
ETHYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
HEXACHLOROBUTADIENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
ISOPROPYLBENZENE (CUMENE)	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
P-ISOPROPYLTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
4-METHYL-2-PENTANONE (MIBK)	524.2	µg/L	10	4J	7J	8J	5J
METHYLENE CHLORIDE	524.2	µg/L	1.8 (a)	<1.8	<1.8	7.2 (b)	10 (b)
METHYL-T-BUTYL ETHER (MTBE)	524.2	µg/L	1	<1	<1	<1	<1
NAPHTHALENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
N-PROPYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
STYRENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,1,2-TETRACHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2,2-TETRACHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TETRACHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,3-TRICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,4-TRICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,1-TRICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2-TRICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRICHLOROFLUOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,3-TRICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,4-TRIMETHYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,3,5-TRIMETHYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
VINYL CHLORIDE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
O-XYLENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
M/P-XYLENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5

# APCL Analytical Report

Component Analyzed	Method	Unit	PQL	Analysis Result			
				MW-12-3 03-03102-5	MW-12-4 03-03102-6	MW-12-5 03-03102-7	TB-11-5/7/03 03-03102-8
BICARBONATE	SM2320B	mg/L	2	180	204	181	-
CARBONATE	SM2320B	mg-CaCO <sub>3</sub> /L	2	< 2	< 2	< 2	-
PH	9040B	pH unit	0.01	8.03	7.88	8.00	-
SOLIDS, TOTAL DISSOLVED (TDS)	160.1	mg/L	10	248	277	252	-
CHROMIUM (VI)	7196	mg/L	0.01	<0.01	<0.01	<0.01	-
Dilution Factor				1	1	1	1
PERCHLORATE	314.0	µg/L	4	2.8J	3.6J	2.2J	-
Dilution Factor				2.5	2.5	2.5	1
CHLORIDE CL <sup>-</sup>	300.0	mg/L	0.2	15.7	14.9	15.6	-
NITRATE AS N	300.0	mg/L	0.04	0.53	1.1	1.1	-
SULFATE SO <sub>4</sub> <sup>-2</sup>	300.0	mg/L	0.5	29.8	30.3	22.0	-
Dilution Factor				1	1	1	1
ARSENIC	200.9	µg/L	5	< 5	< 5	< 5	-
CALCIUM	200.7	µg/L	200	45,300	56,900	39,500	-
IRON	200.7	µg/L	50	64.2	74.4	286	-
MAGNESIUM	200.7	µg/L	100	14,600	13,500	11,100	-
POTASSIUM	200.7	µg/L	400	2,760	2,120	2,080	-
SODIUM	200.7	µg/L	2000	23,600	23,200	35,100	-
<b>VOLATILE ORGANIC COMPOUNDS</b>							
Dilution Factor				1	1	1	1
BENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOCHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMODICHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOFORM	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
BROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2-BUTANONE	524.2	µg/L	10	<10	<10	<10	<10
N-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
SEC-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TERT-BUTYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CARBON TETRACHLORIDE	524.2	µg/L	0.5	2.5	1.5	0.6	<0.5
CHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLORODIBROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CHLOROFORM	524.2	µg/L	0.5	1.1	0.7	0.4J	<0.5
CHLOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2-CHLOROTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
4-CHLOROTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DIBROMO-3-CHLOROPROPANE	524.2	µg/L	1.1 (a)	<1.1	<1.1	<1.1	<1.1
1,2-DIBROMOETHANE (EDB)	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
DIBROMOMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,3-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,4-DICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
DICHLORODIFLUOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5

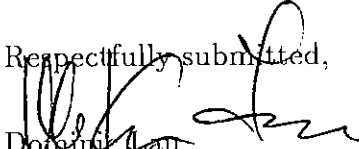
# APCL Analytical Report

Component Analyzed	Method	Unit	PQL	Analysis Result			
				MW-12-3 03-03102-5	MW-12-4 03-03102-6	MW-12-5 03-03102-7	TB-11-5/7/03 03-03102-8
1,1-DICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CIS-1,2-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRANS-1,2-DICHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,3-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
2,2-DICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
CIS-1,3-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRANS-1,3-DICHLOROPROPENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
ETHYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
HEXACHLOROBUTADIENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
ISOPROPYLBENZENE (CUMENE)	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
P-ISOPROPYLTOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
4-METHYL-2-PENTANONE (MIBK)	524.2	µg/L	10	<10	<10	7J	<10
METHYLENE CHLORIDE	524.2	µg/L	1.8 (a)	8.3 (b)	11 (b)	13 (b)	6.8 (b)
METHYL-T-BUTYL ETHER (MTBE)	524.2	µg/L	1	<1	<1	<1	<1
NAPHTHALENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
N-PROPYLBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
STYRENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,1,2-TETRACHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2,2-TETRACHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TETRACHLOROETHENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TOLUENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,3-TRICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,4-TRICHLOROBENZENE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,1-TRICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2-TRICHLOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
TRICHLOROETHENE	524.2	µg/L	0.5	<0.5	0.3J	<0.5	<0.5
TRICHLOROFLUOROMETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,2,3-TRICHLOROPROPANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5
1,1,2,2-TRICHLORO-1,1,2,2-TRIFLUOROETHANE	524.2	µg/L	0.5	<0.5	<0.5	<0.5	<0.5

# APCL Analytical Report

Component Analyzed	Method	Unit	PQL	Analysis Result			
				MW-12-3 03-03102-5	MW-12-4 03-03102-6	MW-12-5 03-03102-7	TB-11-5/7/03 03-03102-8
1,2,4-TRIMETHYLBENZENE	524.2	µg/L	0.5	< 0.5	< 0.5	< 0.5	< 0.5
1,3,5-TRIMETHYLBENZENE	524.2	µg/L	0.5	< 0.5	< 0.5	< 0.5	< 0.5
VINYL CHLORIDE	524.2	µg/L	0.5	< 0.5	< 0.5	< 0.5	< 0.5
O-XYLENE	524.2	µg/L	0.5	< 0.5	< 0.5	< 0.5	< 0.5
M/P-XYLENE	524.2	µg/L	0.5	< 0.5	< 0.5	< 0.5	< 0.5

PQL: Practical Quantitation Limit. MDL: Method Detection Limit. CRDL: Contract Required Detection Limit  
N.D.: Not Detected or less than the practical quantitation limit. "-": Analysis is not required.  
J: Reported between PQL and MDL.  
Listed Dilution Factors (DF) are relative to the method default DF. All unlisted DFs are 1.0  
(<sup>a</sup>) MDL reported.  
(<sup>b</sup>) Laboratory contamination suspected.

Respectfully submitted,  
  
Dominic Lau  
Laboratory Director  
Applied P & Ch Laboratory

**Level C Data Package Deliverables**

# **General Information**

**Project: 04-4428.10 JPL GW Mon-2Q03**

**APCL Service ID: 03-3102**



**Applied Earth & Ch Laboratory**

**13760 Magnolia Ave. Chino, CA 91710**

**Telephone (909)590-1828**

**Fax (909)590-1498**



# Case Narrative

**Project: JPL GW Mon 2Q03/NW-12 E. of Bl 302/04-4428.10**

**For GEOFON, Inc.**

**APCL Service No: 03-3102**

## 1. Sample Identification

The sample identifications are listed in the following table:

GEOFON, Inc. Sample ID	APCL Sample ID
MW-12-5	03-03102-7
MW-12-4	03-03102-6
MW-12-3	03-03102-5
MW-12-2	03-03102-4
MW-12-1	03-03102-3
TB-11-5/7/03	03-03102-8
EB-11-5/7/03	03-03102-2
DUPE-6-2Q03	03-03102-1

## 2. Analytical Methodology

Samples are analyzed by EPA methods

- 524.2 (Volatile Organic Compounds ),
- 7196 (Chromium (VI) ),
- 314.0 (Perchlorate, low level ),
- 300.0 (Anions by IC ),
- SM2320B (Carbonate ),
- SM2320B (Bicarbonate ),
- 9040B (pH ),
- 160.1 (Solids, Total Dissolved (TDS) ),
- 200.7 (Metals by ICP ),
- 200.9 (Arsenic, As, by GFAA ),

## 3. Holding Time

All samples were extracted, digested and analyzed within the holding times defined by the appropriate EPA methods of the analyses.

## 4. Preservation

All samples were preserved and stored according to the appropriate EPA methods.

## 5. Tele-log

None

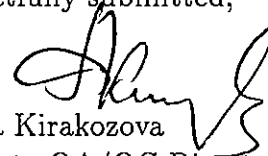
## 6. Anomaly

(1) EPA 524.2:

Methylene Chloride in the amount of 6.7 ug/L was detected in the Method Blank of batch 03G2404. Methylene Chloride was also detected in the associated field samples, due to lab contamination.

"I certify that these data are technically accurate, complete, and in compliance with the terms and conditions of the contract, for other than the conditions detailed above. Release of the data contained in the hardcopy data package and its electronic data deliverable submitted on diskette had been authorized by the Laboratory Manager or her/his designee, as verified by the following signature."

Respectfully submitted,



Regina Kirakozova  
Associate QA/QC Director  
Applied P & Ch Laboratory



INGORPORATED  
2632 GOLDEN SPRINGS DR., SUITE 270  
DIAMOND BAR, CA 91765 • (909) 396-7662 • FAX (909) 396-1455

CHAIN-OF-CUSTODY RECORD

LABORATORY COPY

MW-12 0030

GEOFON, LAB COORDINATOR

LAB COORDINATOR'S PHONE

LAB COORDINATOR'S FAX

LABORATORY SERVICE ID

LABORATORY CONTACT

MAIL REPORT (COMPANY NAME)

Brad Shojaee

(909) 396-7662

(909) 396-1455

—

Kenny Chan

GEOFON, INC.

PROJECT NAME: SDP LWR MW-2903

PROJECT LOCATION: MW-12 (E. of B1302)

PROJECT NUMBER: 04-4428.10

LABORATORY PHONE: (909) 590-1828

LABORATORY FAX: (909) 590-1498

RECIPIENT NAME: Leo W. Williamson

PROJECT CONTACT: Leo W. Williamson

PROJECT PHONE NUMBER: (714) 920-8729

PROJECT FAX: (909) 396-1455

LABORATORY ADDRESS: 13760 Magnolia Ave

CITY, STATE AND ZIP CODE: Chino, CA 91710

ADDRESS: 22632 Golden Springs, Dr. #270

PROJECT ADDRESS: 4800 Oak Grove Dr.

CITY, STATE AND ZIP CODE: Pasadena, CA

CLIENT: US NAVY SWOIC

CITY, STATE AND ZIP CODE: Chino, CA 91710

LABORATORY CONTACT: (909) 396-1455

CITY, STATE AND ZIP CODE: Diamond Bar, CA 91765

PROJECT MANAGER: Asrar Fakhreem

PROJECT PHONE NUMBER: (909) 396-7662

PROJECT MANAGER'S FAX: (909) 396-1455

LABORATORY SERVICE ID

LABORATORY CONTACT

RECIPIENT NAME

Item	Sample Identifier	Matrix	Date	Time	Preserved	# of Cont.	QC Level	T.A.T.
1	MW-12-5	H <sub>2</sub> O	5/7/03	750	3+1+	1+1	III	Normal
2	MW-12-4			830				
3	MW-12-3			940				
4	MW-12-2			1020				
5	MW-12-1			1100				
6								
7	TB-11-5/7/03	H <sub>2</sub> O	—	—	KCl	1	III	Normal
8	EB-11-5/7/03		5/7/03	850	HCl NONE	3+1+	IV	
9	DUPE-6-2903							
10								

ANALYSES: Na/K/Cal/As/My/Ir  
 524.2 (VOCs)  
 200.1 & 200.9 (Minerals)  
 716 (Hex Chrome)  
 SIM 2320B (Carb. Ion)  
 300.0 (Cl Ion)  
 314.0 (Sulfide)  
 160.1 (P)  
 200.8 (Ca P)

3102

SAMPLES COLLECTED BY: Leo W. Williamson

RELINQUISHED BY: [Signature]

COURIER AND AIR BILL NUMBER

RECEIVED BY: [Signature]

DATE: 5/7/03

TIME: 1:30

COOLER TEMPERATURE UPON RECEIPT

SAMPLE'S CONDITION UPON RECEIPT

Distribution: White - Laboratory (To be returned with Analytical Report); Goldenrod - Project File; Yellow - Project Data Manager

# Sample Receiving Checklist

APCL Service ID: **3102** Client Name/Project: Geofon JPL

### 1. Sample Arrival

Date/Time Received 5/7/03 1225 Date/Time Opened 5/7/03 1225 By (name): Kenny Chan  
Custody Transfer:  Client  Golden State  UPS  US Mail  FedEx  APCL Empl: Adam Abad

### 2. Chain-of-Custody (CoC)

With Samples?  Faxed?  Client has Copy?  Signed, dated? By: \_\_\_\_\_  
 Project ID?  Analyses Clear?  Hold Samples? #on Hold \_\_\_\_\_ # Received 8  
 CoC/Docs Zip-Locked under lid?  Compos.#: \_\_\_\_\_  #Samples OK?  
 Discrepancies?  Client notified?  Response (attach docs): \_\_\_\_\_

### 3. Shipping Container/Cooler

Cooler Used? # of 1 Cooled by:  Ice  Blue Ice  Dry Ice  None  
Temp °C 4.1  
(Cooler temperature measured from temp blank if present, otherwise measured from the cooler).  
Cooler Custody Seal?  Absent  Intact  Tampered?

### 4. Sample Preservation

pH <2  pH >12  
If Not, pH = \_\_\_\_\_ Preserved by:  Client  APCL  Third Party \_\_\_\_\_

### 5. Holding-time Requirements

pH 24hr  BACT 6/24hr  Cr<sup>VJ</sup> 24hr  NO<sub>3</sub><sup>-</sup> 48hr  BOD 48hr  
 Cl<sub>2</sub> ASAP  Turbidity 48hr  DO ASAP  Fe(II) ASAP  
 HT Expired?  Client notified?

### 6. Sample Container Condition

Intact?  Broken?  Documented? Number: \_\_\_\_\_  
Type:  plastic  glass  Tube: brass/SS  Tedlar Bag  
 Quantity OK?  Leaking?  Anomaly?  
 Caps tight?  Air Bubbles?  Anomaly?  
Labels:  Unique ID?  Date/Time  Preserved?

### 7. Turn Around Time

RUSH TAT: \_\_\_\_\_  Std (7-10 days)  Not Marked

### 8. Sample Matrix

Drinking H<sub>2</sub>O  Other Liq  Soil  Wipe  Polymer  Air  Other: \_\_\_\_\_  
 Ground H<sub>2</sub>O  Sludge  Filter  Oil/Petro  Paint  W. Water  Extract  Unknown

### 9. Pre-Login Check List Completed & OK?

ALL OK? (if not, attach docs)  Client Contact? (Name: \_\_\_\_\_) Date/Time: \_\_\_\_\_  
Received/Checked by: Kenny Chan Date: 7 May 2003 Time: 7:40 a.m.

\* HT: Samples must be analyzed for results to reflect total concentrations. Results generated outside required of holding times are considered minimal values and may be used to define waste as hazardous but not as non-hazardous.

Applied P & Ch Laboratory

13760 Magnolia Ave. Chino CA 91710

Tel: (909) 590-1828 Fax: (909) 590-1498

# Sample Login: Check List

**03-03102 (0470\_ 139) (2202777\_ 139)**

05/07/03

## Part 1: General Information

---

- |                          |                          |                      |  |
|--------------------------|--------------------------|----------------------|--|
| <input type="checkbox"/> | Company Information      | Name:                | <i>GEOFON, Inc.</i>  |
|                          |                          | Address:             | <i>22632 Golden Spring Dr Ste 270 ,Diamond Bar ,CA 91765</i> |
| <input type="checkbox"/> | Project Information      | Project Description: | <i>JPL</i>   |
|                          |                          | Project #:           | <i>04-4428.10</i>  |
| <input type="checkbox"/> | Billing Information      | P.O. #:              |  |
|                          |                          | Bill Address:        | <i>22632 Golden Spring Dr Ste 270 ,Diamond Bar ,CA 91765</i> |
|                          |                          | Lab Project ID:      |  |
|                          |                          | Client Database #:   | <i>3</i>   |
| <input type="checkbox"/> | Receiving Information    | Who Received Sample? | <i>Kenny Chan</i>  |
|                          |                          | Receiving Date/Time: | <i>05/07/03 1225</i>   |
|                          |                          | COC No.              |  |
| <input type="checkbox"/> | Shipping Information     | Shipping Company     | <i>APCL pick up</i>  |
|                          |                          | Packing Information: | <i>Cooler/Ice Chester</i>                                    |
|                          |                          | Cooler Temperature:  | <i>4.1 °C</i>  |
| <input type="checkbox"/> | Container Information    | Container Provider:  | <i>Client</i>  |
| <input type="checkbox"/> | Sampling Information     | Sampling Person:     |  |
|                          |                          | Sampling Company:    | <i>Client</i>  |
| <input type="checkbox"/> | Turn-Around-Time Option: |                      | <i>Rush 5 working day(s)</i>                                 |
| <input type="checkbox"/> | QC Option:               |                      | <i>NEESA C</i>   |
| <input type="checkbox"/> | Disposal Option:         |                      | <i>Not specify</i>   |
-

## Part 2: Sample Information

Seq. #	Sample ID (on COC)	Sample Sub-ID	APCL Sample ID	Matrix	Cont- tainer	Preser- vative	Vol, ml Am. g	# of Replica	Condition G, L, B	Collected mmdyy	Hold ?	Composite Group	TAT Days
1	MW-12-5	VOC	03-03102-7- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-5	Metal	03-03102-7- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-5	Anion	03-03102-7- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
2	MW-12-4	VOC	03-03102-6- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-4	Metal	03-03102-6- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-4	Anion	03-03102-6- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
3	MW-12-3	VOC	03-03102-5- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-3	Metal	03-03102-5- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-3	Anion	03-03102-5- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
4	MW-12-2	VOC	03-03102-4- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-2	Metal	03-03102-4- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-2	Anion	03-03102-4- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
5	MW-12-1	VOC	03-03102-3- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-1	Metal	03-03102-3- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	MW-12-1	Anion	03-03102-3- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
6	TB-11-5/7/03	VOC	03-03102-8	W	V	C	40	2	G	050703	N	0	7 <input type="checkbox"/>
7	EB-11-5/7/03	VOC	03-03102-2- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	EB-11-5/7/03	Metal	03-03102-2- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	EB-11-5/7/03	Anion	03-03102-2- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>
8	DUPE-6-2Q03	VOC	03-03102-1- $\alpha$	W	V	C	40	3	G	050703	N	0	7 <input type="checkbox"/>
	DUPE-6-2Q03	Metal	03-03102-1- $\beta$	W	P	N	500	1	G	050703	N	0	7 <input type="checkbox"/>
	DUPE-6-2Q03	Anion	03-03102-1- $\gamma$	W	P		1000	1	G	050703	N	0	7 <input type="checkbox"/>

## Part 3: Analysis Information

- Test Items:
- 524.2 Volatile Organic Compounds
  - 7196A Chromium (VI)
  - 314.0/300.0 Perchlorate, low level
  - 300.0 Chloride  $\text{Cl}^-$  by IC
  - 300.0 Sulfate ( $\text{SO}_4^{--}$ ), by IC
  - 300.0/SM4500NO<sub>3</sub> Nitrate ( $\text{NO}_3^-$ ) as N by IC
  - SM2320B Carbonate
  - SM2320B Bicarbonate
  - 9040B/150.1 pH
  - 160.1 Solids, Total Dissolved (TDS)
  - 200.7/6010B Sodium, Na, by ICP
  - 200.7/6010B Calcium, Ca, by ICP

Level C Data Package Deliverables

# Volatile Organics



Applied P & Ch Laboratory

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/16/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: 03G2404-MB-01	Lab Sample ID: 03G2404-MB-01	Received Date: 05/16/2003
Sample Type: Method Blank	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: G2404K01	Prep. No: -	Anal. Time: 14:07
Methanol Vol. -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	<0.5	U
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	<0.5	U
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROETHENE (TOTAL)	540-59-0	µg/L	0.5	<0.5	U
32	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
33	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
34	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
35	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
36	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
37	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
38	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
39	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U



#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U
41	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
42	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	U
43	METHYLENE CHLORIDE	75-09-2	µg/L	1.8	6.7	
44	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
45	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
46	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
47	STYRENE	100-42-5	µg/L	0.5	<0.5	U
48	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
49	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
50	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
51	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
52	1,2,3-TRICHLOROENZENE	87-61-6	µg/L	0.5	<0.5	U
53	1,2,4-TRICHLOROENZENE	120-82-1	µg/L	0.5	<0.5	U
54	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
55	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
56	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
57	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
58	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
59	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
60	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
61	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
62	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
63	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
64	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	4-BROMO-FLUOROBENZENE (BFB)	460-00-4		70-129	104	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	103	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	105	
4	TOLUENE-D8	2037-26-5		73-129	108	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROENZENE-D5	3114-55-4		50-200	87	
2	1,4-DICHLOROENZENE-D4	3855-82-1		50-200	99	
3	FLUOROBENZENE	462-06-6		50-200	99	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

Qualifier: U - Not Detected or less than MDL

E - Exceed calibration range

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

B - A positive value was found in the method blank

D - Diluted

Data Filename: C:\HPCHEM\1\DATA\03G2404\G2404K01.D Sample : f=1  
 Method : C:\HPCHEM\1\METHODS\E524G004.M Inst. : GCMS-G  
 Acq. Time : May 16 14:07 2003 RF via : Multiple Level Calibration  
 Method Update: Fri May 16 12:11 2003 Operator: Eddie  
 Quant. Time : May 16 14:28 2003 Multiplr: 1.000000  
 Print Time : Fri May 16 14:38 2003  
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppb	C,ppb	Quality	Note
Internal Standards											
1	1 Fluorobenzene I1	9.51	9.44	0.007	96	70	646.001	10.00		0.07	
47	47 Cl-benzene-d5, I2	13.11	13.05	0.004	82	119	178.183	10.00		0.06	
62	62 1,4-DCB-d4 150 15	15.62	15.56	0.004	152	150	174.849	10.00		0.06	

System Monitoring Compounds (Surrogate)											
27	27 Di-Br-F-Methane (	7.96	7.90	0.004	111	113	502.090	20.93		20.9	104.66%
29	29 1,2-di-Cl-ethane-	8.53	8.47	0.004	65	102	222.912	20.45		20.4	102.23%
55	55 toluene-d8 (S2)	11.60	11.54	0.004	100	99	699.779	21.50		21.5	107.51%
70	70 4-Br-1-F-Bz (S3)	14.34	14.28	0.004	174	95	298.013	20.68		20.7	103.41%

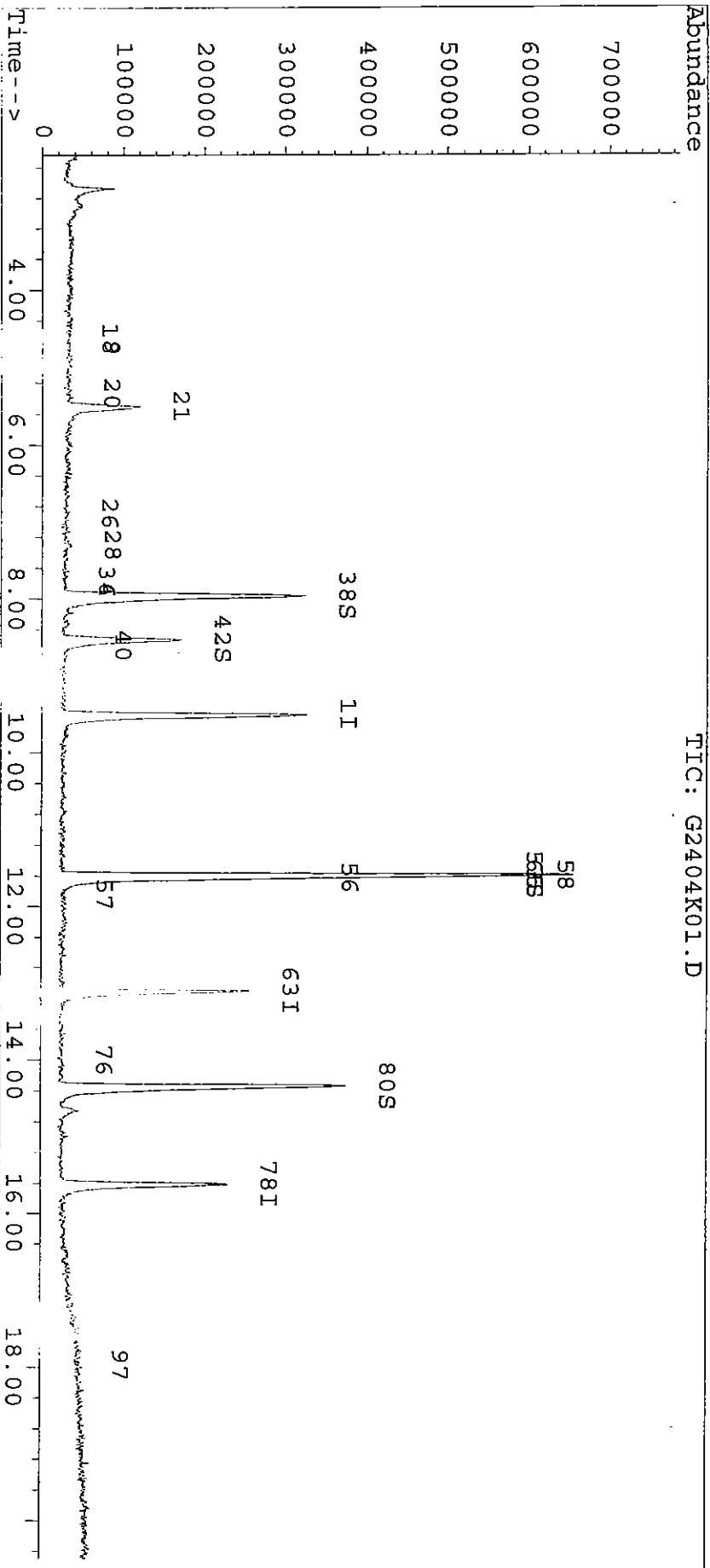
Target Compounds												
<<<	I1	: ISTD ID = 1	>>>									Qvalue
111	111	isopropyl alcoho	4.73	4.73	0.000	45	43	1.133	4.38	<del>4.4</del>	1	
101	101	Acetonitrilex10	4.64	4.65	-0.002	41	40	0.250	10.12	<del>10.1</del>	1	
113	113	Tert butyl alcoho	5.33	5.32	0.001	59	57	1.202	2.02	<del>2.0</del>	79	
18	18	methylene chlorid	5.50	5.45	0.006	84	49	106.423	6.72	<del>6.7</del>	95	
98	98	Vinyl acetate x5	6.89	6.90	-0.002	43	86	0.748	12.55	<del>12.5</del>	69	
91	91	2-butanone MEKx10	7.30	7.31	-0.002	43	72	7.793	1.29	<del>1.3</del>	71	
201	201	Ethyl acetate x2	7.79	7.78	0.002	43	61	3.879	4.77	<del>4.8</del>	52	
117	117	Iso-butyl alcoho	7.79	7.78	0.002	43	42	3.879	15.97	<del>16.0</del>	51	
30	30	12-dichloroethane	8.62	8.57	0.006	62	64	0.272	0.90	<del>0.9</del>	1	
107	107	Et methacrylate	11.64	11.73	-0.010	69	99	0.513	1.65	<del>1.6</del>	58	
93	93	2-Hexanone x5	11.84	11.86	-0.002	43	58	0.659	5.79	<del>5.8</del>	87	
48	48	112-tri-Cl-Et	11.62	11.42	0.021	97	83	22.264	2.93	<del>2.9</del>	8	
<<< I2 : ISTD ID = 47 >>>												
49	49	1,3-di-Cl-propane	11.60	11.68	-0.006	76	78	8.608	0.68	<del>0.7</del>	99	
110	110	t-1,4-dichloro-2	14.00	14.04	-0.003	89	53	0.673	0.48	<del>0.5</del>	9	
<<< I3 : ISTD ID = 62 >>>												
88	88	naphthalene	17.97	17.92	0.004	128	129	0.282	1.37	<del>1.4</del>	69	

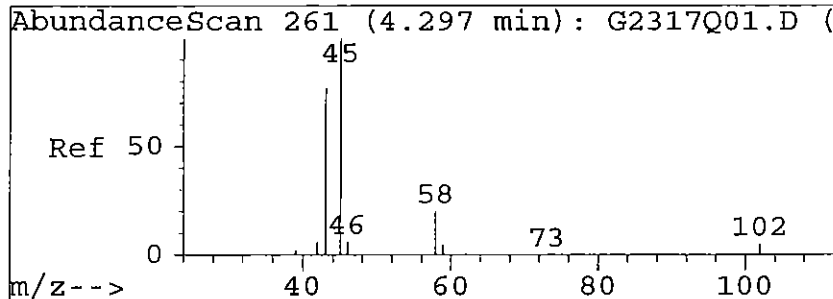
# = qualifier out of range, m = manual integration, ? = RT coelution, \* = DRRT > 0.06

Data File : C:\HPCHEM\1\DATA\03G2404\G2404K01.D  
Acq On : 16 May 03 2:07 pm  
Sample : f=1  
Misc :  
Quant Time: May 16 14:28 2003

Vial: 10  
Operator: Eddie  
Inst : GCMS-G  
Multiplier: 1.00  
Quant Results File: quant.res

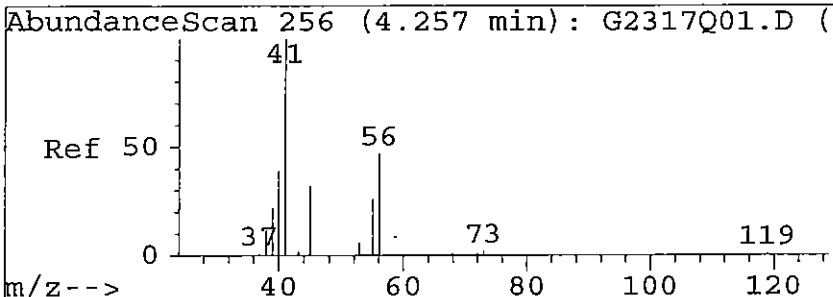
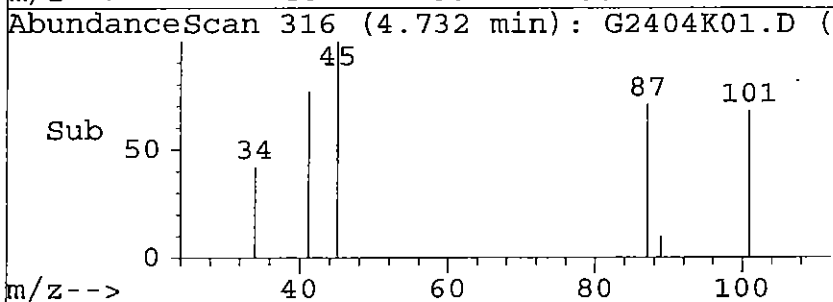
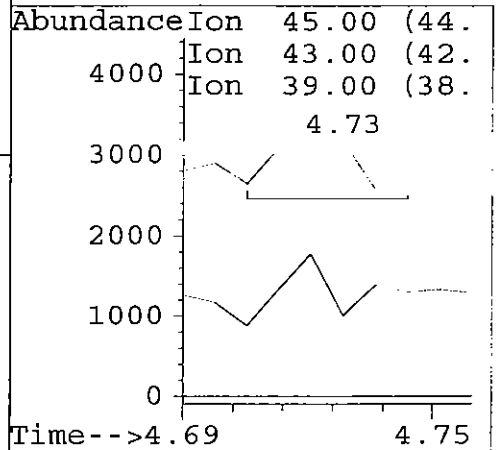
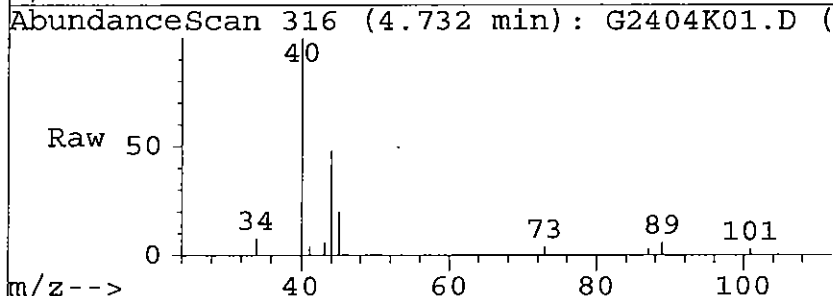
Method : C:\HPCHEM\1\METHODS\E524G004.M  
Title : \*\*Applied P & Ch Lab\*\* EPA 524.2  
Last Update : Fri May 16 12:11:07 2003  
Response via : Multiple Level Calibration





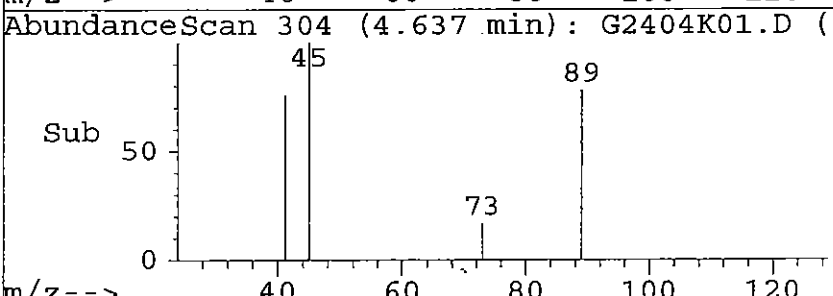
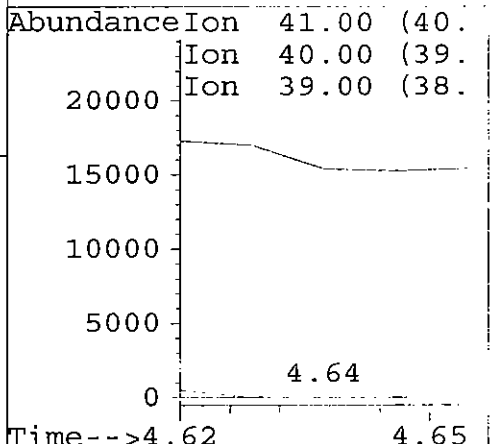
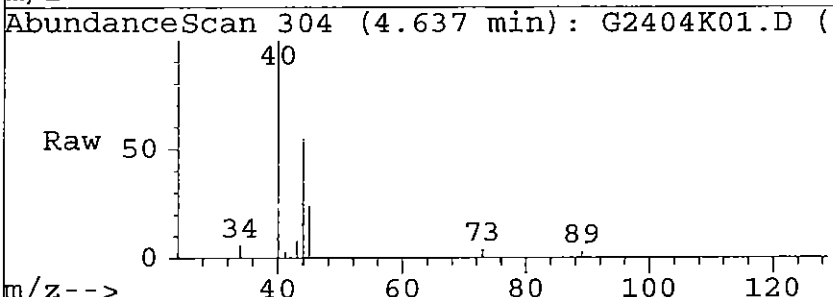
#9  
 111 isopropyl alcohol x10  
 Concen: 4.38 ppb  
 RT: 4.73 min Scan# 316  
 Delta R.T. 0.01 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

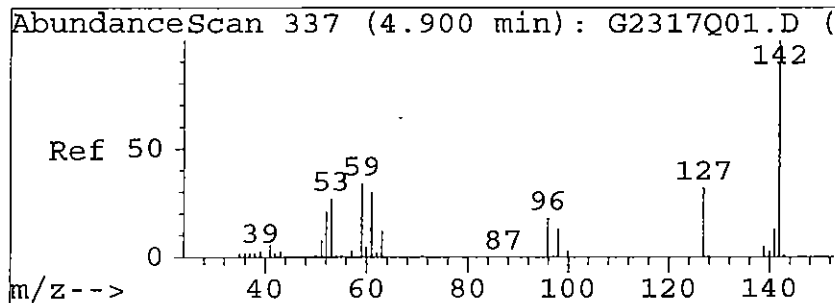
Tgt Ion	Resp	Lower	Upper
45	1133		
43	326.0	77.0	115.5#
39	0.0	13.3	19.9#
0	0.0	0.0	0.0



#18  
 101 Acetonitrile x10  
 Concen: 10.12 ppb  
 RT: 4.64 min Scan# 304  
 Delta R.T. -0.02 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

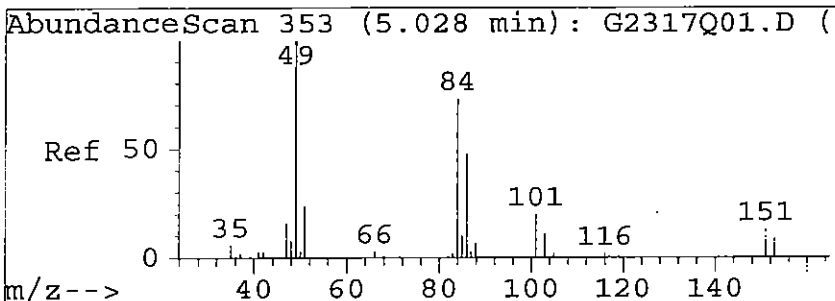
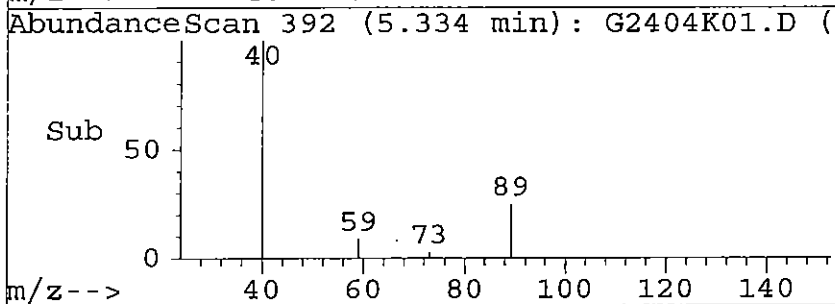
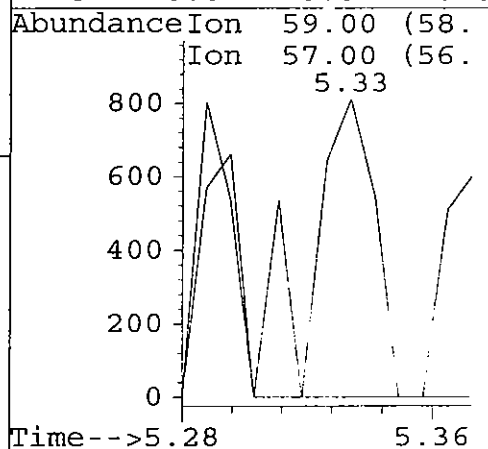
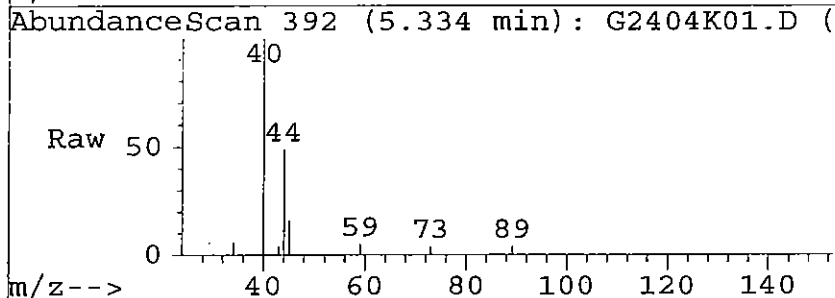
Tgt Ion	Resp	Lower	Upper
41	250		
40	100	151.7	227.5#
39	0.0	17.3	25.9#
0	0.0	0.0	0.0





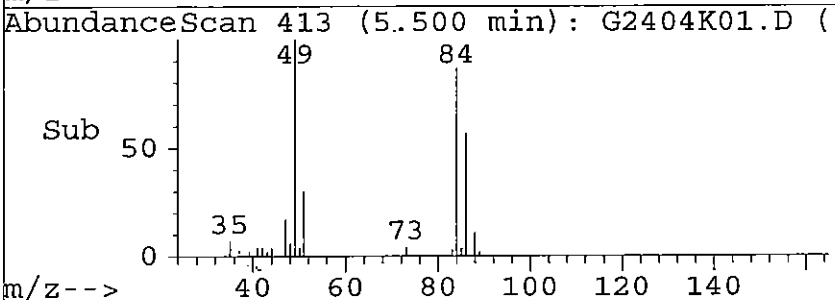
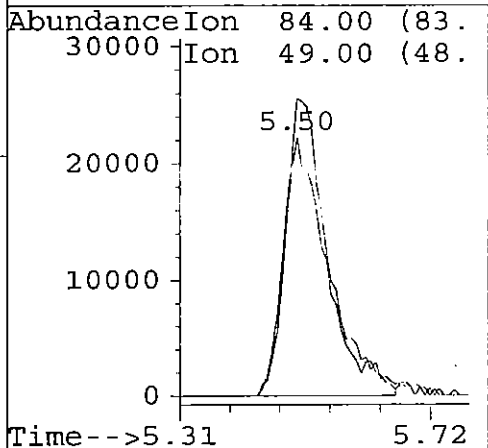
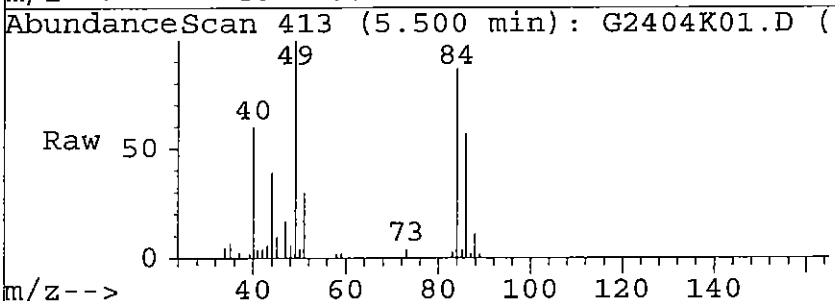
#20  
 113 Tert butyl alcohol x10  
 Concen: 2.02 ppb  
 RT: 5.33 min Scan# 392  
 Delta R.T. 0.01 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

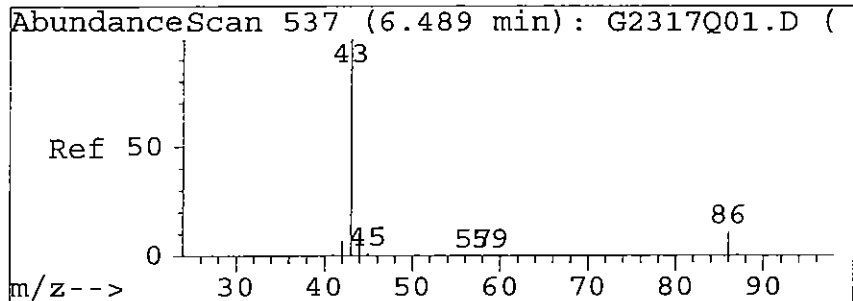
Tgt Ion	Resp	Lower	Upper
59	1202		
57	0.0	5.8	8.7#
0	0.0	0.0	0.0
0	0.0	0.0	0.0



#21  
 18 methylene chloride 49 84  
 Concen: 6.72 ppb  
 RT: 5.50 min Scan# 413  
 Delta R.T. 0.05 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

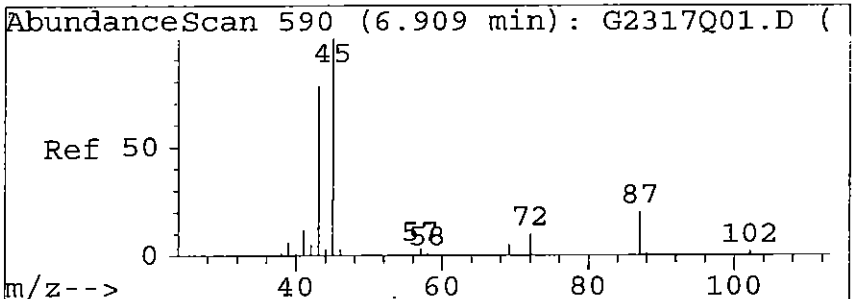
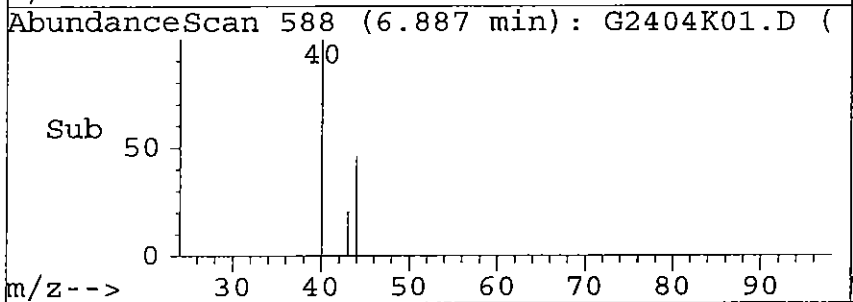
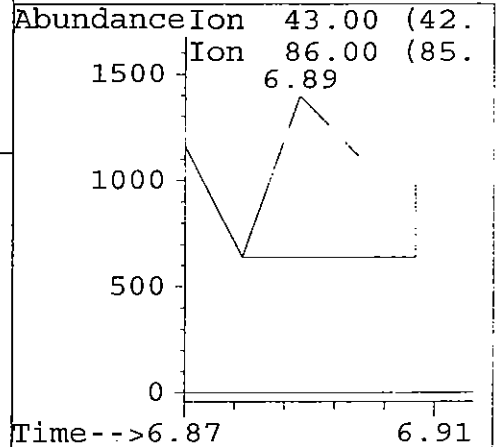
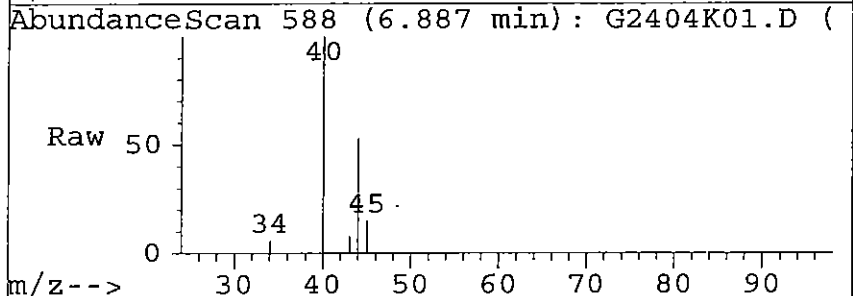
Tgt Ion	Resp	Lower	Upper
84	106423		
49	120.4	62.9	188.5
0	0.0	0.0	0.0
0	0.0	0.0	0.0





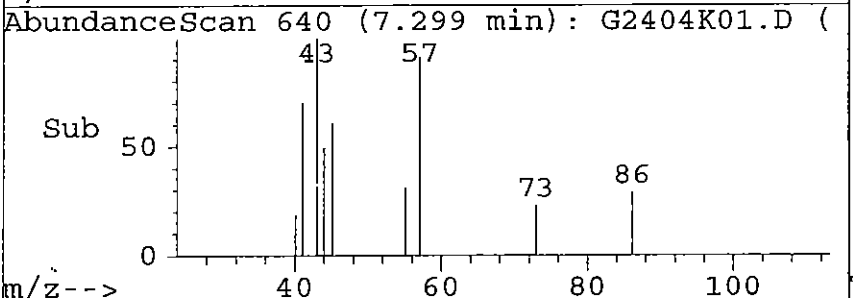
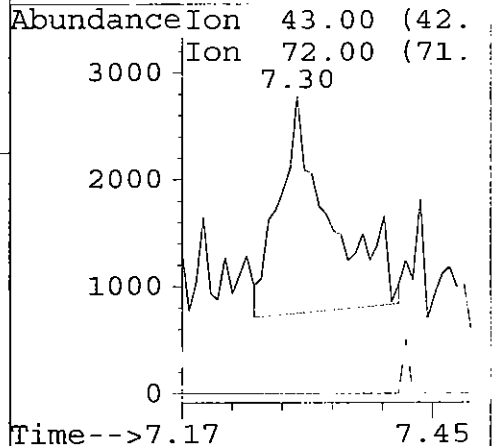
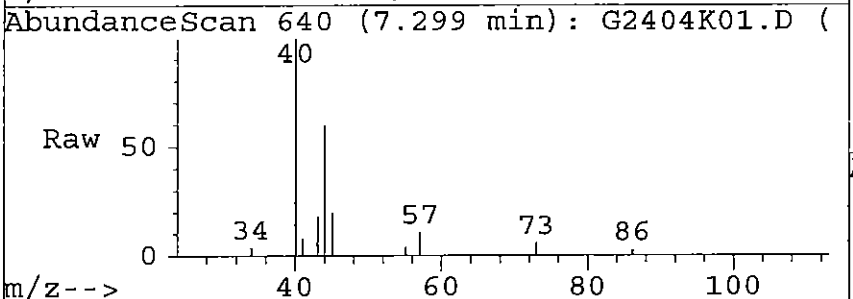
#26  
 98 Vinyl acetate x5  
 Concen: 12.55 ppb  
 RT: 6.89 min Scan# 588  
 Delta R.T. -0.02 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

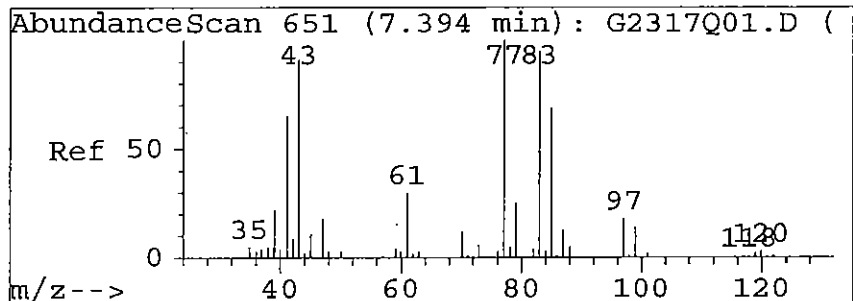
Tgt Ion	43	86	0	0
Resp:	748	0.0	0.0	0.0
Lower	9.5	0.0	0.0	0.0
Upper	14.2#	0.0	0.0	0.0



#28  
 91 2-butanone MEKx10  
 Concen: 1.29 ppb  
 RT: 7.30 min Scan# 640  
 Delta R.T. -0.02 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

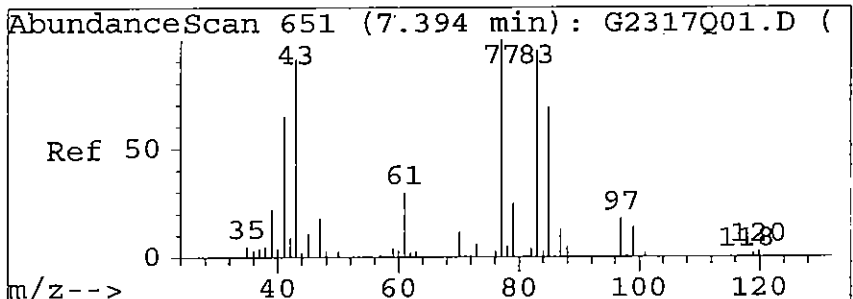
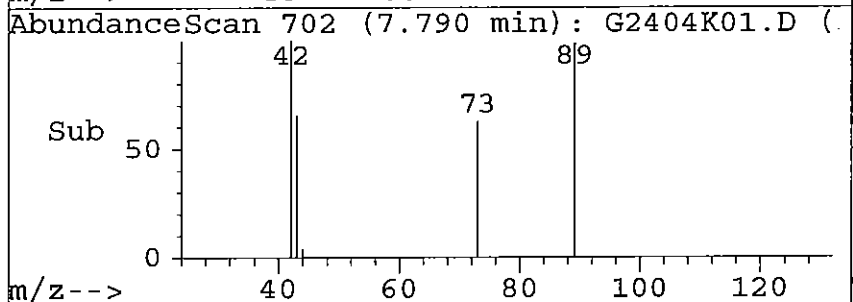
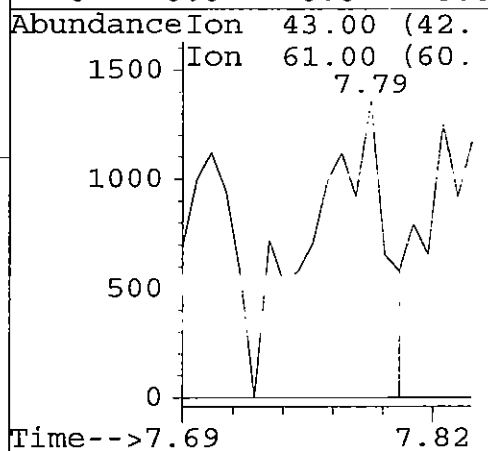
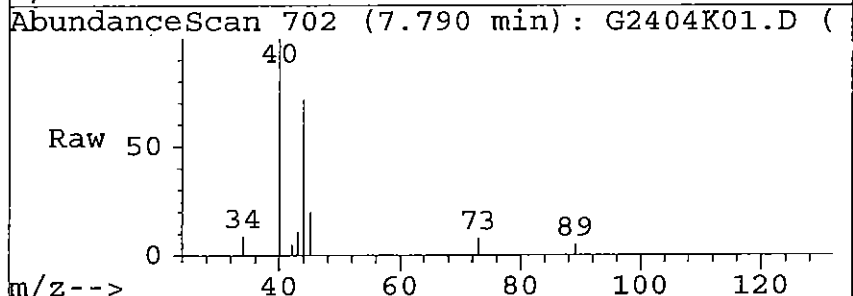
Tgt Ion	43	72	0	0
Resp:	7793	0.0	0.0	0.0
Lower	9.0	0.0	0.0	0.0
Upper	13.5#	0.0	0.0	0.0





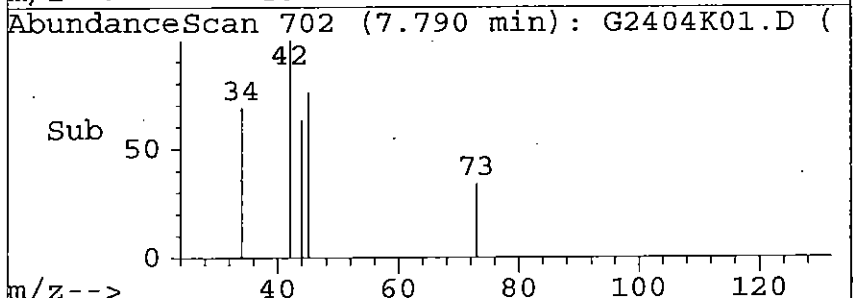
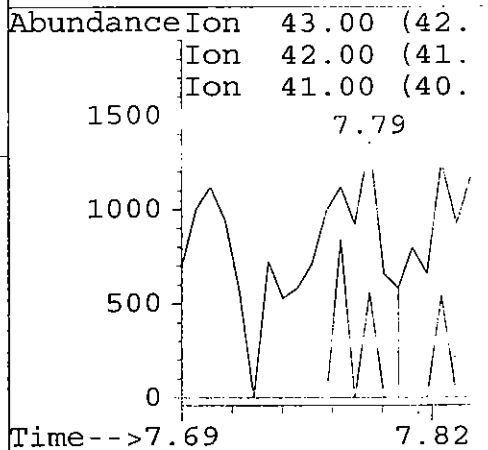
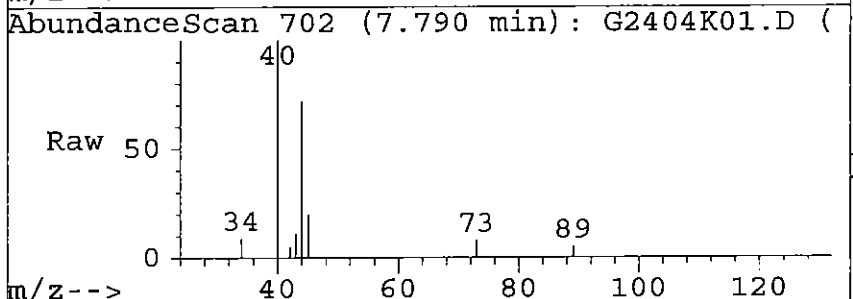
#34  
 201 Ethyl acetate x2  
 Concen: 4.77 ppb  
 RT: 7.79 min Scan# 702  
 Delta R.T. 0.01 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

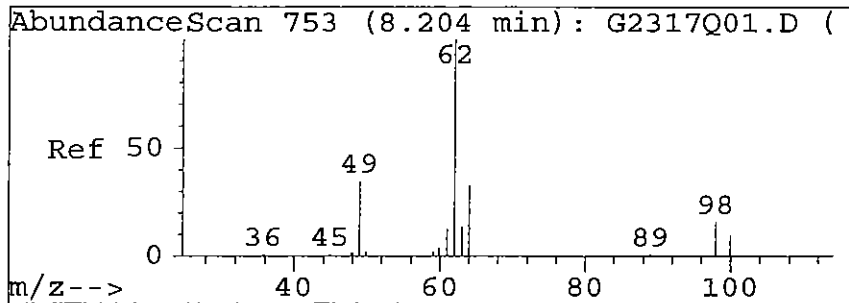
Tgt Ion:43	Resp:	3879
Ion Ratio	Lower	Upper
43	100	
61	0.0	19.1 28.6#
0	0.0	0.0 0.0
0	0.0	0.0 0.0



#36  
 117 Iso-butyl alcohol X10  
 Concen: 15.97 ppb  
 RT: 7.79 min Scan# 702  
 Delta R.T. 0.01 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

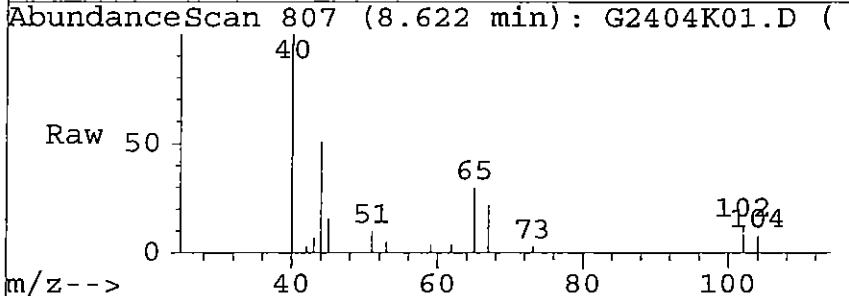
Tgt Ion:43	Resp:	3879
Ion Ratio	Lower	Upper
43	100	
42	6.8	7.3 10.9#
41	10.3	39.7 59.5#
0	0.0	0.0 0.0



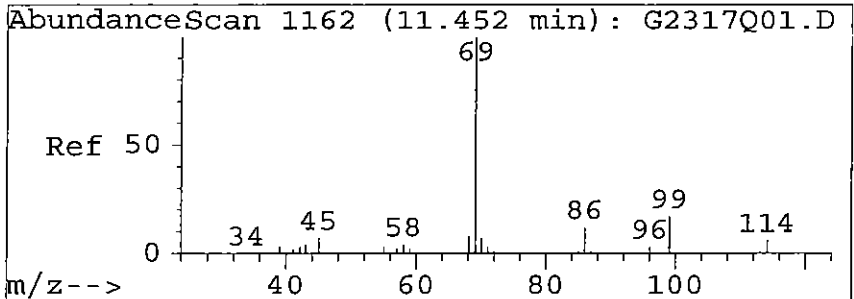
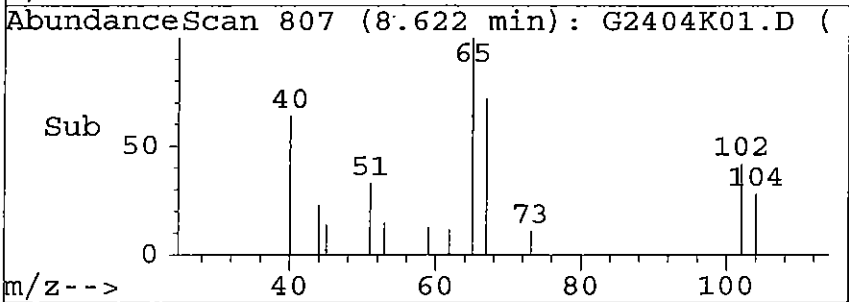
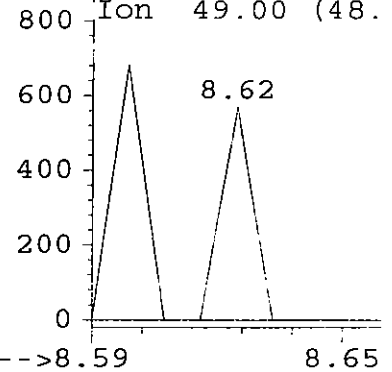


#40  
 30 12-dichloroethane 64 62  
 Concen: 0.90 ppb  
 RT: 8.62 min Scan# 807  
 Delta R.T. 0.05 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

Tgt Ion	62	Resp	272
Ion Ratio	Lower	Upper	
62	100		
64	0.0	13.1	53.1#
49	2080.1	16.2	56.2#
0	0.0	0.0	0.0

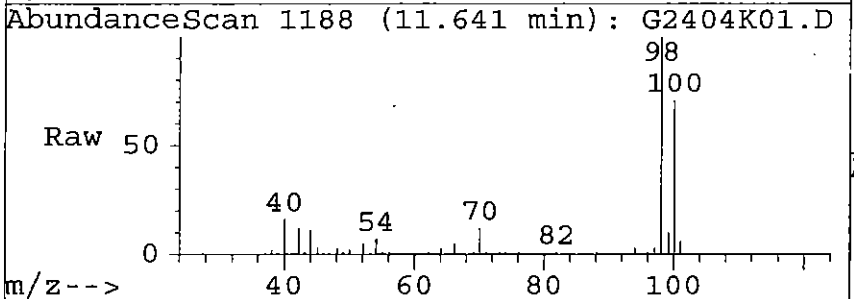


Abundance	Ion	62.00 (61.
	Ion	64.00 (63.
	Ion	49.00 (48.

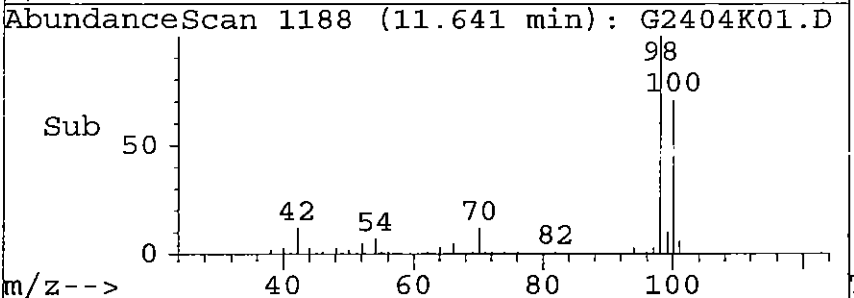
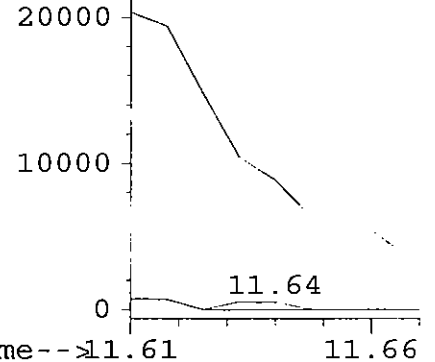


#56  
 107 Et methacrylate  
 Concen: 1.65 ppb  
 RT: 11.64 min Scan# 1188  
 Delta R.T. -0.09 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

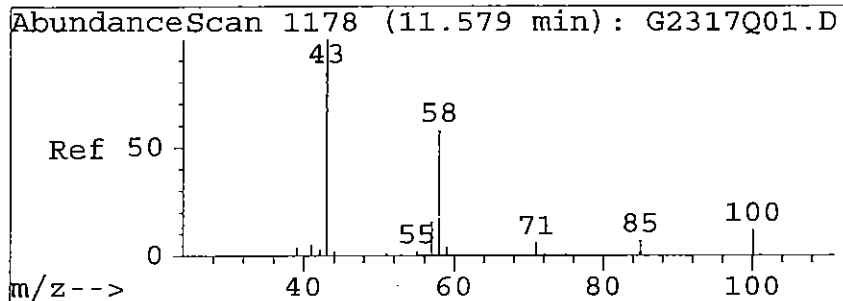
Tgt Ion	69	Resp	513
Ion Ratio	Lower	Upper	
69	100		
99	0.0	9.4	28.0#
0	0.0	0.0	0.0
0	0.0	0.0	0.0



Abundance	Ion	69.00 (68.
	Ion	99.00 (98.

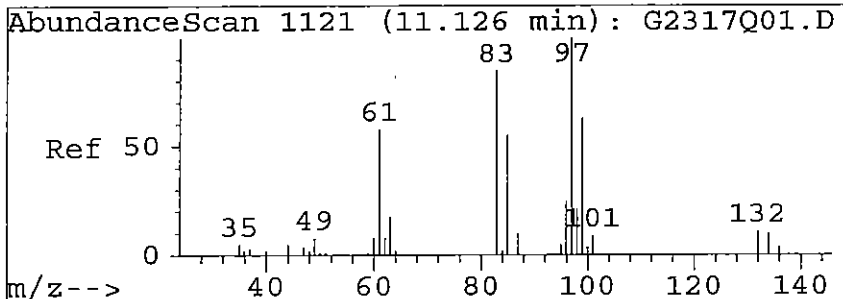
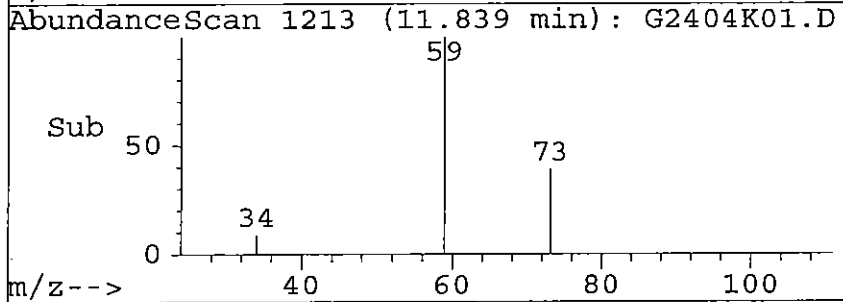
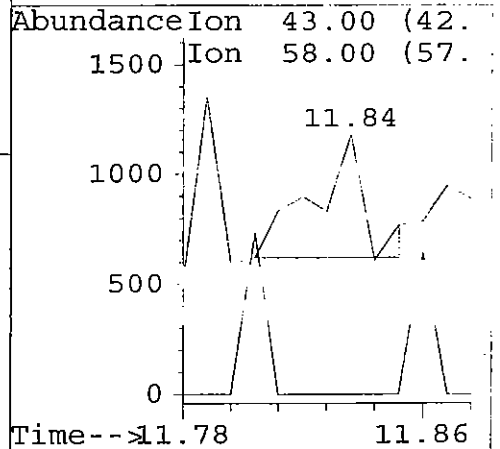
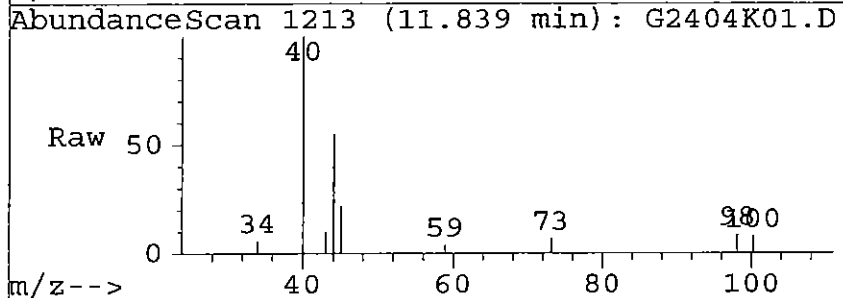






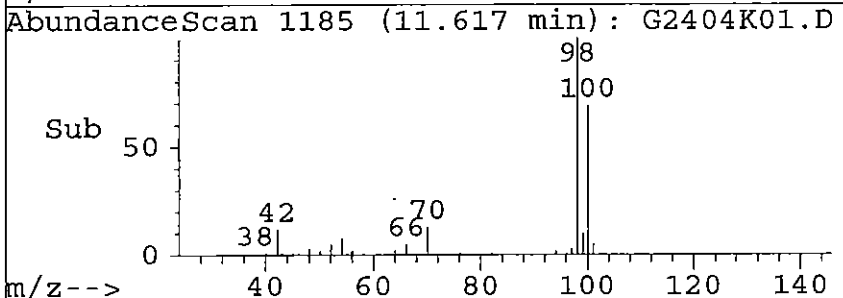
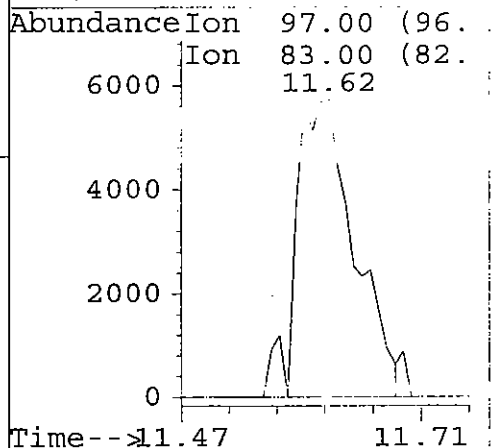
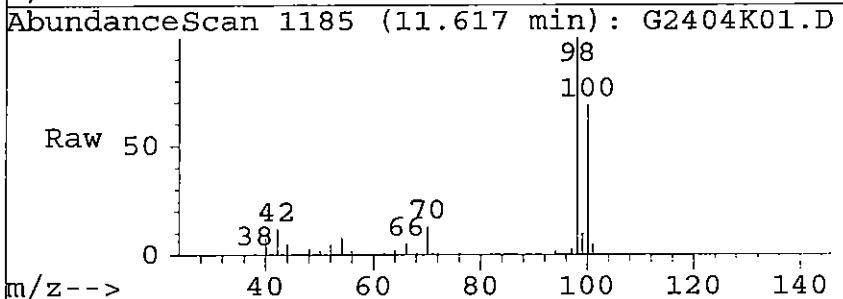
#57  
 93 2-Hexanone x5  
 Concen: 5.79 ppb  
 RT: 11.84 min Scan# 1213  
 Delta R.T. -0.02 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

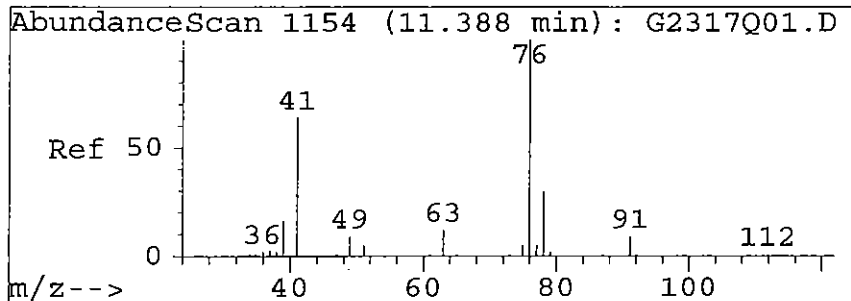
Tgt Ion	43	Resp	659
Ion Ratio	100	Lower	Upper
58	46.3	44.5	66.8
0	0.0	0.0	0.0
0	0.0	0.0	0.0



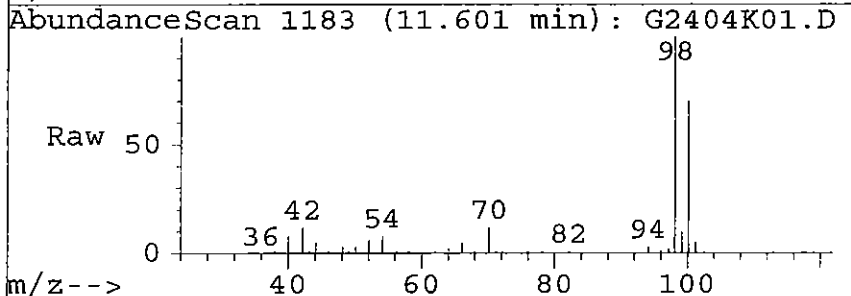
#58  
 48 112-tri-Cl-Et 97 83  
 Concen: 2.93 ppb  
 RT: 11.62 min Scan# 1185  
 Delta R.T. 0.20 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm

Tgt Ion	97	Resp	22264
Ion Ratio	100	Lower	Upper
83	0.0	41.5	124.4#
0	0.0	0.0	0.0
0	0.0	0.0	0.0

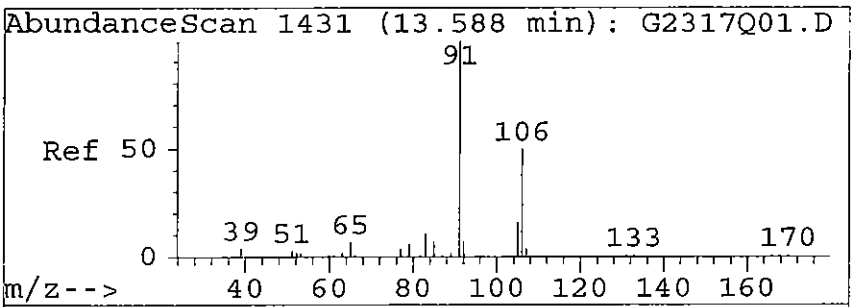
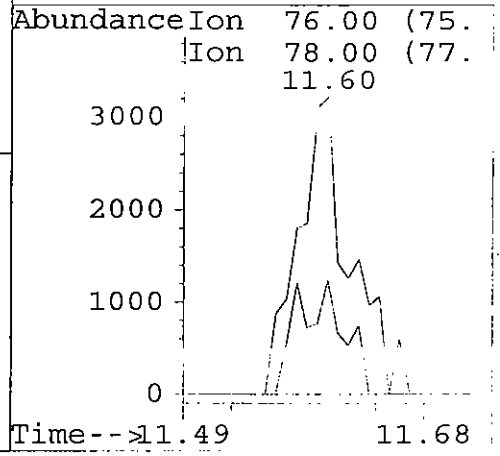
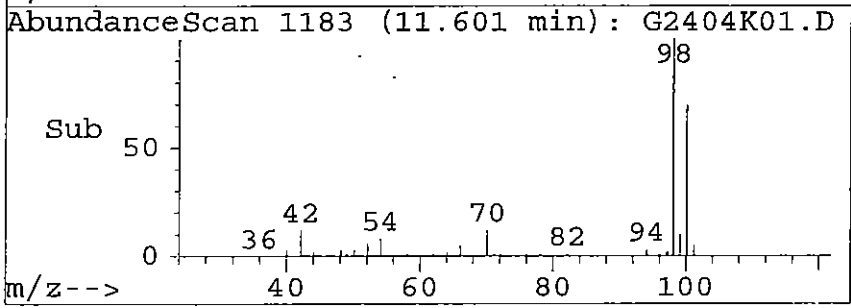




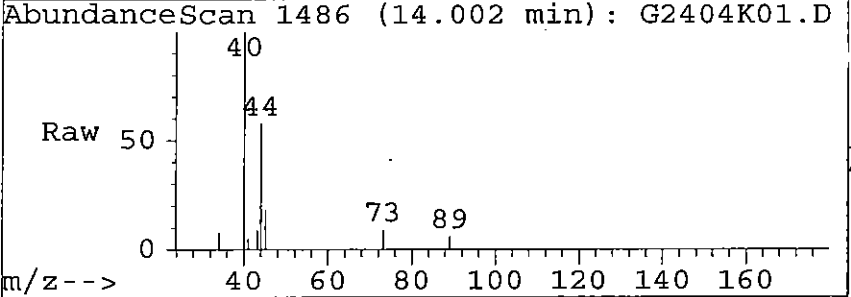
#65  
 49 1,3-di-cl-propane 76 78  
 Concen: 0.68 ppb  
 RT: 11.60 min Scan# 1183  
 Delta R.T. -0.07 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm



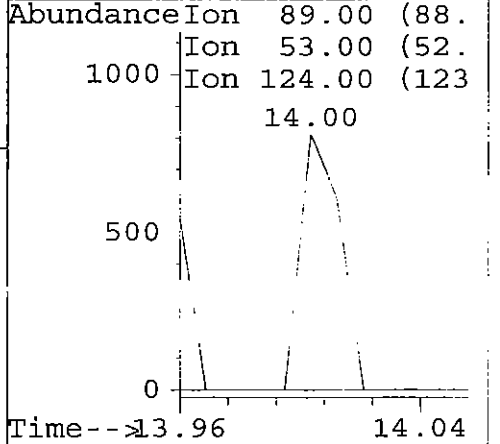
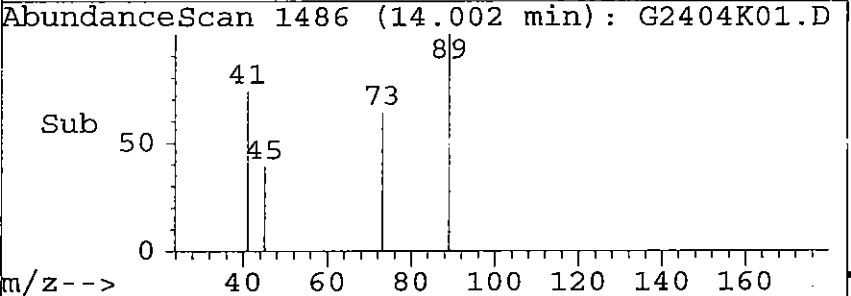
Tgt Ion	Resp	Lower	Upper
76	100		
78	35.5	27.9	41.8
0	0.0	0.0	0.0
0	0.0	0.0	0.0



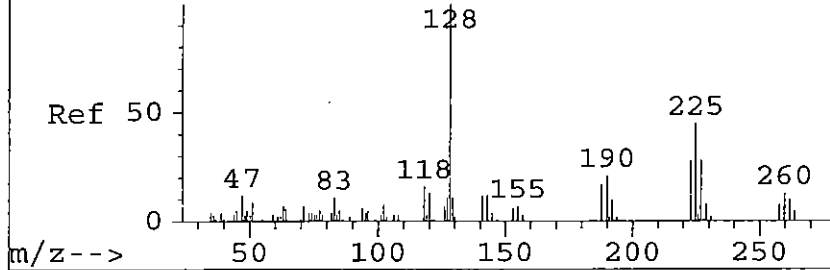
#76  
 110 t-1,4-dichloro-2-butene  
 Concen: 0.48 ppb  
 RT: 14.00 min Scan# 1486  
 Delta R.T. -0.04 min  
 Lab File: G2404K01.D  
 Acq: 16 May 03 2:07 pm



Tgt Ion	Resp	Lower	Upper
89	100		
53	0.0	83.4	125.1#
124	0.0	29.4	44.1#
0	0.0	0.0	0.0



AbundanceScan 1933 (17.578 min): G2317Q01.D



#97

88 naphthalene 128 129

Concen: 1.37 ppb

RT: 17.97 min Scan# 1987

Delta R.T. 0.06 min

Lab File: G2404K01.D

Acq: 16 May 03 2:07 pm

Tgt Ion:128 Resp: 282

Ion Ratio Lower Upper

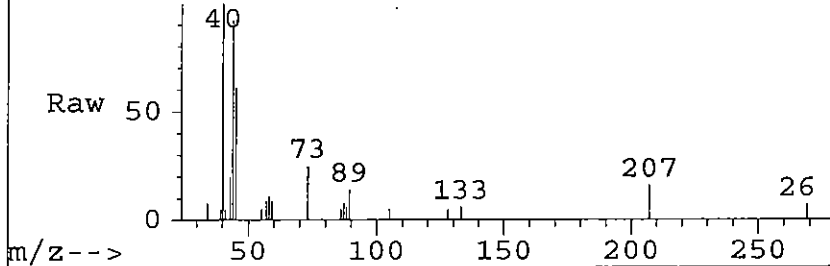
128 100

129 0.0 6.2 18.4#

0 0.0 0.0 0.0

0 0.0 0.0 0.0

AbundanceScan 1987 (17.974 min): G2404K01.D



Abundance Ion 128.00 (127

Ion 129.00 (128

17.97

600

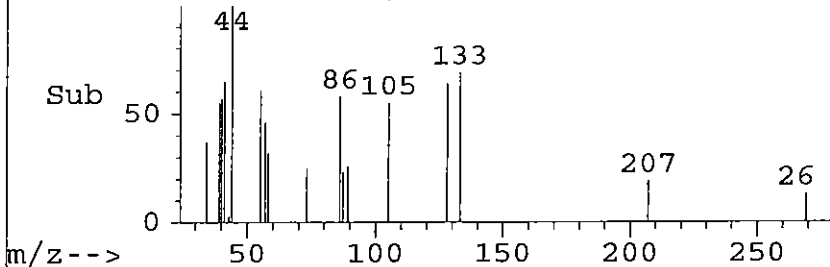
400

200

0

Time-->17.94 18.00

AbundanceScan 1987 (17.974 min): G2404K01.D



Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: <b>DUPE-6-2Q03</b>	Lab Sample ID: 03-3102-1	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-01	Prep. No: -	Anal. Time: 17:59
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	2.6	
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	1.2	
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROETHENE (TOTAL)	540-59-0	µg/L	0.5	<0.5	U
32	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
33	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
34	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
35	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
36	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
37	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
38	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
39	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U
41	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
42	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	4	J
43	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	<1.8	U
44	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
45	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
46	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
47	STYRENE	100-42-5	µg/L	0.5	<0.5	U
48	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
49	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
50	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
51	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
52	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
53	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
54	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
55	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
56	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
57	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
58	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
59	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
60	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
61	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
62	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
63	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
64	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U

## Surrogates

		Control Limit, %	Surro. Rec.%	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4	70-129	94
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	88
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	94
4	TOLUENE-D8	2037-26-5	73-129	96
# of out-of-control			0	

## Internal Standard

		Control Limit, %	IS Rec.%	
1	CHLOROBENZENE-D5	3114-55-4	50-200	107
2	1,4-DICHLOROBENZENE-D4	3855-82-1	50-200	118
3	FLUOROBENZENE	462-06-6	50-200	123
# of out-of-control			0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a)MDL reported.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted

Data Filename: C:\HPCHEM\1\DATA\03G2404\3102-01.D Sample : F=1 8  
 Method : C:\HPCHEM\1\METHODS\E524G004.M Inst. : GCMS-G  
 Acq. Time : May 16 17:59 2003 RF via : Multiple Level Calibration  
 Method Update: Fri May 16 12:11 2003 Operator: Eddie  
 Quant. Time : May 19 10:20 2003 Multiplr: 1.000000  
 Print Time : Mon May 19 10:20 2003  
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppb	C,ppb	Quality	Note
Internal Standards											
1	Fluorobenzene I1	9.47	9.52	-0.005	96	70	797.559	10.00		-0.05	
47	Cl-benzene-d5, I2	13.08	13.12	-0.003	82	119	217.915	10.00		-0.04	
62	1,4-DCB-d4 150 15	15.60	15.63	-0.002	152	150	209.628	10.00		-0.03	

System Monitoring Compounds (Surrogate)											
27	Di-Br-F-Methane (	7.94	7.97	-0.002	111	113	557.153	18.81	18.8	94.07%	
29	1,2-di-Cl-ethane-	8.50	8.54	-0.003	65	102	235.511	17.50	17.5	87.49%	
55	toluene-d8 (S2)	11.57	11.61	-0.003	100	99	774.823	19.28	19.3	96.42%	
70	4-Br-1-F-Bz (S3)	14.32	14.36	-0.002	174	95	324.004	18.76	18.8	93.78%	

Qvalue

Target Compounds											
<<< I1 : ISTD ID = 1 >>>											
111	111 isopropyl alcoho	4.76	4.76	0.000	45	43	0.839	2.63	<del>2.6</del>	<del>10</del>	#
102	102 Acrolein x10	4.80	4.67	0.013	56	55	0.238	4.44	<del>4.4</del>	<del>14</del>	#
101	101 Acetonitrilex10	4.72	4.71	0.000	41	40	1.364	12.11	<del>12.1</del>	<del>1</del>	#
113	113 Tert butyl alcoho	5.38	5.38	0.000	59	57	0.307	0.42	<del>0.4</del>	<del>79</del>	#
18	18 methylene chlorid	5.48	5.50	-0.003	84	49	25.220	11.29	1.3	94	#
98	98 Vinyl acetate x5	6.99	6.97	0.002	43	86	0.936	12.55	12.5	69	#
91	91 2-butanone MEKx10	7.37	7.37	0.000	43	72	6.272	0.84	0.8	71	#
25	25 chloroform	7.77	7.81	-0.005	83	85	48.408	1.18	1.2	98	#
201	201 Ethyl acetate x2	7.82	7.83	0.000	43	61	1.343	4.44	<del>4.4</del>	<del>52</del>	#?
117	117 Iso-butyl alcoho	7.82	7.83	0.000	43	42	1.343	14.30	<del>14.3</del>	<del>35</del>	#?
30	30 12-dichloroethane	8.61	8.63	-0.003	62	64	0.601	0.92	<del>0.9</del>	<del>1</del>	#
37	37 CC14	9.16	9.19	-0.003	117	119	73.832	2.58	2.6	96	#
118	118 TAME	9.47	9.47	0.000	73	43	13.982	0.40	<del>0.4</del>	<del>70</del>	#
107	107 Et methacrylate	11.60	11.80	-0.020	69	99	0.799	1.66	1.7	58	#?
93	93 2-Hexanone x5	11.96	11.93	0.003	43	58	0.435	5.71	<del>5.7</del>	<del>95</del>	#?
48	48 112-tri-Cl-Et	11.59	11.49	0.010	97	83	25.317	2.72	<del>2.7</del>	<del>9</del>	#?
<<< I2 : ISTD ID = 47 >>>											
54	54 MIBK	10.94	10.99	-0.004	43	58	27.296	3.88	3.9	76	#
49	49 1,3-di-cl-propane	11.56	11.75	-0.015	76	78	9.220	0.60	<del>0.6</del>	<del>88</del>	? #

# = qualifier out of range, m = manual integration, ? = RT coelution, \* = DRRT > 0.06

Quantitation Report: \*\*Applied P' & Ch Lab\*\* EPA 524.2

Data Filename: C:\HPCHEM\1\DATA\03G2404\3102-01.D Sample : f=1 8  
 Method : C:\HPCHEM\1\METHODS\E524G004.M Inst. : GCMS-G  
 Acq. Time : May 16 17:59 2003 RF via : Multiple Level Calibration  
 Method Update: Fri May 16 12:11 2003 Operator: Eddie  
 Quant. Time : May 19 10:20 2003 Multiplr: 1.000000  
 Print Time : Mon May 19 10:20 2003  
 Miscellaneous :

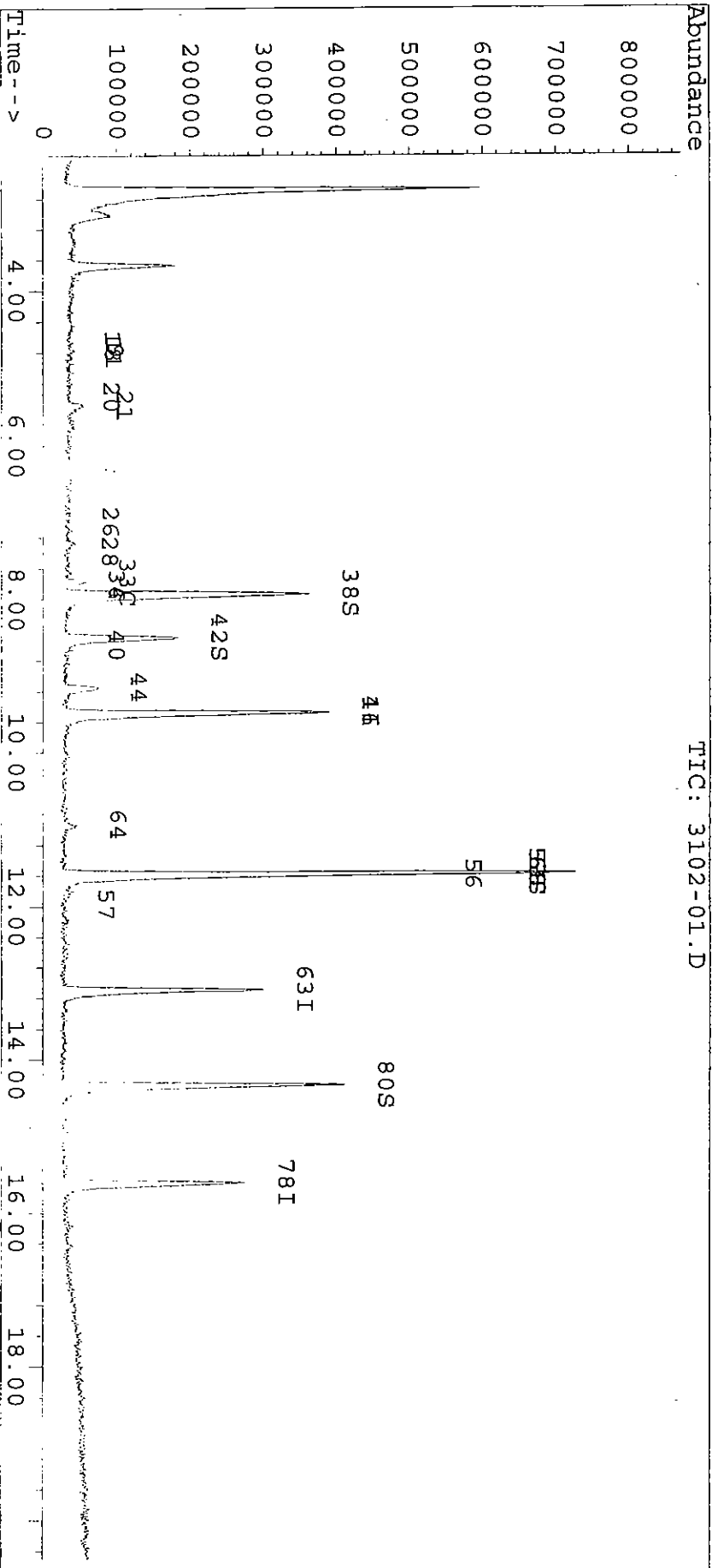
ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppb	c,ppb	Quality	Note
----	----------------	------	-----	------	------	----	---------	--------	-------	---------	------

# = qualifier out of range, m = manual integration, ? = RT coelution, \* = DRRT > 0.06

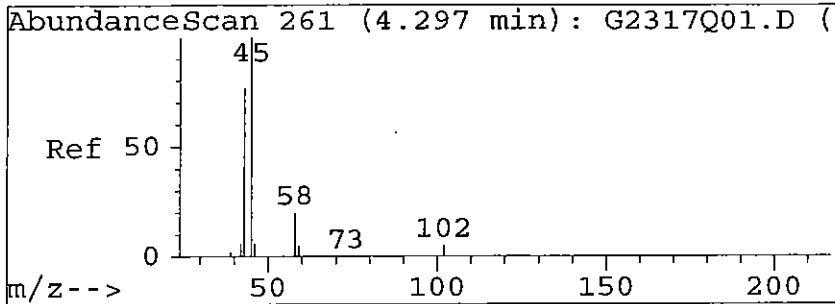
Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2404\3102-01.D  
Acq On : 16 May 03 5:59 pm  
Sample : f=1 8  
Misc :  
Quant Time: May 19 10:20 2003  
Operator: Eddie  
Inst : GCMS-G  
Multiplier: 1.00  
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\E524G004.M  
Title : \*\*Applied P & Ch Lab\*\* EPA 524.2  
Last Update : Fri May 16 12:11:07 2003  
Response via : Multiple Level Calibration

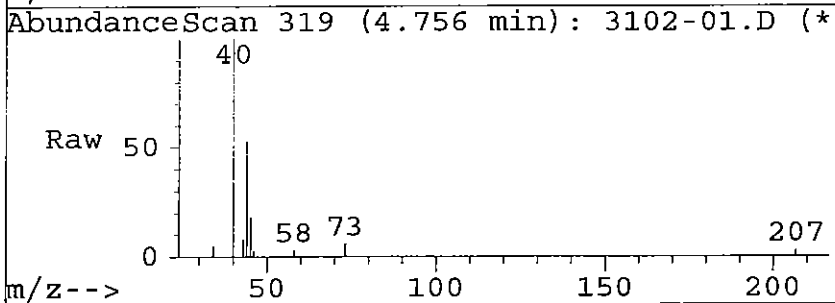




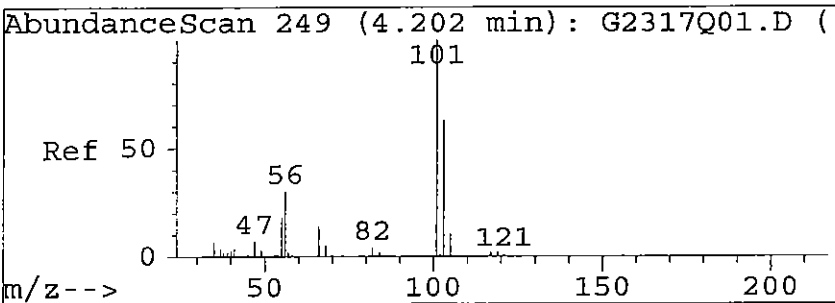
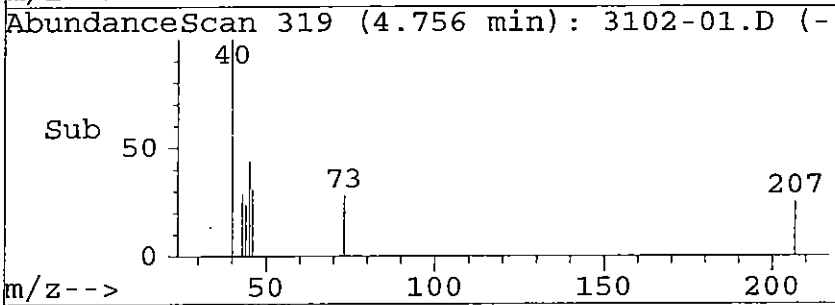
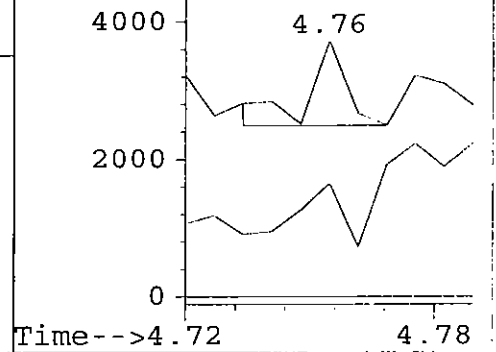


#9  
 111 isopropyl alcohol x10  
 Concen: 2.63 ppb  
 RT: 4.76 min Scan# 319  
 Delta R.T. 0.00 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
45	100		
43	0.0	77.0	115.5#
39	0.0	13.3	19.9#
0	0.0	0.0	0.0

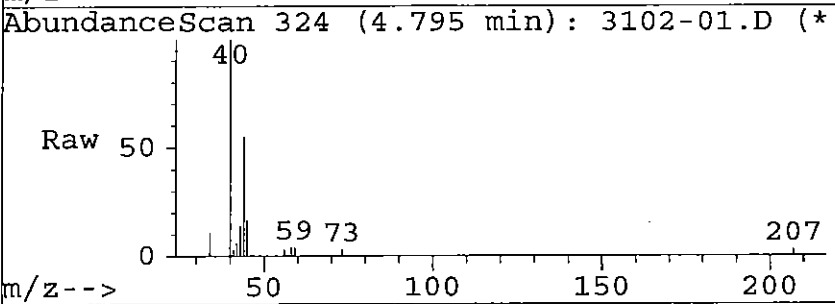


Abundance	Ion	Ion	Ion
45.00	(44.		
43.00	(42.		
39.00	(38.		

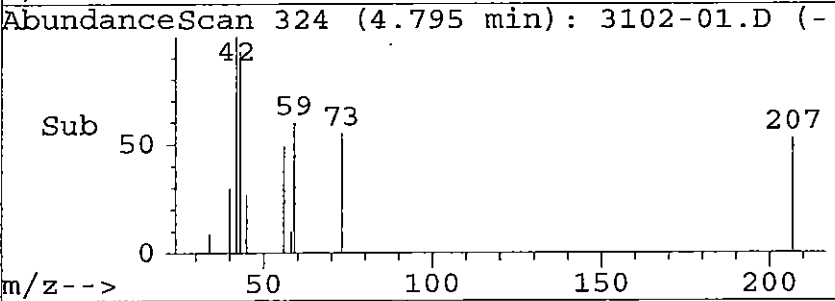
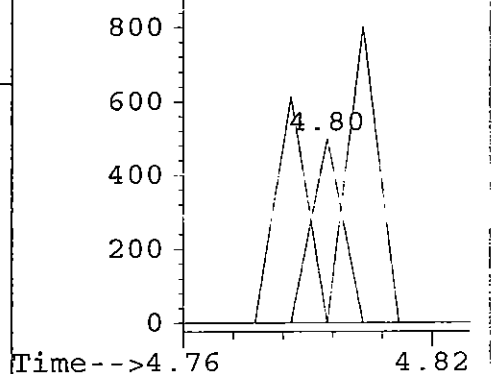


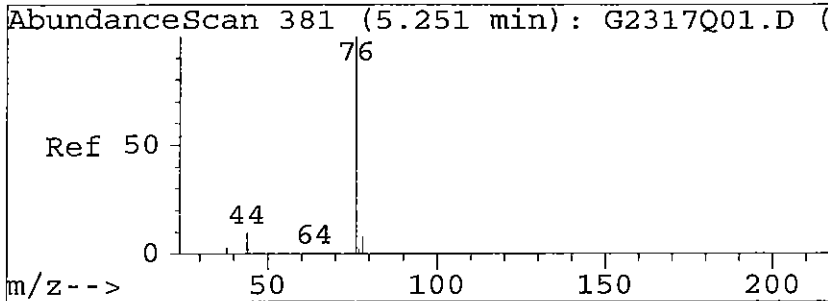
#11  
 102 Acrolein x10  
 Concen: 4.44 ppb  
 RT: 4.80 min Scan# 324  
 Delta R.T. 0.13 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
56	100		
55	122.7	29.2	87.5#
0	0.0	0.0	0.0
0	0.0	0.0	0.0



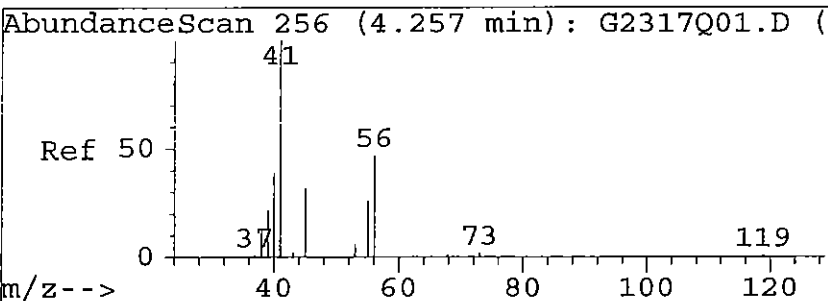
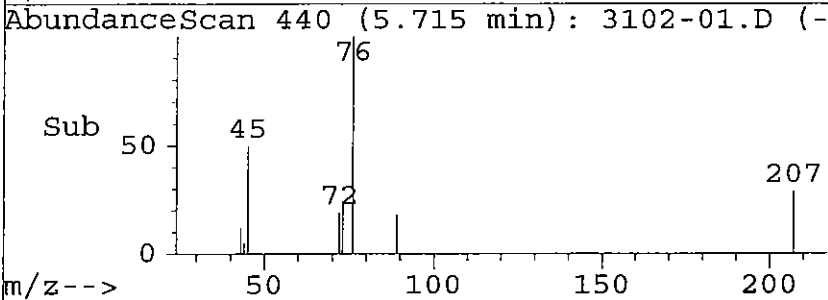
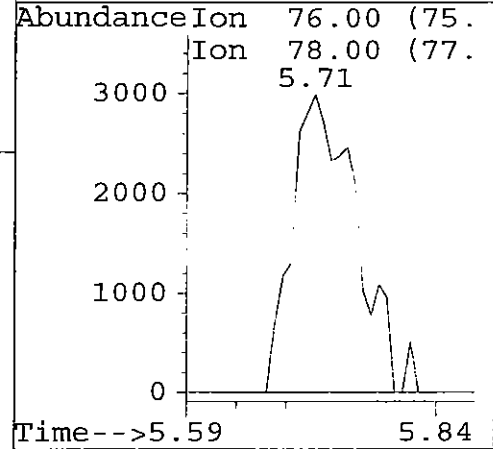
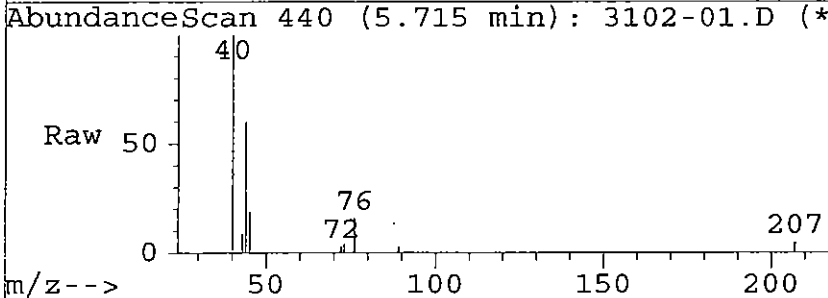
Abundance	Ion	Ion
56.00	(55.	
55.00	(54.	





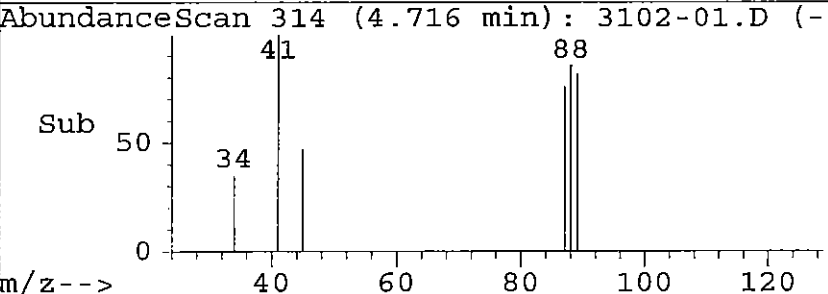
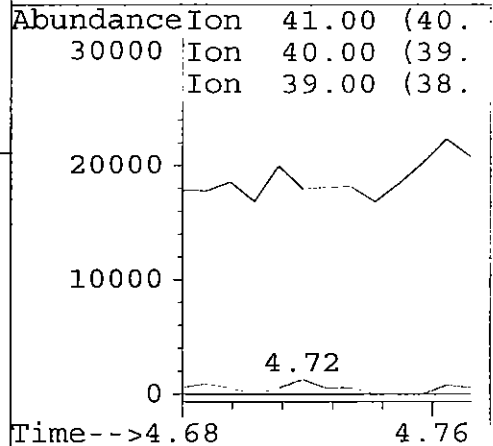
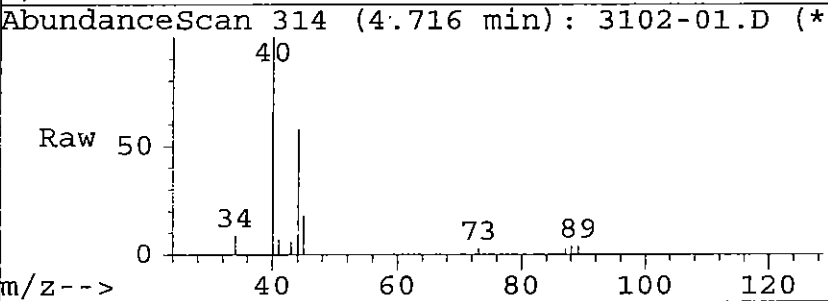
#13  
 104 Carbon disulfide  
 Concen: 0.26 ppb  
 RT: 5.71 min Scan# 440  
 Delta R.T. -0.02 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

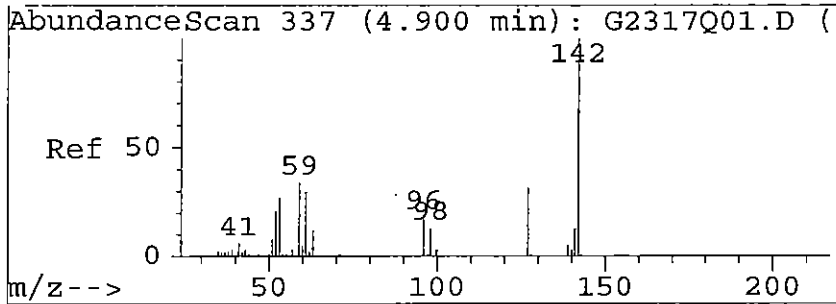
Tgt Ion	Resp	Lower	Upper
76	12991		
78	0.0	4.5	13.4#
0	0.0	0.0	0.0
0	0.0	0.0	0.0



#18  
 101 Acetonitrilex10  
 Concen: 12.11 ppb  
 RT: 4.72 min Scan# 314  
 Delta R.T. 0.01 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

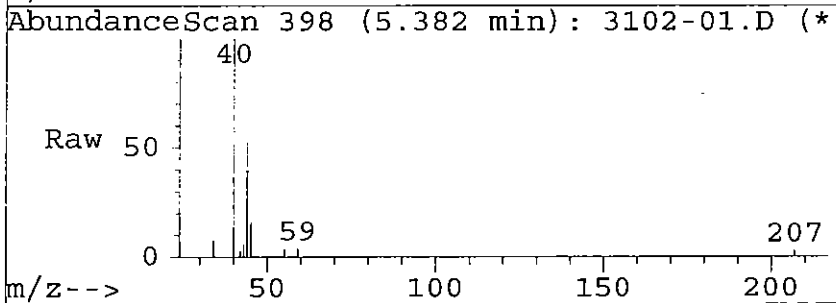
Tgt Ion	Resp	Lower	Upper
41	1364		
40	349.3	151.7	227.5#
39	0.0	17.3	25.9#
0	0.0	0.0	0.0



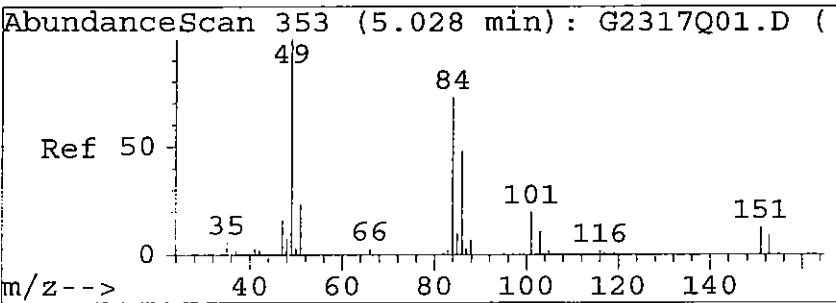
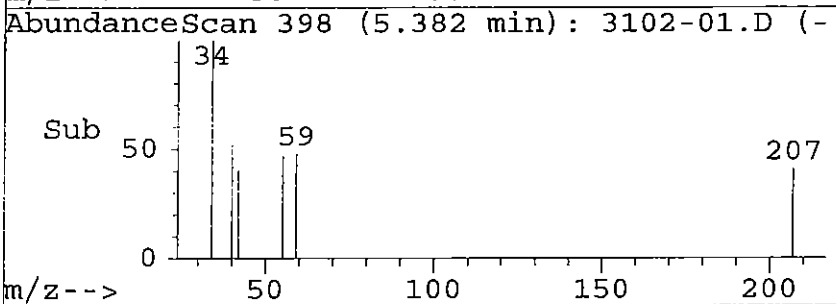
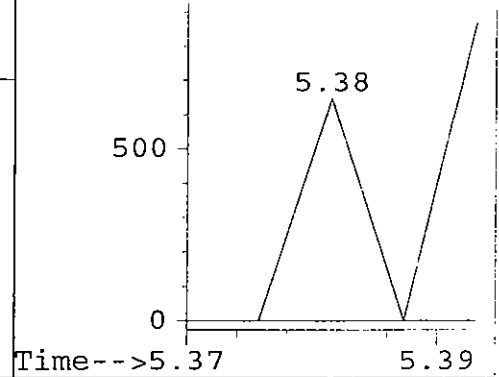


#20  
 113 Tert butyl alcohol x10  
 Concen: 0.42 ppb  
 RT: 5.38 min Scan# 398  
 Delta R.T. -0.00 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
59	100		
57	0.0	5.8	8.7#
0	0.0	0.0	0.0
0	0.0	0.0	0.0

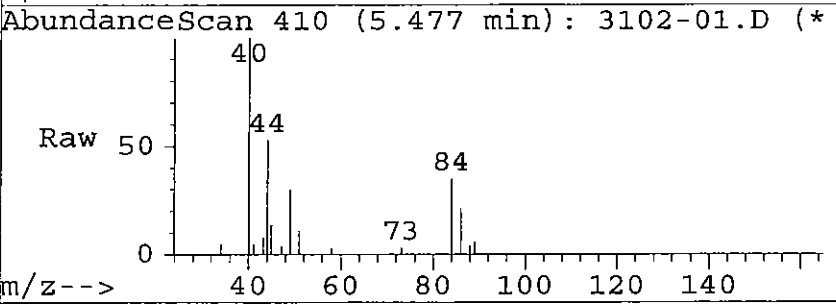


Abundance Ion 59.00 (58.  
 1000 Ion 57.00 (56.

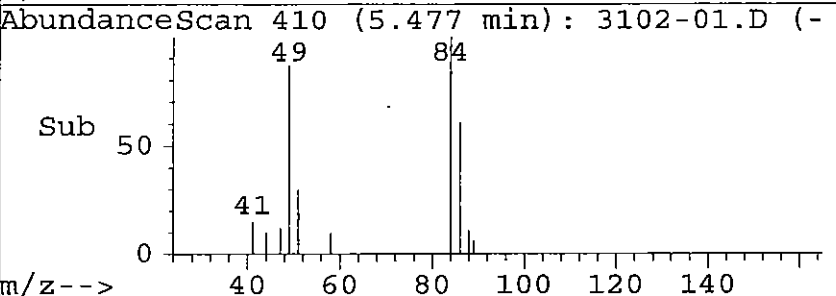
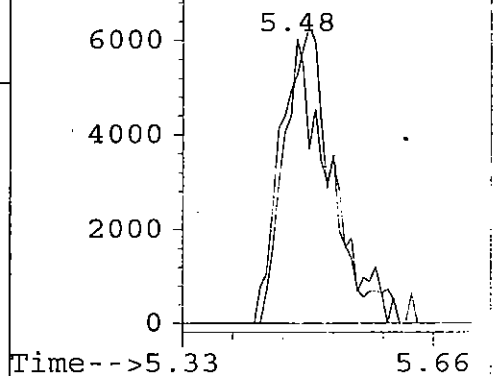


#21  
 18 methylene chloride 49 84  
 Concen: 1.29 ppb  
 RT: 5.48 min Scan# 410  
 Delta R.T. -0.03 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

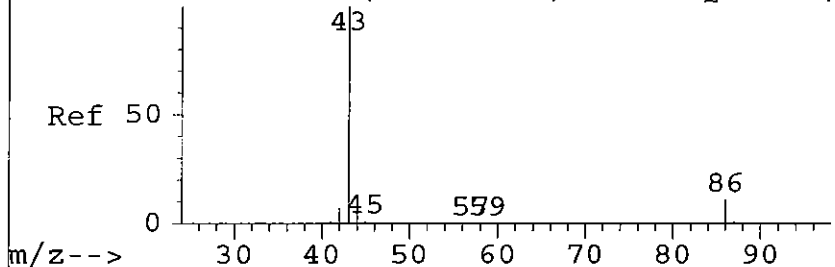
Tgt Ion	Resp	Lower	Upper
84	100		
49	118.9	62.9	188.5
0	0.0	0.0	0.0
0	0.0	0.0	0.0



Abundance Ion 84.00 (83.  
 Ion 49.00 (48.



AbundanceScan 537 (6.489 min): G2317Q01.D (

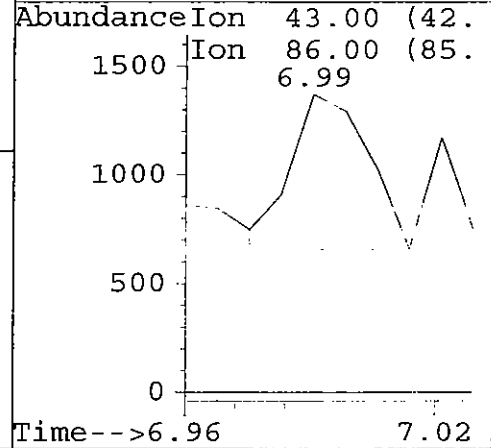
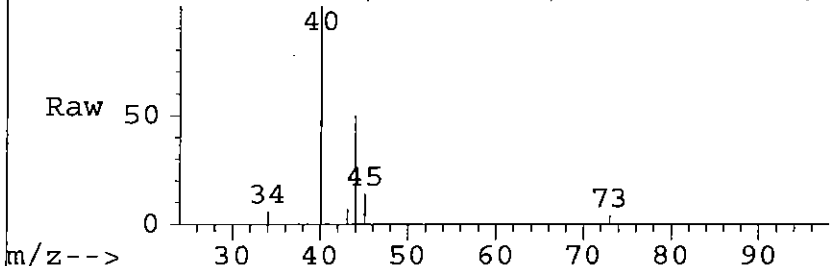


#26

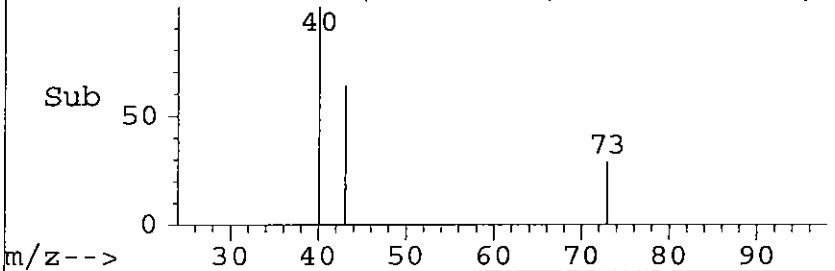
98 Vinyl acetate x5  
 Concen: 12.55 ppb  
 RT: 6.99 min Scan# 601  
 Delta R.T. 0.02 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	43	86	0	0	Resp:	936	Lower	Upper
Ion Ratio	100	0.0	0.0	0.0				
		9.5	0.0	0.0				14.2#

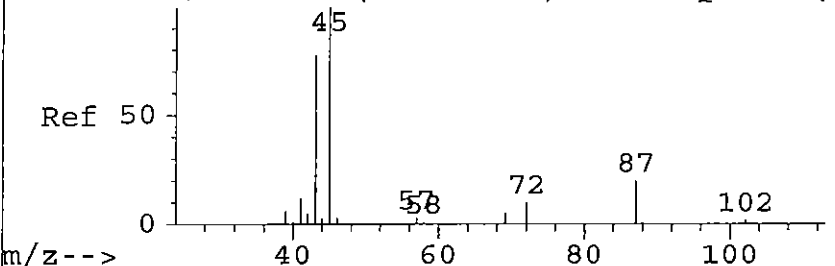
AbundanceScan 601 (6.990 min): 3102-01.D (\*)



AbundanceScan 601 (6.990 min): 3102-01.D (-



AbundanceScan 590 (6.909 min): G2317Q01.D (

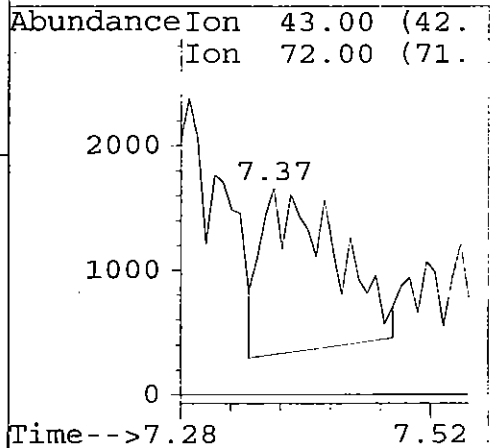
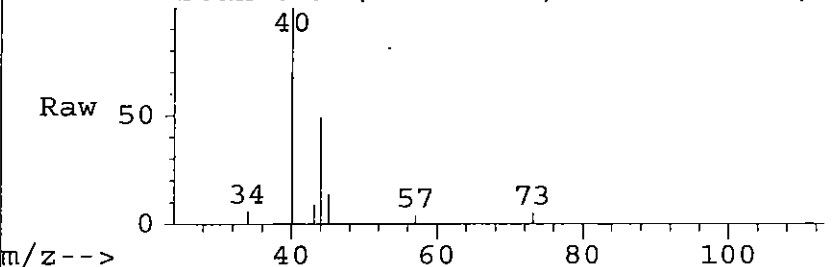


#28

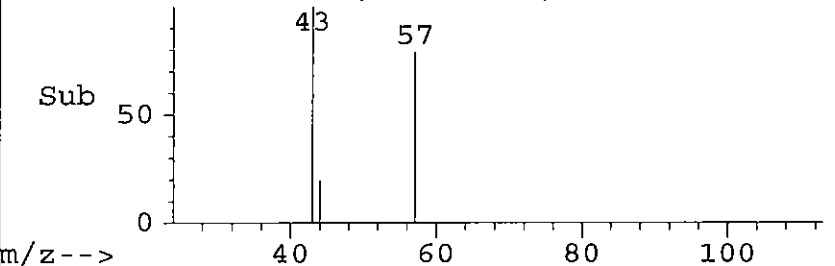
91 2-butanone MEKx10  
 Concen: 0.84 ppb  
 RT: 7.37 min Scan# 649  
 Delta R.T. -0.00 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

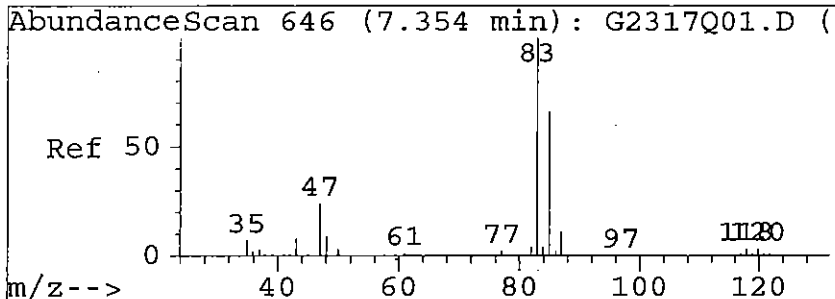
Tgt Ion	43	72	0	0	Resp:	6272	Lower	Upper
Ion Ratio	100	0.0	0.0	0.0				
		9.0	0.0	0.0				13.5#

AbundanceScan 649 (7.371 min): 3102-01.D (\*)

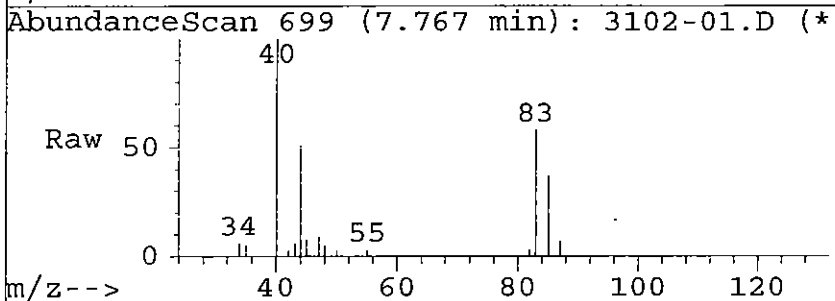


AbundanceScan 649 (7.371 min): 3102-01.D (-

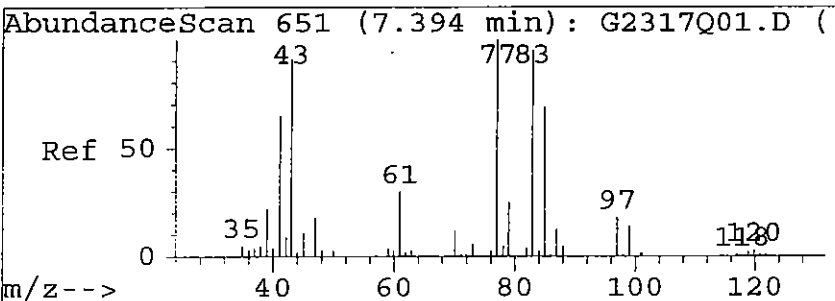
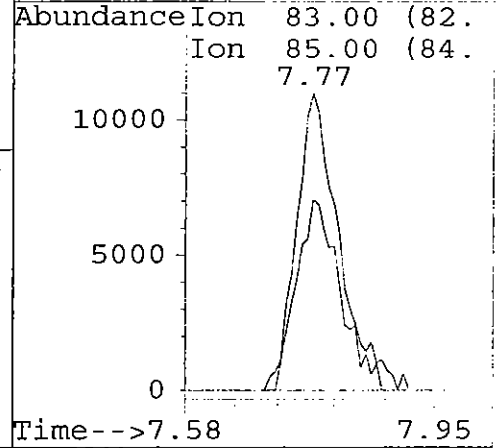
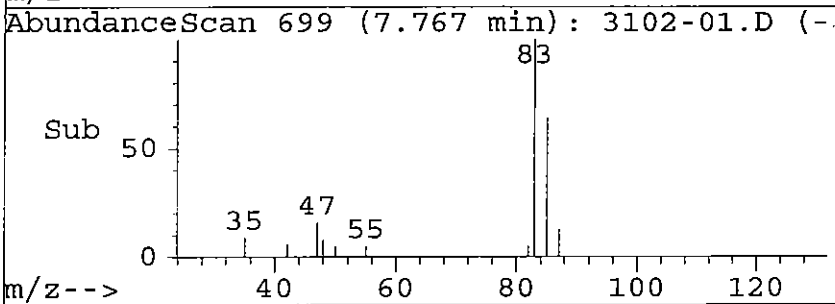




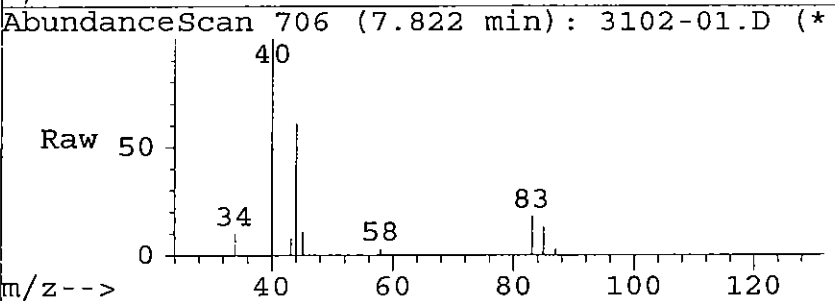
#33  
 25 chloroform 83 85  
 Concen: 1.18 ppb  
 RT: 7.77 min Scan# 699  
 Delta R.T. -0.05 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm



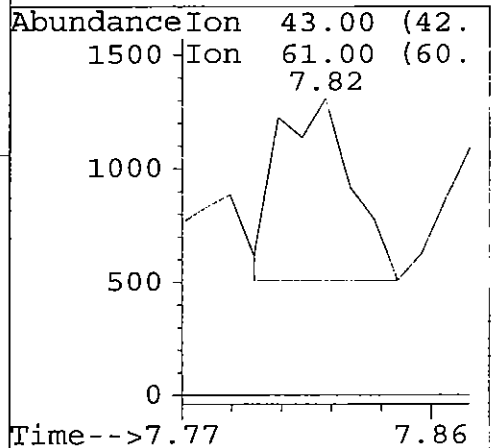
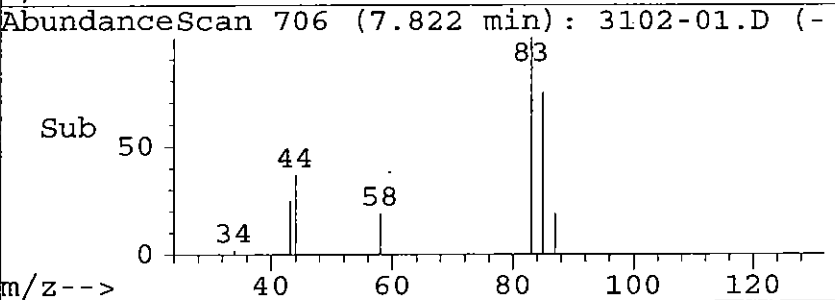
Tgt Ion	83	Resp	48408
Ion	Ratio	Lower	Upper
83	100		
85	65.6	33.5	100.3
0	0.0	0.0	0.0
0	0.0	0.0	0.0

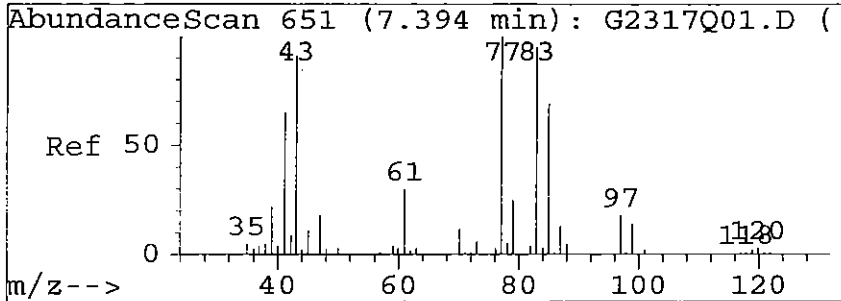


#34  
 201 Ethyl acetate x2  
 Concen: 4.44 ppb  
 RT: 7.82 min Scan# 706  
 Delta R.T. -0.01 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

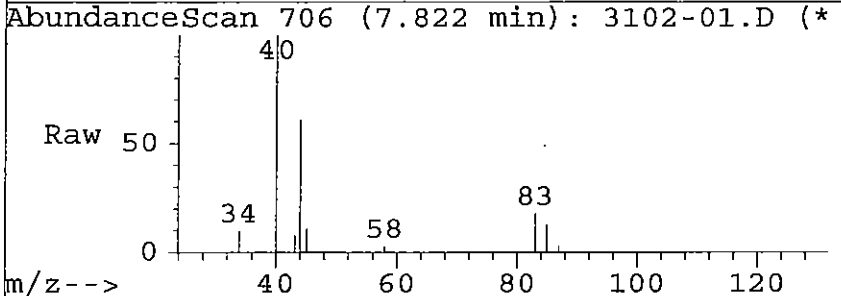


Tgt Ion	43	Resp	1343
Ion	Ratio	Lower	Upper
43	100		
61	0.0	19.1	28.6#
0	0.0	0.0	0.0
0	0.0	0.0	0.0

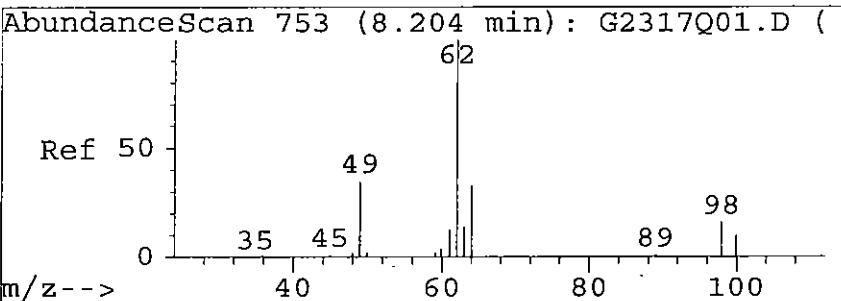
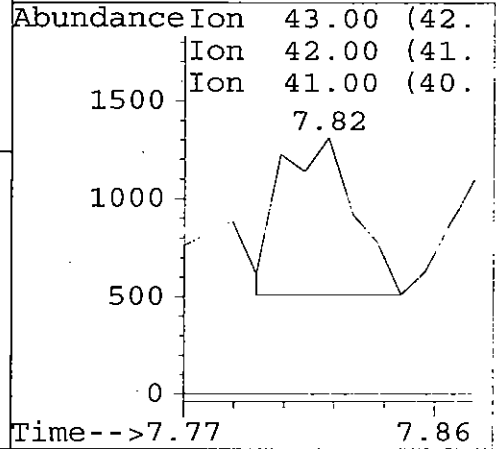
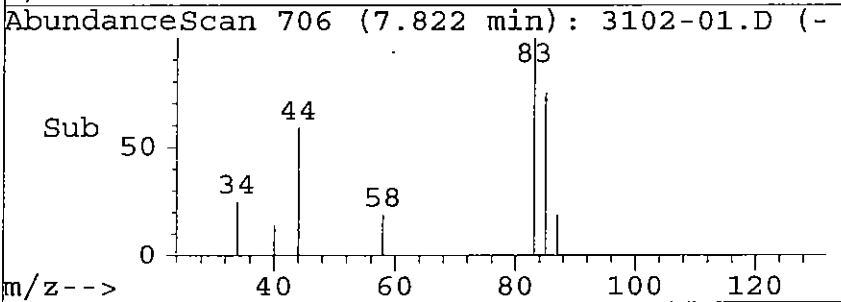




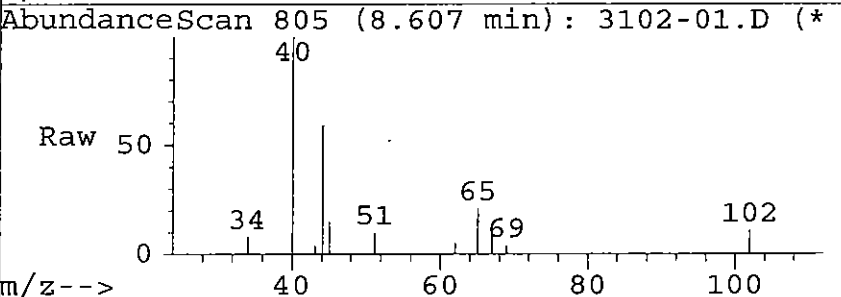
#36  
 117 Iso-butyl alcohol X10  
 Concen: 14.30 ppb  
 RT: 7.82 min Scan# 706  
 Delta R.T. -0.01 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm



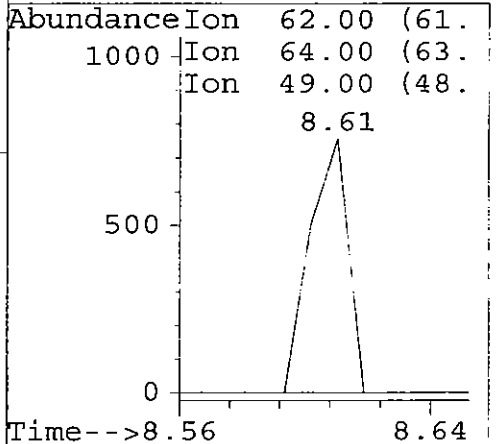
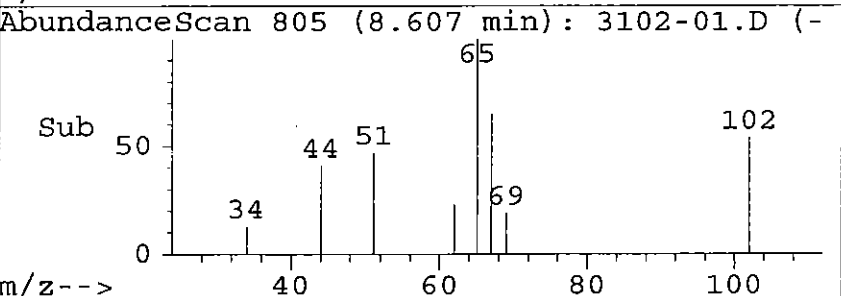
Tgt Ion	Resp	Lower	Upper
43	1343		
42	0.0	7.3	10.9#
41	0.0	39.7	59.5#
0	0.0	0.0	0.0

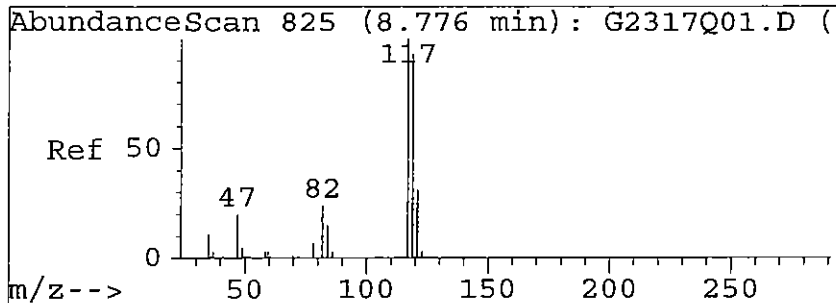


#40  
 30 12-dichloroethane 64 62  
 Concen: 0.92 ppb  
 RT: 8.61 min Scan# 805  
 Delta R.T. -0.03 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm



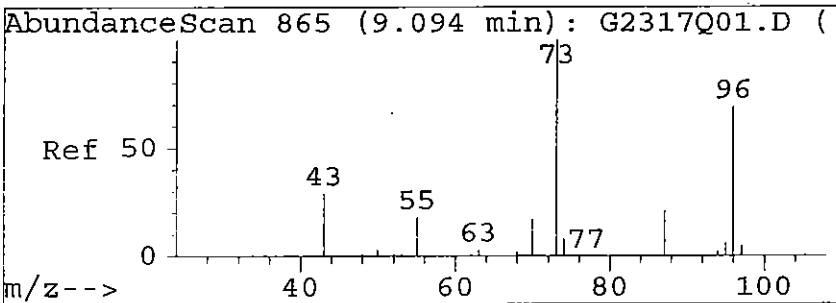
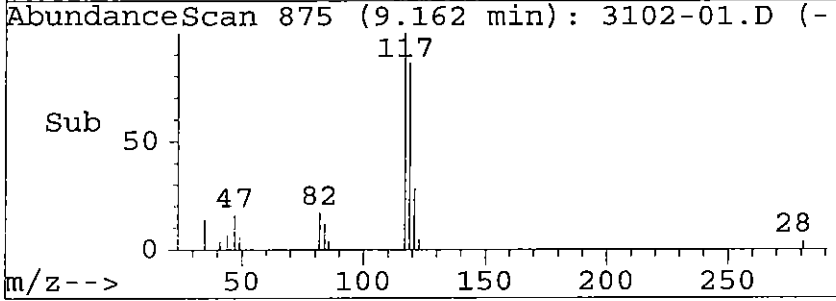
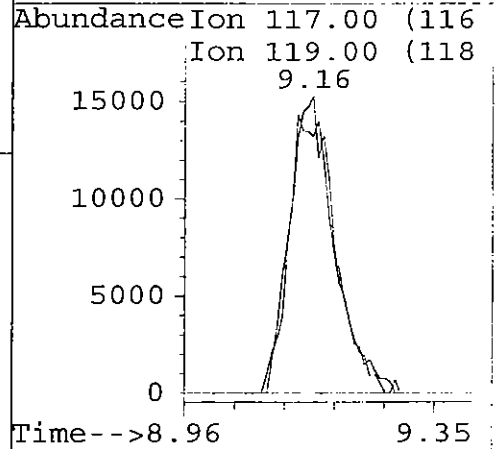
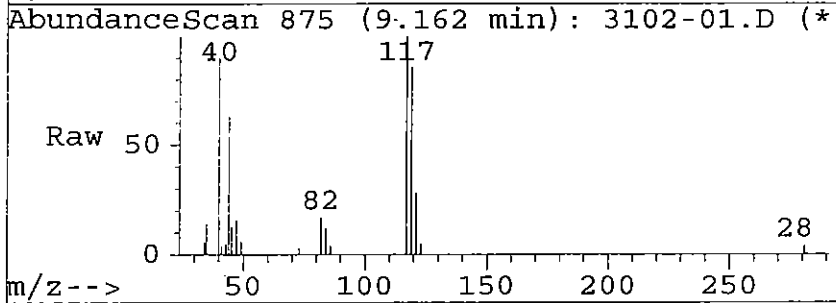
Tgt Ion	Resp	Lower	Upper
62	601		
64	49.8	13.1	53.1
49	1052.7	16.2	56.2#
0	0.0	0.0	0.0





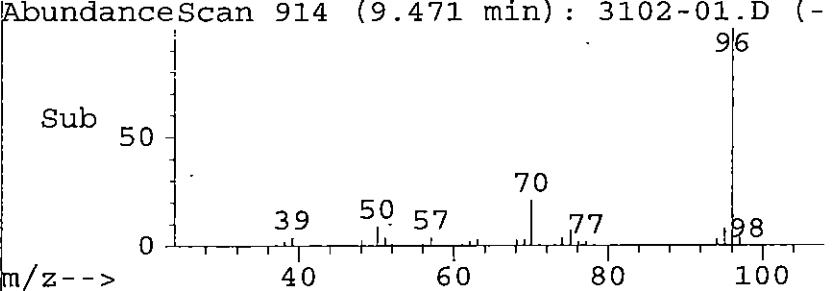
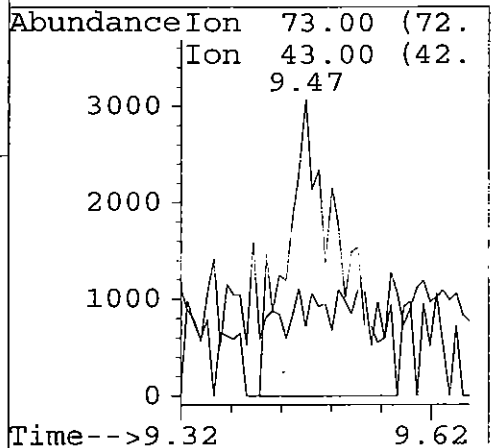
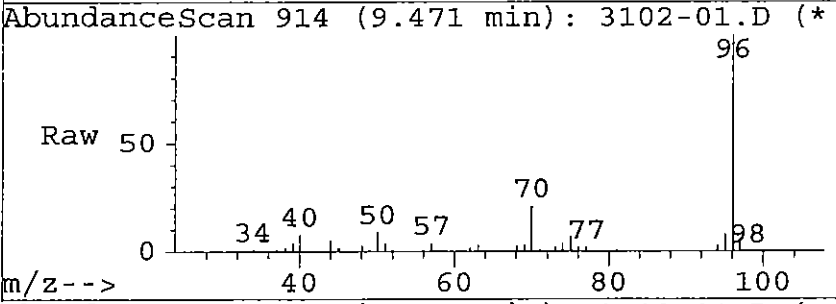
#44  
 37 CCl4 117 119  
 Concen: 2.58 ppb  
 RT: 9.16 min Scan# 875  
 Delta R.T. -0.03 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
117	73832		
119	96.2	80.3	120.3
0	0.0	0.0	0.0
0	0.0	0.0	0.0

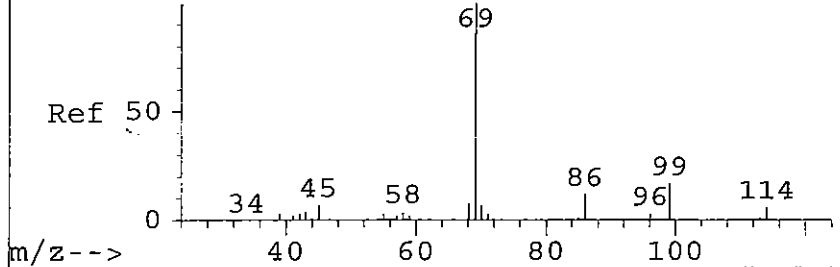


#46  
 118 TAME  
 Concen: 0.40 ppb  
 RT: 9.47 min Scan# 914  
 Delta R.T. -0.00 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
73	13982		
43	14.6	25.0	37.5#
0	0.0	0.0	0.0
0	0.0	0.0	0.0



AbundanceScan 1162 (11.452 min): G2317Q01.D

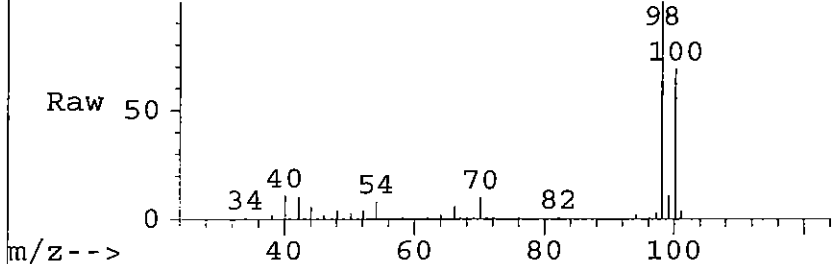


#56

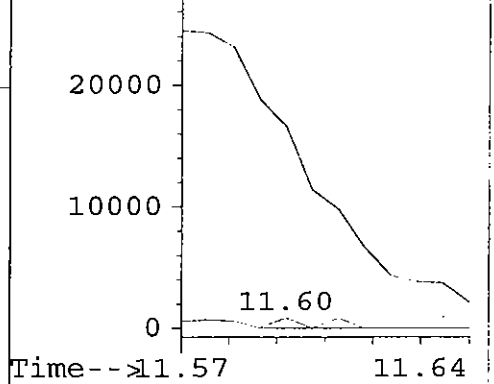
107 Et methacrylate  
 Concen: 1.66 ppb  
 RT: 11.60 min Scan# 1183  
 Delta R.T. -0.20 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
69	100		
99	0.0	9.4	28.0#
0	0.0	0.0	0.0
0	0.0	0.0	0.0

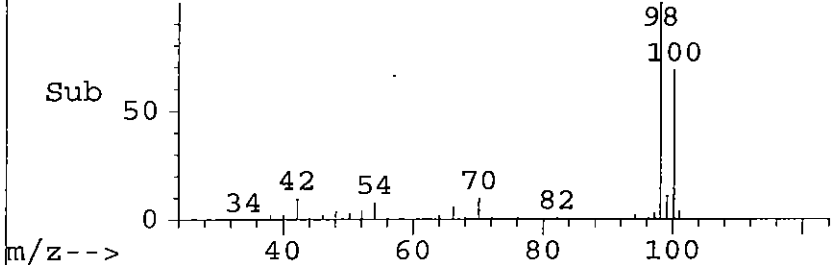
AbundanceScan 1183 (11.602 min): 3102-01.D



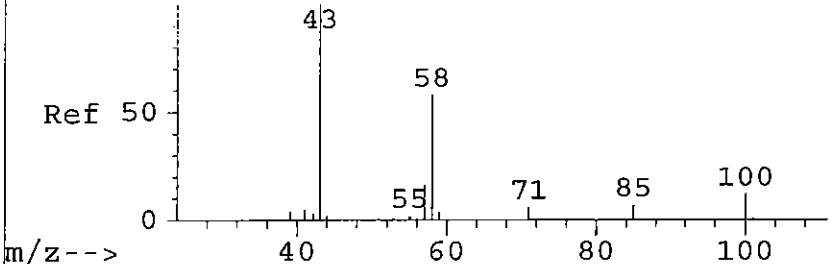
Abundance Ion 69.00 (68.  
 Ion 99.00 (98.



AbundanceScan 1183 (11.602 min): 3102-01.D



AbundanceScan 1178 (11.579 min): G2317Q01.D

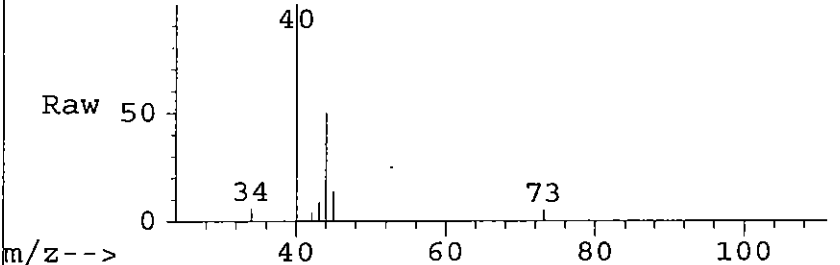


#57

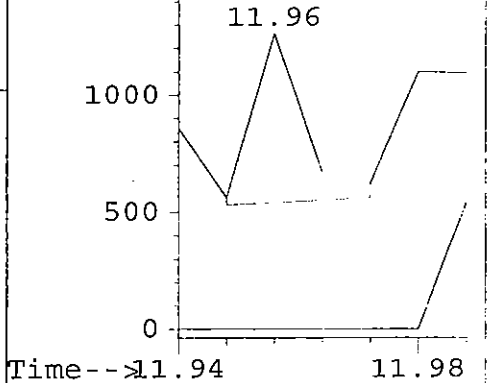
93 2-Hexanone x5  
 Concen: 5.71 ppb  
 RT: 11.96 min Scan# 1228  
 Delta R.T. 0.03 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
43	100		
58	59.1	44.5	66.8
0	0.0	0.0	0.0
0	0.0	0.0	0.0

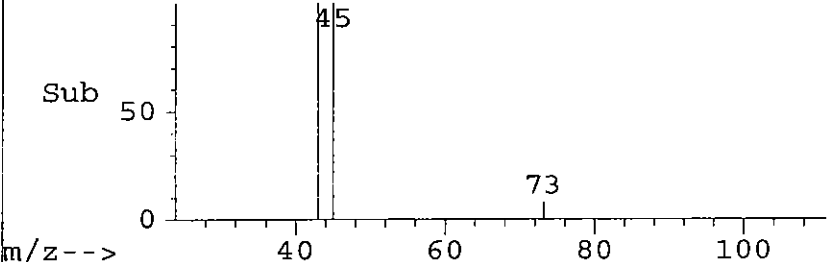
AbundanceScan 1228 (11.959 min): 3102-01.D



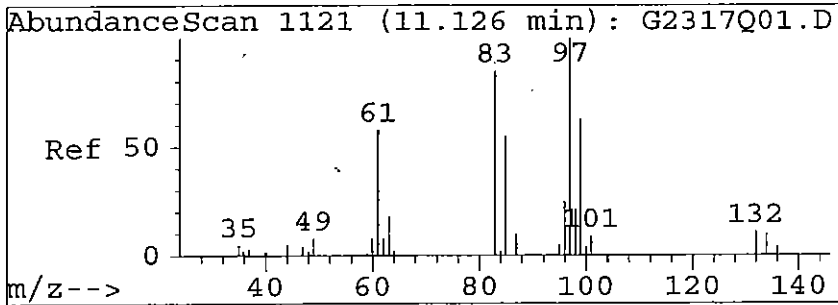
Abundance Ion 43.00 (42.  
 Ion 58.00 (57.



AbundanceScan 1228 (11.959 min): 3102-01.D

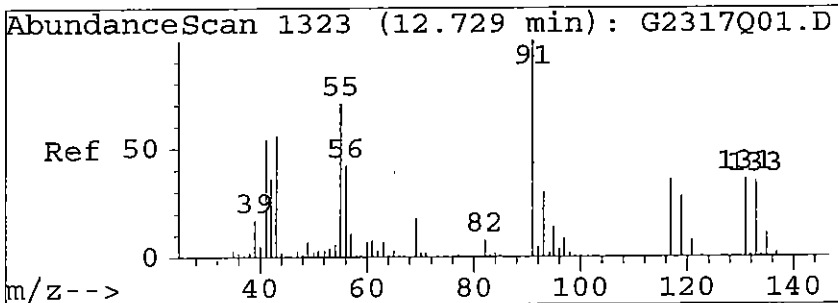
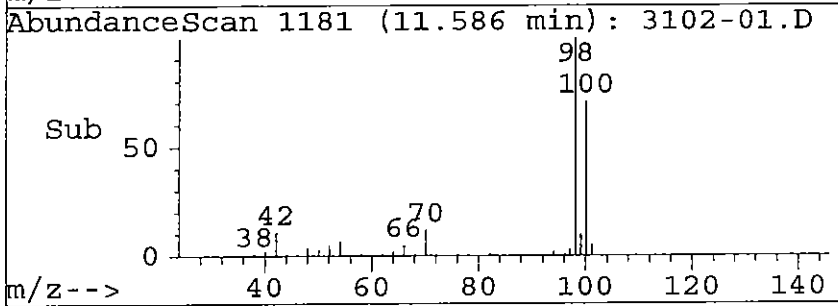
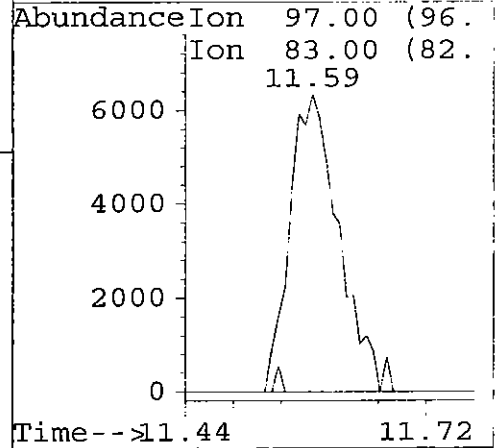
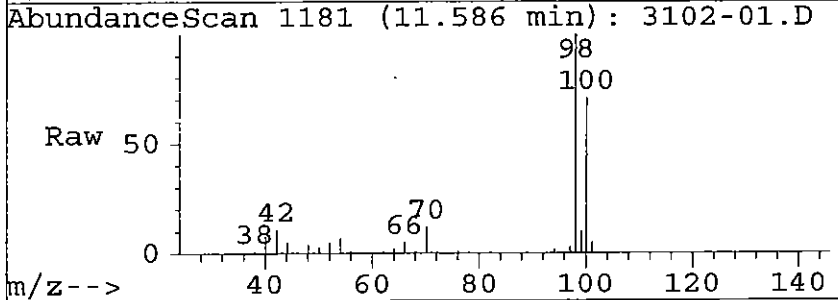






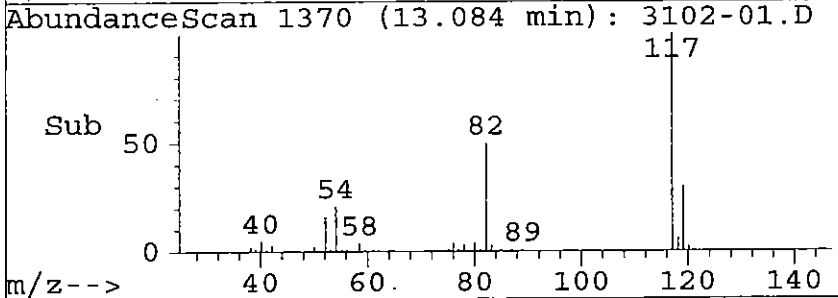
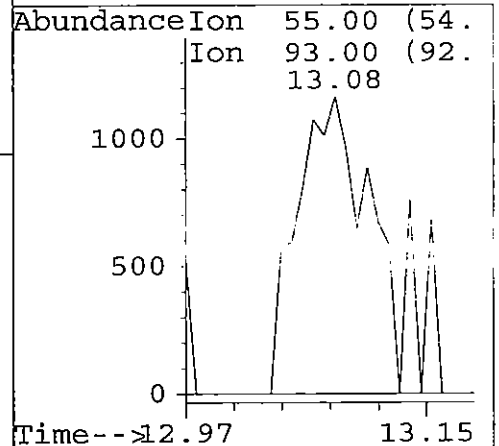
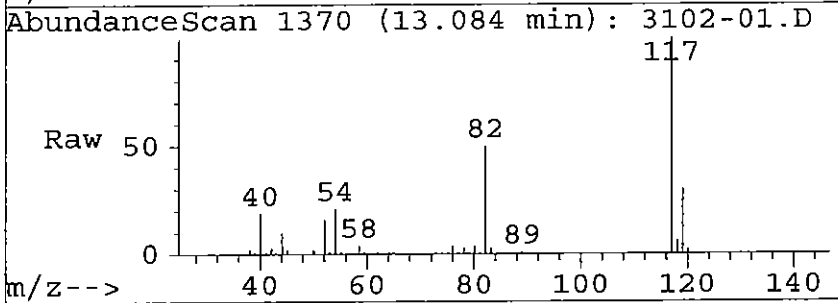
#58  
 48 112-tri-Cl-Et 97 83  
 Concen: 2.72 ppb  
 RT: 11.59 min Scan# 1181  
 Delta R.T. 0.10 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
97	100		
83	1.0	41.5	124.4#
0	0.0	0.0	0.0
0	0.0	0.0	0.0

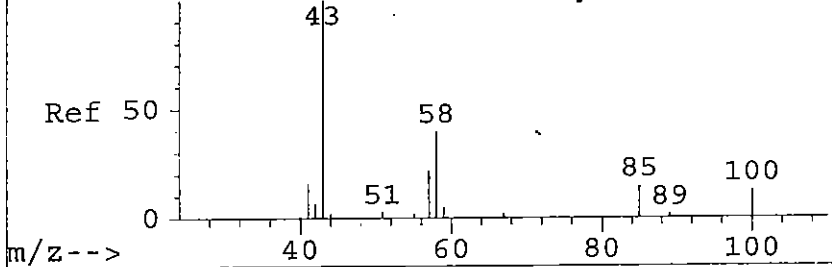


#62  
 105 1-Chlorohexane  
 Concen: 0.21 ppb  
 RT: 13.08 min Scan# 1370  
 Delta R.T. 0.00 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Resp	Lower	Upper
55	100		
93	0.0	0.0	94.3
0	0.0	0.0	0.0
0	0.0	0.0	0.0



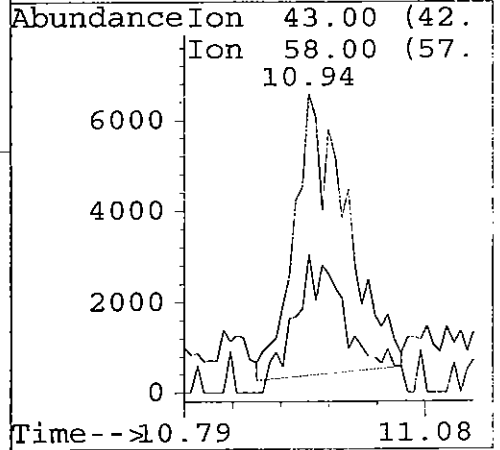
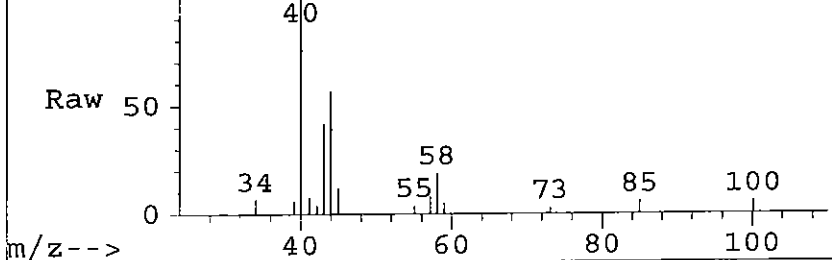
AbundanceScan 1056 (10.610 min): G2317Q01.D



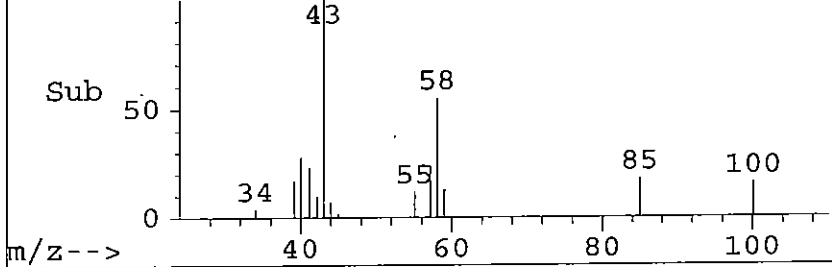
#64  
 54 MIBK  
 Concen: 3.88 ppb  
 RT: 10.94 min Scan# 1099  
 Delta R.T. -0.06 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Ratio	Lower	Upper
43	100		
58	52.0	17.5	57.5
0	0.0	0.0	0.0
0	0.0	0.0	0.0

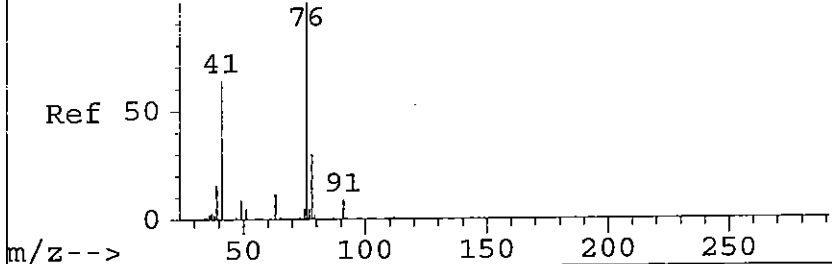
AbundanceScan 1099 (10.937 min): 3102-01.D



AbundanceScan 1099 (10.937 min): 3102-01.D



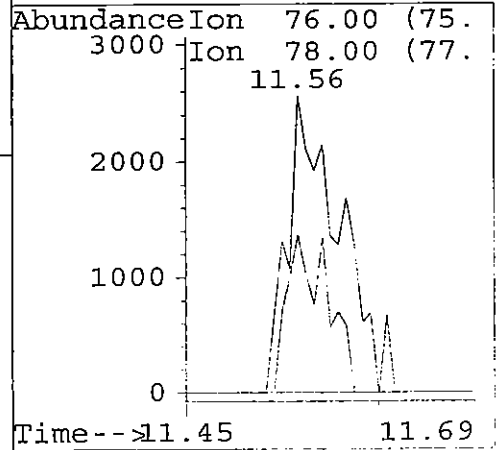
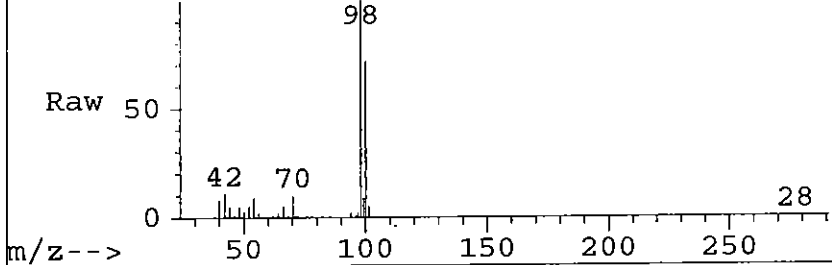
AbundanceScan 1154 (11.388 min): G2317Q01.D



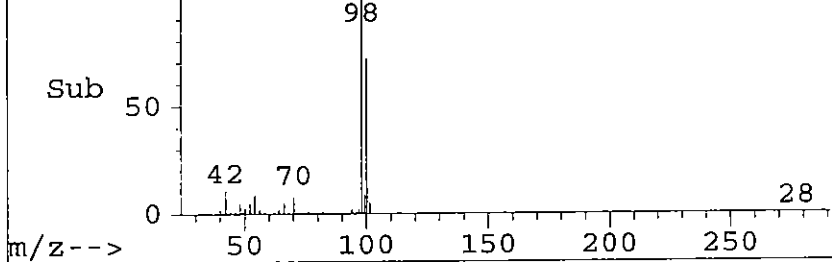
#65  
 49 1,3-di-cl-propane 76 78  
 Concen: 0.60 ppb  
 RT: 11.56 min Scan# 1178  
 Delta R.T. -0.19 min  
 Lab File: 3102-01.D  
 Acq: 16 May 03 5:59 pm

Tgt Ion	Ratio	Lower	Upper
76	100		
78	41.6	27.9	41.8
0	0.0	0.0	0.0
0	0.0	0.0	0.0

AbundanceScan 1178 (11.562 min): 3102-01.D



AbundanceScan 1178 (11.562 min): 3102-01.D



Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: <b>EB-11-5/7/03</b>	Lab Sample ID: 03-3102-2	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-02	Prep. No: -	Anal. Time: 18:28
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	<0.5	U
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	<0.5	U
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROETHENE (TOTAL)	540-59-0	µg/L	0.5	<0.5	U
32	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
33	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
34	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
35	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
36	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
37	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
38	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
39	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U
41	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
42	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	7	J
43	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	<1.8	U
44	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
45	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
46	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
47	STYRENE	100-42-5	µg/L	0.5	<0.5	U
48	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
49	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
50	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
51	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
52	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
53	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
54	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
55	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
56	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
57	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
58	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
59	1,1,2-TRICHLORO-1,1,2,2-TRIFLUOROETHANE	76-13-1	µg/L	0.5	<0.5	U
60	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
61	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
62	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
63	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
64	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4		70-129	96	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	94	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	101	
4	TOLUENE-D8	2037-26-5		73-129	103	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROBENZENE-D5	3114-55-4		50-200	103	
2	1,4-DICHLOROBENZENE-D4	3855-82-1		50-200	118	
3	FLUOROBENZENE	462-06-6		50-200	116	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a)MDL reported.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: MW-12-1	Lab Sample ID: 03-3102-3	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-03	Prep. No: -	Anal. Time: 18:57
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	<0.5	U
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	<0.5	U
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROETHENE (TOTAL)	540-59-0	µg/L	0.5	<0.5	U
32	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
33	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
34	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
35	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
36	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
37	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
38	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
39	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U
41	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
42	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	8	J
43	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	7.2 (b)	B
44	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
45	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
46	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
47	STYRENE	100-42-5	µg/L	0.5	<0.5	U
48	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
49	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
50	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
51	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
52	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
53	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
54	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
55	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
56	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
57	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
58	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
59	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
60	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
61	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
62	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
63	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
64	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4		70-129	94	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	90	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	97	
4	TOLUENE-D8	2037-26-5		73-129	101	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROBENZENE-D5	3114-55-4		50-200	105	
2	1,4-DICHLOROBENZENE-D4	3855-82-1		50-200	117	
3	FLUOROBENZENE	462-06-6		50-200	115	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a) MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: MW-12-2	Lab Sample ID: 03-3102-4	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-04	Prep. No: -	Anal. Time: 19:26
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	<0.5	U
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	<0.5	U
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROETHENE (TOTAL)	540-59-0	µg/L	0.5	<0.5	U
32	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
33	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
34	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
35	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
36	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
37	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
38	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
39	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U
41	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
42	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	5	J
43	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	10 (b)	B
44	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
45	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
46	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
47	STYRENE	100-42-5	µg/L	0.5	<0.5	U
48	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
49	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
50	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
51	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
52	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
53	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
54	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
55	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
56	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
57	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
58	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
59	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
60	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
61	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
62	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
63	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
64	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4		70-129	91	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	92	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	99	
4	TOLUENE-D8	2037-26-5		73-129	104	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROBENZENE-D5	3114-55-4		50-200	104	
2	1,4-DICHLOROETHANE-D4	3855-82-1		50-200	122	
3	FLUOROBENZENE	462-06-6		50-200	115	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a) MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted



Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: MW-12-3	Lab Sample ID: 03-3102-5	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-05	Prep. No: -	Anal. Time: 19:55
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	2.5	
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	1.1	
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
32	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
33	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
34	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
35	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
36	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
37	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
38	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U
39	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
41	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	U
42	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	8.3 (b)	B
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
45	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
46	STYRENE	100-42-5	µg/L	0.5	<0.5	U
47	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
48	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
49	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
50	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
51	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
52	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
53	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
54	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
55	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
56	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
57	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
58	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
59	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
60	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
61	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
62	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
63	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4		70-129	97	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	93	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	98	
4	TOLUENE-D8	2037-26-5		73-129	102	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROBENZENE-D5	3114-55-4		50-200	111	
2	1,4-DICHLOROETHANE-D4	3855-82-1		50-200	123	
3	FLUOROBENZENE	462-06-6		50-200	120	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a)MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

E - Exceed calibration range

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

B - A positive value was found in the method blank

D - Diluted

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: MW-12-4	Lab Sample ID: 03-3102-6	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-06	Prep. No: -	Anal. Time: 20:24
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	1.5	
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	0.7	
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
32	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
33	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
34	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
35	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
36	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
37	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
38	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U
39	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
41	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	U
42	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	11 (b)	B
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
45	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
46	STYRENE	100-42-5	µg/L	0.5	<0.5	U
47	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
48	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
49	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
50	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
51	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
52	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
53	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
54	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
55	TRICHLOROETHENE	79-01-6	µg/L	0.5	0.3	J
56	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
57	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
58	1,1,2,2-TETRACHLOROETHANE	76-13-1	µg/L	0.5	<0.5	U
59	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
60	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
61	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
62	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
63	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U
<b>Surrogates</b>				<b>Control Limit, %</b>	<b>Surro. Rec.%</b>	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4		70-129	99	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	96	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	100	
4	TOLUENE-D8	2037-26-5		73-129	103	
# of out-of-control					0	
<b>Internal Standard</b>				<b>Control Limit, %</b>	<b>IS Rec.%</b>	
1	CHLOROBENZENE-D5	3114-55-4		50-200	105	
2	1,4-DICHLOROETHANE-D4	3855-82-1		50-200	118	
3	FLUOROBENZENE	462-06-6		50-200	117	
# of out-of-control					0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a)MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: MW-12-5	Lab Sample ID: 03-3102-7	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-07	Prep. No: -	Anal. Time: 20:53
Methanol Vol: -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYL BENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYL BENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYL BENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	0.6	
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	0.4	J
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
32	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
33	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
34	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
35	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
36	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
37	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
38	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U
39	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
41	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	7	J
42	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	13 (b)	B
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
45	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
46	STYRENE	100-42-5	µg/L	0.5	<0.5	U
47	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
48	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
49	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
50	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
51	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
52	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
53	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
54	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
55	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
56	TRICHLOROFUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
57	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
58	112TRICHLORO-122TRIFLUOROETHANE	76-13-1	µg/L	0.5	<0.5	U
59	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
60	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
61	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
62	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
63	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U

**Surrogates**

		Control Limit, %	Surro. Rec.%	
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4	70-129	100
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	95
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	101
4	TOLUENE-D8	2037-26-5	73-129	105
# of out-of-control			0	

**Internal Standard**

		Control Limit, %	IS Rec.%	
1	CHLOROBENZENE-D5	3114-55-4	50-200	99
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	112
3	FLUOROBENZENE	462-06-6	50-200	112
# of out-of-control			0	

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a) MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

E - Exceed calibration range

B - A positive value was found in the method blank

D - Diluted

Applied P & Ch Laboratory  
**Organic Analysis Results for Method 524.2**

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/07/2003
Project ID: JPL	Service ID: 33102	Collected by:
Sample ID: <b>TB-11-5/7/03</b>	Lab Sample ID: 03-3102-8	Received Date: 05/07/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: G
Batch No: 03G2404	Prep. Date: 05/16/03	Anal. Date: 05/16/03
Data File Name: 3102-08	Prep. No: -	Anal. Time: 21:22
Methanol Vol. -	Sample Amount: 25 mL	Dilution Factor: 1
Test Level: Low	Sparge Size: 25 mL	Heated Purge: (Y/N) N

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	BENZENE	71-43-2	µg/L	0.5	<0.5	U
2	BROMOBENZENE	108-86-1	µg/L	0.5	<0.5	U
3	BROMOCHLOROMETHANE	74-97-5	µg/L	0.5	<0.5	U
4	BROMODICHLOROMETHANE	75-27-4	µg/L	0.5	<0.5	U
5	BROMOFORM	75-25-2	µg/L	0.5	<0.5	U
6	BROMOMETHANE	74-83-9	µg/L	0.5	<0.5	U
7	2-BUTANONE	78-93-3	µg/L	10	<10	U
8	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
9	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
10	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	<0.5	U
12	CHLOROBENZENE	108-90-7	µg/L	0.5	<0.5	U
13	CHLORODIBROMOMETHANE	124-48-1	µg/L	0.5	<0.5	U
14	CHLOROETHANE	75-00-3	µg/L	0.5	<0.5	U
15	CHLOROFORM	67-66-3	µg/L	0.5	<0.5	U
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
18	4-CHLOROTOLUENE	106-43-4	µg/L	0.5	<0.5	U
19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	1.1 (a)	<1.1	U
20	1,2-DIBROMOETHANE (EDB)	106-93-4	µg/L	0.5	<0.5	U
21	DIBROMOMETHANE	74-95-3	µg/L	0.5	<0.5	U
22	1,2-DICHLOROBENZENE	95-50-1	µg/L	0.5	<0.5	U
23	1,3-DICHLOROBENZENE	541-73-1	µg/L	0.5	<0.5	U
24	1,4-DICHLOROBENZENE	106-46-7	µg/L	0.5	<0.5	U
25	DICHLORODIFLUOROMETHANE	75-71-8	µg/L	0.5	<0.5	U
26	1,1-DICHLOROETHANE	75-34-3	µg/L	0.5	<0.5	U
27	1,2-DICHLOROETHANE	107-06-2	µg/L	0.5	<0.5	U
28	1,1-DICHLOROETHENE	75-35-4	µg/L	0.5	<0.5	U
29	CIS-1,2-DICHLOROETHENE	156-59-2	µg/L	0.5	<0.5	U
30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
31	1,2-DICHLOROPROPANE	78-87-5	µg/L	0.5	<0.5	U
32	1,3-DICHLOROPROPANE	142-28-9	µg/L	0.5	<0.5	U
33	2,2-DICHLOROPROPANE	594-20-7	µg/L	0.5	<0.5	U
34	1,1-DICHLOROPROPENE	563-58-6	µg/L	0.5	<0.5	U
35	CIS-1,3-DICHLOROPROPENE	10061-01-5	µg/L	0.5	<0.5	U
36	TRANS-1,3-DICHLOROPROPENE	10061-02-6	µg/L	0.5	<0.5	U
37	ETHYLBENZENE	100-41-4	µg/L	0.5	<0.5	U
38	HEXACHLOROBUTADIENE	87-68-3	µg/L	0.5	<0.5	U
39	ISOPROPYLBENZENE (CUMENE)	98-82-8	µg/L	0.5	<0.5	U

#	Component Name	CAS No	Unit	RL	Result	Qualifier
40	P-ISOPROPYLTOLUENE	99-87-6	µg/L	0.5	<0.5	U
41	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	U
42	METHYLENE CHLORIDE	75-09-2	µg/L	1.8 (a)	6.8 (b)	B
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	NAPHTHALENE	91-20-3	µg/L	0.5	<0.5	U
45	N-PROPYLBENZENE	103-65-1	µg/L	0.5	<0.5	U
46	STYRENE	100-42-5	µg/L	0.5	<0.5	U
47	1,1,1,2-TETRACHLOROETHANE	630-20-6	µg/L	0.5	<0.5	U
48	1,1,2,2-TETRACHLOROETHANE	79-34-5	µg/L	0.5	<0.5	U
49	TETRACHLOROETHENE	127-18-4	µg/L	0.5	<0.5	U
50	TOLUENE	108-88-3	µg/L	0.5	<0.5	U
51	1,2,3-TRICHLOROBENZENE	87-61-6	µg/L	0.5	<0.5	U
52	1,2,4-TRICHLOROBENZENE	120-82-1	µg/L	0.5	<0.5	U
53	1,1,1-TRICHLOROETHANE	71-55-6	µg/L	0.5	<0.5	U
54	1,1,2-TRICHLOROETHANE	79-00-5	µg/L	0.5	<0.5	U
55	TRICHLOROETHENE	79-01-6	µg/L	0.5	<0.5	U
56	TRICHLOROFLUOROMETHANE	75-69-4	µg/L	0.5	<0.5	U
57	1,2,3-TRICHLOROPROPANE	96-18-4	µg/L	0.5	<0.5	U
58	1,1,2,2-TRICHLORO-1,1,2,2-TRIFLUOROETHANE	76-13-1	µg/L	0.5	<0.5	U
59	1,2,4-TRIMETHYLBENZENE	95-63-6	µg/L	0.5	<0.5	U
60	1,3,5-TRIMETHYLBENZENE	108-67-8	µg/L	0.5	<0.5	U
61	VINYL CHLORIDE	75-01-4	µg/L	0.5	<0.5	U
62	O-XYLENE	95-47-6	µg/L	0.5	<0.5	U
63	M/P-XYLENE	108-38-3	µg/L	0.5	<0.5	U

**Surrogates**

			Control Limit, %	Surro. Rec.%
1	1-BROMO-4-FLUOROBENZENE (4-BROMOFL)	460-00-4	70-129	97
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	91
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	99
4	TOLUENE-D8	2037-26-5	73-129	103
# of out-of-control				0

**Internal Standard**

			Control Limit, %	IS Rec.%
1	CHLOROBENZENE-D5	3114-55-4	50-200	108
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	121
3	FLUOROBENZENE	462-06-6	50-200	119
# of out-of-control				0

Not Detected is shown as PQL, with dilution and moisture corrected if applicable.

(a)MDL reported.

(b) Laboratory contamination suspected.

Qualifier: U - Not Detected or less than MDL

J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)

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