

**Laboratory Data Consultants, Inc.  
Data Validation Report**

**Project/Site Name:** NASA JPL  
**Collection Date:** October 28, 2003  
**LDC Report Date:** December 12, 2003  
**Matrix:** Water  
**Parameters:** Volatiles  
**Validation Level:** EPA Level III & IV  
**Laboratory:** Applied P & Ch Laboratory  
**Sample Delivery Group (SDG):** 03-5852

**Sample Identification**

Dupe-1-4Q03\*\*  
EB-5-10-28-03  
MW-24-1  
MW-24-2  
MW-24-3  
MW-24-4  
MW-24-5  
TB-5-10-28-03

\*\*Indicates sample underwent EPA Level IV review

## Introduction

This data review covers 8 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 524.2 for Volatiles.

This review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review (October 1999) as there are no current guidelines for the method stated above.

A table summarizing all data qualification is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

Blank results are summarized in Section V.

Field duplicates are summarized in Section XVI.

Samples indicated by a double asterisk on the front cover underwent a EPA Level IV review. A EPA Level III review was performed on all of the other samples. Raw data were not evaluated for the samples reviewed by Level III criteria since this review is based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. GC/MS Instrument Performance Check

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

## III. Initial Calibration

Initial calibration was performed using required standard concentrations.

Percent relative standard deviations (%RSD) were less than or equal to 20.0% for selected compounds.

A curve fit, based on the initial calibration, was established for quantitation for selected compounds. The coefficient of determination ( $r^2$ ) was greater than or equal to 0.990 .

## IV. Continuing Calibration

Continuing calibration was performed at the required frequencies.

All of the continuing calibration percent differences (%D) between the initial calibration RRF and the continuing calibration RRF were less than or equal to 30.0% with the following exceptions:

Date	Compound	%D	Associated Samples	Flag	A or P
11/1/03	Chloromethane	34.05	All samples in SDG 03-5852	J (all detects)	P
	Bromomethane	59.05		UJ (all non-detects)	
				J (all detects)	
				UJ (all non-detects)	

## V. Blanks

Method blanks were reviewed for each matrix as applicable. No volatile contaminants were found in the method blanks with the following exceptions:

Method Blank ID	Analysis Date	Compound TIC (RT in minutes)	Concentration	Associated Samples
03G4668MB01	11/1/03	2-Butanone	0.61 ug/L	All samples in SDG 03-5852

Sample concentrations were compared to concentrations detected in the method blanks. The sample concentrations were either not detected or were significantly greater (>10X for common contaminants, >5X for other contaminants) than the concentrations found in the associated method blanks.

#### **VI. Surrogate Spikes**

Surrogates were added to all samples and blanks as required by the method. All surrogate recoveries (%R) were within QC limits.

#### **VII. Matrix Spike/Matrix Spike Duplicates**

Although matrix spike (MS) and matrix spike duplicate (MSD) samples were not required by the method, MS and MSD samples were reported by the laboratory. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits.

#### **VIII. Laboratory Control Samples (LCS)**

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

#### **IX. Regional Quality Assurance and Quality Control**

Not applicable.

#### **X. Internal Standards**

All internal standard areas and retention times were within QC limits.

#### **XI. Target Compound Identifications**

All target compound identifications were within validation criteria for samples on which a EPA Level IV review was performed. Raw data were not evaluated for the samples reviewed by Level III criteria.

#### **XII. Compound Quantitation and CRQLs**

All compound quantitation and CRQLs were within validation criteria for samples on which a EPA Level IV review was performed. Raw data were not evaluated for the samples reviewed by Level III criteria.

#### **XIII. Tentatively Identified Compounds (TICs)**

Tentatively identified compounds were not reported by the laboratory.

#### **XIV. System Performance**

The system performance was within validation criteria for samples on which a EPA Level IV review was performed. Raw data were not evaluated for the samples reviewed by Level III criteria.

#### **XV. Overall Assessment of Data**

Data flags have been summarized at the end of the report.

#### **XVI. Field Duplicates**

Samples Dupe-1-4Q03\*\* and MW-24-4 were identified as field duplicates. No volatiles were detected in any of the samples.

#### **XVII. Field Blanks**

Sample TB-5-10-28-03 was identified as a trip blank. No volatile contaminants were found in this blank.

Sample EB-5-10-28-03 was identified as an equipment blank. No volatile contaminants were found in this blank.

**NASA JPL**  
**Volatiles - Data Qualification Summary - SDG 03-5852**

SDG	Sample	Compound	Flag	A or P	Reason
03-5852	Dupe-1-4Q03** EB-5-10-28-03 MW-24-1 MW-24-2 MW-24-3 MW-24-4 MW-24-5 TB-5-10-28-03	Chloromethane  Bromomethane	J (all detects) UJ (all non-detects) J (all detects) UJ (all non-detects)	P	Continuing calibration (%D)

**NASA JPL**  
**Volatiles - Laboratory Blank Data Qualification Summary - SDG 03-5852**

No Sample Data Qualified in this SDG

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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: DUPE-1-4Q03	Lab Sample ID: 03-5852-1	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-01	Prep. No: -	Anal. Time: 07:35
Methanol Vol: -	Sample Amount: 25.0 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 <i>UJ</i>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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 <i>UJ</i>
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	0.5	<0.5	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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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	98	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	97	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	99	
4	TOLUENE-D8	2037-26-5		73-129	105	
# 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	101	
2	1,4-DICHLOROBENZENE-D4	3855-82-1		50-200	107	
3	FLUOROBENZENE	462-06-6		50-200	108	
# 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

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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: <b>EB-5-10-28-03</b>	Lab Sample ID: 03-5852-2	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-02	Prep. No: -	Anal. Time: 08:01
Methanol Vol: -	Sample Amount: 25.0 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 <i>UT</i>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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 <i>U</i>
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	0.5	<0.5	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

*1/2/03*

#	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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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
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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-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	97	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	99	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	99	
4	TOLUENE-D8	2037-26-5		73-129	105	
# 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	101	
2	1,4-DICHLOROETHANE-D4	3855-82-1		50-200	108	
3	FLUOROBENZENE	462-06-6		50-200	108	
# of out-of-control					0	

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

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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

*M/DH*

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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: MW-24-1	Lab Sample ID: 03-5852-3	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-03	Prep. No: -	Anal. Time: 08:27
Methanol Vol: -	Sample Amount: 25.0 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
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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 <sup>W</sup>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	< 0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	< 0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	< 0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	< 10	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	19.1	
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	6.8	
16	CHLOROMETHANE	74-87-3	µg/L	0.5	< 0.5	U <sup>W</sup>
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	0.5	< 0.5	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	0.7	
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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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	1.6	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	3.7	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-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	97	
2	1,2-DICHLOROETHANE-D4	17060-07-0		70-129	98	
3	DIBROMOFLUOROMETHANE	1868-53-7		70-122	99	
4	TOLUENE-D8	2037-26-5		73-129	105	
# 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	100	
2	1,4-DICHLOROETHANE-D4	3855-82-1		50-200	105	
3	FLUOROBENZENE	462-06-6		50-200	107	
# 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

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

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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: MW-24-2	Lab Sample ID: 03-5852-4	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-04	Prep. No: -	Anal. Time: 08:54
Methanol Vol. -	Sample Amount: 25.0 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 <i>UJ</i>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	U
11	CARBON TETRACHLORIDE	56-23-5	µg/L	0.5	3.4	
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	1.4	
16	CHLOROMETHANE	74-87-3	µg/L	0.5	<0.5	U <i>UJ</i>
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	0.5	<0.5	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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
43	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
44	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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	0.6	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	98
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	97
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	99
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	100
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	106
3	FLUOROBENZENE	462-06-6	50-200	108
# 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  
 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: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: MW-24-3	Lab Sample ID: 03-5852-5	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-05	Prep. No: -	Anal. Time: 09:20
Methanol Vol: -	Sample Amount: 25.0 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 UT
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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 UT
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	0.5	<0.5	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-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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
42	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
43	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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	98
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	97
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	99
4	TOLUENE-D8	2037-26-5	73-129	104
# of out-of-control			0	

## Internal Standard

		Control Limit, %	IS Rec.%	
1	CHLOROBENZENE-D5	3114-55-4	50-200	100
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	105
3	FLUOROBENZENE	462-06-6	50-200	107
# 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

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

10/12/03



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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: MW-24-4	Lab Sample ID: 03-5852-6	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-06	Prep. No: -	Anal. Time: 09:46
Methanol Vol: -	Sample Amount: 25.0 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 <i>WT</i>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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 <i>WT</i>
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	0.5	<0.5	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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
42	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
43	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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-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	98
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	99
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	100
4	TOLUENE-D8	2037-26-5	73-129	104
	# of out-of-control			0

## Internal Standard

		Control Limit, %	IS Rec.%	
1	CHLOROBENZENE-D5	3114-55-4	50-200	101
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	107
3	FLUOROBENZENE	462-06-6	50-200	107
	# 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

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

9/12/06

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

Client Name: GEOFON, Inc.	Project No: 04.4428.10	Collection Date: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: MW-24-5	Lab Sample ID: 03-5852-7	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-07	Prep. No: -	Anal. Time: 10:12
Methanol Vol. -	Sample Amount: 25.0 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	UW
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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	UW
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	0.5	<0.5	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-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	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
42	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
43	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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-TRICHLORO-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	98
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	98
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	100
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	107
3	FLUOROBENZENE	462-06-6	50-200	107
# 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

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: 10/28/2003
Project ID: JPL	Service ID: 35852	Collected by: Jr
Sample ID: TB-5-10-28-03	Lab Sample ID: 03-5852-8	Received Date: 10/28/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 524.2	Prep. Method: 5030	Instrument ID: GC/MS: A
Batch No: 03G4668	Prep. Date: 11/01/03	Anal. Date: 11/01/03
Data File Name: 5852-08	Prep. No: -	Anal. Time: 04:59
Methanol Vol: -	Sample Amount: 25.0 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 <i>UJ</i>
7	N-BUTYLBENZENE	104-51-8	µg/L	0.5	<0.5	U
8	SEC-BUTYLBENZENE	135-98-8	µg/L	0.5	<0.5	U
9	TERT-BUTYLBENZENE	98-06-6	µg/L	0.5	<0.5	U
10	2-BUTANONE	78-93-3	µg/L	10	<10	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 <i>UJ</i>
17	2-CHLOROTOLUENE	95-49-8	µg/L	0.5	<0.5	U
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19	1,2-DIBROMO-3-CHLOROPROPANE	96-12-8	µg/L	0.5	<0.5	U
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30	TRANS-1,2-DICHLOROETHENE	156-60-5	µg/L	0.5	<0.5	U
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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
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41	METHYLENE CHLORIDE	75-09-2	µg/L	0.5	<0.5	U
42	METHYL-T-BUTYL ETHER (MTBE)	1634-04-4	µg/L	1	<1	U
43	4-METHYL-2-PENTANONE (MIBK)	108-10-1	µg/L	10	<10	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	96
2	1,2-DICHLOROETHANE-D4	17060-07-0	70-129	99
3	DIBROMOFLUOROMETHANE	1868-53-7	70-122	98
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	104
2	1,4-DICHLOROETHANE-D4	3855-82-1	50-200	111
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.

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

LDC #: 11272D1

## VALIDATION COMPLETENESS WORKSHEET

SDG #: 03-5852

Level III/IV

Laboratory: Applied P &amp; Ch Laboratory

Date: 12/12/03

Page: 1 of 1

Reviewer: [Signature]

2nd Reviewer: [Signature]

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 10/28/03
II.	GC/MS Instrument performance check	A	
III.	Initial calibration	A	70 RSD. Y <sup>2</sup> NO RPF'S
IV.	Continuing calibration	W	
V.	Blanks	W	
VI.	Surrogate spikes	A	
VII.	Matrix spike/Matrix spike duplicates	A	MW-24-4 (03-5892)
VIII.	Laboratory control samples	A	LCS
IX.	Regional Quality Assurance and Quality Control	N	
X.	Internal standards	A	
XI.	Target compound identification	A	Not reviewed for Level III validation.
XII.	Compound quantitation/CRQLs	A	Not reviewed for Level III validation.
XIII.	Tentitatively identified compounds (TICs)	N	Not reviewed for Level III validation.
XIV.	System performance	A	Not reviewed for Level III validation.
XV.	Overall assessment of data	A	
XVI.	Field duplicates	ND	D = 1 + 6
XVII.	Field blanks	ND	EB = 2. TB = 8

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples: \*\* Indicates sample underwent Level IV validation.

1	Dupe-1-4Q03**	11	03G4668MB01	21		31	
2	EB-5-10-28-03	12		22		32	
3	MW-24-1	13		23		33	
4	MW-24-2	14		24		34	
5	MW-24-3	15		25		35	
6	MW-24-4	16		26		36	
7	MW-24-5	17		27		37	
8	TB-5-10-28-03	18		28		38	
9		19		29		39	
10		20		30		40	

LDC #: 1127201  
 SDG #: 03-5852

VALIDATION FINDINGS CHECKLIST

Page: 1 of 2  
 Reviewer: [Signature]  
 2nd Reviewer: [Signature]

**Method: Volatiles (EPA Method 524.2)**

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Cooler temperature criteria was met.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>II. GC/MS Instrument performance check</b>				
Were the BFB performance results reviewed and found to be within the specified criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all samples analyzed within the 12 hour clock criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>III. Initial calibration</b>				
Did the laboratory perform a 5 point calibration prior to sample analysis?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent relative standard deviations (%RSD) $\leq$ 20%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>IV. Continuing calibration</b>				
Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent differences (%D) $\leq$ 30%?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
<b>V. Blanks</b>				
Was a method blank associated with every sample in this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was a method blank analyzed at least once every 12 hours for each matrix and concentration?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>VI. Surrogate spikes</b>				
Were all surrogate %R within QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R outside of criteria?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>VII. Matrix spike/Matrix spike duplicates</b>				
Was a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>VIII. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was an LCS analyzed per analytical batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	



LDC #: 1127201  
 SDG #: 03-5852

VALIDATION FINDINGS CHECKLIST

Page: 2 of 2  
 Reviewer: [Signature]  
 2nd Reviewer: [Signature]

Validation Area	Yes	No	NA	Findings/Comments
<b>IX. Regional Quality Assurance and Quality Control</b>				
Were performance evaluation (PE) samples performed?		/		
Were the performance evaluation (PE) samples within the acceptance limits?			/	
<b>X. Internal standards</b>				
Were internal standard area counts within +/-40% from the associated calibration standard?	/			
Were retention times within - 30% of the last continuing calibration or +/- 50% of the initial calibration?	/			
<b>XI. Target compound identification</b>				
Were relative retention times (RRT's) within $\pm$ 0.06 RRT units of the standard?	/			
Did compound spectra meet specified EPA "Functional Guidelines" criteria?	/			
Were chromatogram peaks verified and accounted for?	/			
<b>XII. Compound quantitation/CRQLs</b>				
Were the correct internal standard (IS), quantitation ion and relative response factor (RRF) used to quantitate the compound?	/			
Were compound quantitation and CRQLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	/			
<b>XIII. Tentatively identified compounds (TICs)</b>				
Were the major ions (> 25 percent relative intensity) in the reference spectrum evaluated in sample spectrum?			/	
Were relative intensities of the major ions within $\pm$ 20% between the sample and the reference spectra?			/	
Did the raw data indicate that the laboratory performed a library search for all required peaks in the chromatograms (samples and blanks)?		/		
<b>XIV. System performance</b>				
System performance was found to be acceptable.	/			
<b>XV. Overall assessment of data</b>				
Overall assessment of data was found to be acceptable.	/			
<b>XVI. Field duplicates</b>				
Field duplicate pairs were identified in this SDG.	/			
Target compounds were detected in the field duplicates.		/		
<b>XVII. Field blanks</b>				
Field blanks were identified in this SDG.	/			
Target compounds were detected in the field blanks.		/		

# TARGET COMPOUND WORKSHEET

METHOD: VOA (EPA Method 524.2)

A. Chloromethane	Q. 1,2-Dichloropropane	GG. Xylenes, total	WW. Bromobenzene	MMM. Naphthalene
B. Bromomethane	R. cis-1,3-Dichloropropene	HH. Vinyl acetate	XX. 1,2,3-Trichloropropane	NNN. 1,2,3-Trichlorobenzene
C. Vinyl chloride	S. Trichloroethene	II. 2-Chloroethylvinyl ether	YY. n-Propylbenzene	OOO. 1,3,5-Trichlorobenzene
D. Chloroethane	T. Dibromochloromethane	JJ. Dichlorodifluoromethane	ZZ. 2-Chlorotoluene	PPP. trans-1,2-Dichloroethene
E. Methylene chloride	U. 1,1,2-Trichloroethane	KK. Trichlorofluoromethane	AAA. 1,3,5-Trimethylbenzene	QQQ. cis-1,2-Dichloroethene
F. Acetone	V. Benzene	LL. Methyl-tert-butyl ether	BBB. 4-Chlorotoluene	RRR. m,p-Xylenes
G. Carbon disulfide	W. trans-1,3-Dichloropropene	MM. 1,2-Dibromo-3-chloropropane	CCC. tert-Butylbenzene	SSS. o-Xylene
H. 1,1-Dichloroethane	X. Bromoform	NN. Diethyl ether	DDD. 1,2,4-Trimethylbenzene	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane
I. 1,1-Dichloroethane	Y. 4-Methyl-2-pentanone	OO. 2,2-Dichloropropane	EEE. sec-Butylbenzene	UUU. Benzyl chloride
J. 1,2-Dichloroethene, total	Z. 2-Hexanone	PP. Bromochloromethane	FFF. 1,3-Dichlorobenzene	VVV. 4-Ethyltoluene
K. Chloroform	AA. Tetrachloroethene	QQ. 1,1-Dichloropropene	GGG. p-Isopropyltoluene	WWW. Ethanol
L. 1,2-Dichloroethane	BB. 1,1,2,2-Tetrachloroethane	RR. Dibromomethane	HHH. 1,4-Dichlorobenzene	XXX. Ethyl ether
M. 2-Butanone	CC. Toluene	SS. 1,3-Dichloropropane	III. n-Butylbenzene	
N. 1,1,1-Trichloroethane	DD. Chlorobenzene	TT. 1,2-Dibromoethane	JJJ. 1,2-Dichlorobenzene	
O. Carbon tetrachloride	EE. Ethylbenzene	UU. 1,1,1,2-Tetrachloroethane	KKK. 1,2,4-Trichlorobenzene	
P. Bromodichloromethane	FF. Styrene	VV. Isopropylbenzene	LLL. Hexachlorobutadiene	

Notes:

### VALIDATION FINDINGS WORKSHEET

#### Continuing Calibration

METHOD: GC/MS VOA (EPA Method 524.2)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Y N N/A Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?

Y N N/A Were all percent differences (%D) ≤ 30%?

#	Date	Standard ID	Compound	Finding %D (Limit: ≤30.0%)	Associated Samples	Qualifications
	11/1/03	94668001	A	37.05	M+ Bk	<u>Y</u> N <u>N</u> P
			B	59.05		↓

**VALIDATION FINDINGS WORKSHEET**  
**Blanks**

METHOD: GC/MS VOA (EPA Method 524.2)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".  
 Y  N  N/A  Was a method blank associated with every sample in this SDG?  
 Y  N  N/A  Was a method blank analyzed at least once every 12 hours for each matrix and concentration?  
 Y  N  N/A  Was there contamination in the method blanks? If yes, please see the qualifications below.

Blank analysis date: 11/1/03  
 Conc. units: ppb Associated Samples: AA

Compound	Blank ID	Sample Identification
	039-668MBP	
Methylene chloride		
Acetone	0.61	
M		
CRQL		

Blank analysis date: \_\_\_\_\_  
 Conc. units: \_\_\_\_\_ Associated Samples: \_\_\_\_\_

Compound	Blank ID	Sample Identification
Methylene chloride		
Acetone		
CRQL		

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:

Note: Common contaminants such as Methylene chloride, Acetone, 2-Butanone, Carbon disulfide and TICs that were detected in samples within ten times the associated method blank concentration were qualified as not detected, "U". Other contaminants within five times the method blank concentration were also qualified as not detected, "U".

**VALIDATION FINDINGS WORKSHEET**  
**Initial Calibration Calculation Verification**

METHOD: GC/MS VOA (EPA Method 524.2)

The Relative Response Factor (RRF), average RRF, and percent relative standard deviation (%RSD) were recalculated for the compounds identified below using the following calculations:

$RRF = (A_x)(C_s)/(A_s)(C_x)$   
 average RRF = sum of the RRFs/number of standards  
 $\%RSD = 100 * (S/X)$   
 $A_x$  = Area of compound,  
 $C_x$  = Concentration of compound,  
 $S$  = Standard deviation of the RRFs  
 $X$  = Mean of the RRFs  
 $A_s$  = Area of associated internal standard  
 $C_s$  = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference Internal Standard)	Reported		Recalculated		Reported		Recalculated	
				RRF ( / 0 std)	RRF ( / 0 std)	Average RRF (Initial)	Average RRF (Initial)	%RSD	%RSD		
1	1CAL	10/21/03	Methylene chloride (1st Internal Standard)	0.278	0.278	0.271	0.271	4.23	4.18		
			Trichloroethene (2nd internal standard)	0.389	0.389	0.373	0.373	7.19	7.23		
			Bromoform (3rd internal standard)	2.614	2.614	2.542	2.542	4.04	4.04		
2			Methylene chloride (1st Internal Standard)								
			Trichloroethene (2nd internal standard)								
			Bromoform (3rd internal standard)								
3			Methylene chloride (1st Internal Standard)								
			Trichloroethene (2nd internal standard)								
			Bromoform (3rd internal standard)								
4			Methylene chloride (1st Internal Standard)								
			Trichloroethene (2nd internal standard)								
			Bromoform (3rd internal standard)								

Comments: Refer to Initial Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

**VALIDATION FINDINGS WORKSHEET**  
**Continuing Calibration Results Verification**

METHOD: GC/MS VOA (EPA Method 524.2)

The percent difference (%D) of the initial calibration average Relative Response Factors (RRFs) and the continuing calibration RRFs were recalculated for the compounds identified below using the following calculation:

% Difference =  $100 \times (\text{ave. RRF} - \text{RRF}) / \text{ave. RRF}$       Where: ave. RRF = initial calibration average RRF  
 RRF =  $(A_x)(C_s) / (A_s)(C_x)$       RRF = continuing calibration RRF  
 $A_s$  = Area of compound,       $A_x$  = Area of associated internal standard  
 $C_x$  = Concentration of compound,       $C_s$  = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference internal Standard)	Average RRF (Initial)	Reported		Recalculated	
					RRF (CC)	%D	RRF (CC)	%D
1	94668801	11/1/03	Methylene-chloride (1st Internal Standard)	0.271	0.266	1.8	0.266	1.7
			Trichlorethene (2nd internal standard)	0.383	0.384	2.9	0.384	3.0
			Bromoform (3rd internal standard)	2.42	2.421	4.8	2.421	4.8
2			Methylene chloride (1st Internal Standard)					
			Trichlorethene (2nd internal standard)					
			Bromoform (3rd internal standard)					
3			Methylene chloride (1st Internal Standard)					
			Trichlorethene (2nd internal standard)					
			Bromoform (3rd internal standard)					
4			Methylene chloride (1st Internal Standard)					
			Trichlorethene (2nd internal standard)					
			Bromoform (3rd internal standard)					

Comments: Refer to Continuing Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 1127001  
 SDG #: 03-5852

**VALIDATION FINDINGS WORKSHEET**  
**Surrogate Results Verification**

Page: 1 of 1  
 Reviewer: [Signature]  
 2nd reviewer: [Signature]

**METHOD: GC/MS VOA (EPA Method 524.2)**

The percent recoveries (%R) of surrogates were recalculated for the compounds identified below using the following calculation:

% Recovery: SF/SS \* 100

Where: SF = Surrogate Found  
 SS = Surrogate Spiked

Sample ID: 1

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8	20	20.97	105	105	0
Bromofluorobenzene	↓	19.49	98	98	↓
1,2-Dichlorobenzene-d4	↓	19.38	97	97	↓

Sample ID: \_\_\_\_\_

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					

Sample ID: \_\_\_\_\_

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					

Sample ID: \_\_\_\_\_

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					

Sample ID: \_\_\_\_\_

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					

LDC #: 1127001  
 SDG #: 035852

**VALIDATION FINDINGS WORKSHEET**  
**Laboratory Control Sample Results Verification**

Page: 1 of 1  
 Reviewer: [Signature]  
 2nd Reviewer: [Signature]

**METHOD: GC/MS VOA (EPA Method 524.2)**

The percent recoveries (%R) and Relative Percent Difference (RPD) of the laboratory control sample and laboratory control sample duplicate (if applicable) were recalculated for the compounds identified below using the following calculation:

% Recovery = 100 \* SSC/SA

Where: SSC = Spiked sample concentration  
 SA = Spike added

RPD = |LCS - LCSD| \* 2 / (LCS + LCSD)

LCS = Laboratory control sample percent recovery

LCSD = Laboratory control sample duplicate percent recovery

LCS ID: GA668101

Compound	Spike Added ( <u>NA</u> )		Spiked Sample Concentration ( <u>NA</u> )		LCS		LCSD		Percent Recovery		RPD	
	LCS	LCSD	LCS	LCSD	Reported	Recalc.	Reported	Recalc.	Reported	Recalc.	Reported	Recalculated
1,1-Dichloroethene	20	NA	19.3	NA	97	97	97	97				
Trichloroethene			19.1		96	96	96	96				
Benzene			17.9		90	90	90	90				
Toluene			17.7		89	89	89	89				
Chlorobenzene			18.4		92	92	92	92				

Comments: Refer to Laboratory Control Sample findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.



LDC #: 11272D1  
SDG #: 035852

### VALIDATION FINDINGS WORKSHEET

**Sample Calculation Verification**

Page: 1 of 1  
Reviewer:                           
2nd reviewer:                         

**METHOD:** GC/MS VOA (EPA Method 524.2)

Y  N  N/A Were all reported results recalculated and verified for all level IV samples?

Y  N  N/A Were all recalculated results for detected target compounds agree within 10.0% of the reported results?

$$\text{Concentration} = \frac{(A_x)(I_s)(DF)}{(A_{is})(RRF)(V_s)(\%S)}$$

- A<sub>x</sub> = Area of the characteristic ion (EICP) for the compound to be measured
- A<sub>is</sub> = Area of the characteristic ion (EICP) for the specific internal standard
- I<sub>s</sub> = Amount of internal standard added in nanograms (ng)
  
- RRF = Relative response factor of the calibration standard.
- V<sub>s</sub> = Volume or weight of sample pruged in milliliters (ml) or grams (g).
- Df = Dilution factor.
- %S = Percent solids, applicable to soils and solid matrices only.

Example:

Sample I.D.           /          , ND :

Conc. = (            ) (            ) (            )  
          (            ) (            ) (            ) (            )  
          =

#	Sample ID	Compound	Reported Concentration (            )	Calculated Concentration (            )	Qualification

**NASA JPL  
Data Validation Reports  
LDC# 11272**

Chromium

**LDC**

## Laboratory Data Consultants, Inc. Data Validation Report

**Project/Site Name:** NASA JPL  
**Collection Date:** October 22 through October 29, 2003  
**LDC Report Date:** December 18, 2003  
**Matrix:** Water  
**Parameters:** Chromium  
**Validation Level:** EPA Level III & IV  
**Laboratory:** Advanced Technology Laboratories &  
Applied P & Ch Laboratory

**Sample Delivery Group (SDG):** 03-5876/065632

### Sample Identification

MW-20-5	MW-4-3	MW-19-3
MW-20-4	MW-4-2	MW-19-2
MW-20-3	MW-4-1	MW-19-1
MW-20-2	Dupe-3-4-Q03**	MW-4-3MS
MW-20-1	EB-4-10-27-03	MW-4-3MSD
EB-3-10-24-03	MW-24-5	MW-4-3DUP
Dupe-6-4-Q03**	MW-24-4	MW-11-3DUP
MW-11-5	Dupe-1-4-Q03**	MW-18-2MS
MW-11-4	MW-24-3	MW-18-2MSD
MW-11-3	MW-24-2	MW-18-2DUP
MW-11-2	MW-24-1	MW-19-1MS
MW-11-1	EB-5-10-28-03	MW-19-1MSD
EB-1-10-22-03	MW-18-5	MW-19-1DUP
MW-22-5	MW-18-4	MW-19-3MS
MW-22-4	MW-18-3	MW-19-3MSD
MW-22-3	MW-18-2	MW-19-3DUP
MW-22-2	MW-18-1	MW-24-2DUP
MW-22-1	EB-6-10-29-03	
MW-4-5	MW-19-5	
MW-4-4	MW-19-4	

\*\*Indicates sample underwent EPA Level IV review

## Introduction

This data review covers 57 water samples listed on the cover sheet including dilutions and reanalysis as applicable. The analyses were per EPA Method 200.8 for Chromium.

The review follows a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines for Inorganic Data Review (February 1994) as there are no current guidelines for the method stated above.

A table summarizing all data qualification flags is provided at the end of this report. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from specified protocols or is of technical advisory nature.

Blanks are summarized in Section III.

Field duplicates are summarized in Section XIII.

Samples indicated by a double asterisk on the front cover underwent a EPA Level IV review. A EPA Level III review was performed on all of the other samples. Raw data were not evaluated for the samples reviewed by Level III criteria since this review is based on QC data.

The following are definitions of the data qualifiers:

- U Indicates the compound or analyte was analyzed for but not detected at or above the stated limit.
- J Indicates an estimated value.
- R Quality control indicates the data is not usable.
- N Presumptive evidence of presence of the constituent.
- UJ Indicates the compound or analyte was analyzed for but not detected. The sample detection limit is an estimated value.
- A Indicates the finding is based upon technical validation criteria.
- P Indicates the finding is related to a protocol/contractual deviation.
- None Indicates the data was not significantly impacted by the finding, therefore qualification was not required.

## I. Technical Holding Times

All technical holding time requirements were met.

The chain-of-custodies were reviewed for documentation of cooler temperatures. All cooler temperatures met validation criteria.

## II. Calibration

An initial calibration was performed.

The frequency and analysis criteria of the initial calibration verification (ICV) and continuing calibration verification (CCV) were met.

CRDL standards for ICP and AA were analyzed and reported as required.

## III. Blanks

Method blanks were reviewed for each matrix as applicable.

Data qualification by the initial, continuing and preparation blanks (ICB/CCB/PBs) was based on the maximum contaminant concentration in the ICB/CCB/PBs in the analysis of each analyte. No contaminant concentrations were found in the initial, continuing and preparation blanks with the following exceptions:

Method Blank ID	Analyte	Maximum Concentration	Associated Samples
ICB/CCB	Chromium	0.284 ug/L	MW-20-5 MW-20-4 MW-20-3 MW-20-2 MW-20-1 EB-3-10-24-03 Dupe-6-4-Q03** MW-11-5 MW-11-4 MW-11-3 MW-11-2 MW-11-1 EB-1-10-22-03 MW-22-5 MW-22-4 MW-22-3 MW-22-2 MW-22-1 MW-4-5 MW-19-3
PB (prep blank)	Chromium	0.461 ug/L	MW-4-4 MW-19-2 MW-19-1

Method Blank ID	Analyte	Maximum Concentration	Associated Samples
ICB/CCB	Chromium	0.442 ug/L	MW-4-4 MW-4-2 MW-4-1 Dupe-3-4-Q03** EB-4-10-27-03 MW-24-5 MW-24-4 Dupe-1-4-Q03** MW-24-3 MW-24-2 MW-24-1 EB-5-10-28-03 MW-18-5 MW-18-4 MW-18-3 MW-18-2 MW-18-1 EB-6-10-29-03 MW-19-5 MW-19-4 MW-19-2 MW-19-1
PB (prep blank)	Chromium	0.1825 ug/L	MW-4-3
ICB/CCB	Chromium	0.400 ug/L	MW-4-3

Sample concentrations were compared to the maximum contaminant concentrations detected in the ICB/CCB/PBs. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated method blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-20-5	Chromium	1.3 ug/L	1.3U ug/L
MW-20-2	Chromium	1.3 ug/L	1.3U ug/L
Dupe-6-4-Q03**	Chromium	1.4 ug/L	1.4U ug/L
MW-11-4	Chromium	0.8 ug/L	0.8U ug/L
MW-11-2	Chromium	1 ug/L	1U ug/L
MW-22-5	Chromium	0.7 ug/L	0.7U ug/L
MW-22-2	Chromium	0.9 ug/L	0.9U ug/L

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-20-4	Chromium	1.6 ug/L	1.6U ug/L
MW-11-5	Chromium	1.5 ug/L	1.5U ug/L
MW-24-4	Chromium	1.2 ug/L	1.2U ug/L
Dupe-1-4-Q03**	Chromium	1.1 ug/L	1.1U ug/L
MW-24-3	Chromium	1.7 ug/L	1.7U ug/L
MW-18-5	Chromium	1.0 ug/L	1.0U ug/L
MW-18-2	Chromium	1.9 ug/L	1.9U ug/L
MW-18-1	Chromium	1.5 ug/L	1.5U ug/L
MW-19-5	Chromium	1.8 ug/L	1.8U ug/L
MW-19-1	Chromium	1.2 ug/L	1.2U ug/L
MW-20-4	Chromium	2.4 ug/L	2.4U ug/L
MW-4-1	Chromium	2.6 ug/L	2.6U ug/L
MW-24-2	Chromium	2.7 ug/L	2.7U ug/L
MW-18-4	Chromium	2.6 ug/L	2.6U ug/L
MW-19-4	Chromium	2.4 ug/L	2.4U ug/L
MW-4-3	Chromium	0.4 ug/L	0.4U ug/L

#### IV. ICP Interference Check Sample (ICS) Analysis

The frequency of analysis was met.

The criteria for analysis were met.

## V. Matrix Spike Analysis

Matrix spike (MS) and matrix spike duplicate (MSD) samples were reviewed for each matrix as applicable. Percent recoveries (%R) and relative percent differences (RPD) were within QC limits with the following exceptions:

Spike ID (Associated Samples)	Analyte	MS (%R) (Limits)	MSD (%R) (Limits)	RPD (Limits)	Flag	A or P
MW-19-3MS/MSD (MW-20-5 MW-20-4 MW-20-3 MW-20-2 MW-20-1 EB-3-10-24-03 Dupe-6-4-Q03** MW-11-5 MW-11-4 MW-11-3 MW-11-2 MW-11-1 EB-1-10-22-03 MW-22-5 MW-22-4 MW-22-3 MW-22-2 MW-22-1 MW-4-5 MW-19-3)	Chromium	68.4 (80-120)	72.8 (80-120)	-	J (all detects) UJ (all non-detects)	A

## VI. Duplicate Sample Analysis

Duplicate (DUP) sample analyses were reviewed for each matrix as applicable. Results were within QC limits.

## VII. Laboratory Control Samples (LCS)

Laboratory control samples were reviewed for each matrix as applicable. Percent recoveries (%R) were within QC limits.

## VIII. Internal Standards

All internal standard percent recoveries (%R) were within QC limits for samples on which a EPA Level IV review was performed. Raw data were not evaluated for the samples reviewed by Level III criteria.

## IX. Furnace Atomic Absorption QC

Graphite furnace atomic absorption was not utilized in this SDG.



## X. ICP Serial Dilution

ICP serial dilution was not required by the method.

## XI. Sample Result Verification

All sample result verifications met validation criteria for samples on which a EPA Level IV review was performed. Raw data were not evaluated for the samples reviewed by Level III criteria.

## XII. Overall Assessment of Data

Data flags have been summarized at the end of this report.

## XIII. Field Duplicates

Samples MW-20-2 and Dupe-6-4-Q03\*\*, samples MW-4-5 and Dupe-3-4-Q03\*\*, and samples MW-24-4 and Dupe-1-4-Q03\*\* were identified as field duplicates. No chromium was detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-20-2	Dupe-6-4-Q03**	
Chromium	1.3	1.4	7

Analyte	Concentration (ug/L)		RPD
	MW-4-5	Dupe-3-4-Q03**	
Chromium	3.5	5.6	46

Analyte	Concentration (ug/L)		RPD
	MW-24-4	Dupe-1-4-Q03**	
Chromium	1.2	1.1	9

## XIV. Field Blanks

Samples EB-3-10-24-03, EB-1-10-22-03, EB-4-10-27-03, EB-5-10-28-03, and EB-6-10-29-03 were identified as equipment blanks. No chromium was found in these blanks.

**NASA JPL**

**Chromium - Data Qualification Summary - SDG 03-5876/065632**

SDG	Sample	Analyte	Flag	A or P	Reason
03-5876/ 065632	MW-20-5 MW-20-4 MW-20-3 MW-20-2 MW-20-1 EB-3-10-24-03 Dupe-6-4-Q03** MW-11-5 MW-11-4 MW-11-3 MW-11-2 MW-11-1 EB-1-10-22-03 MW-22-5 MW-22-4 MW-22-3 MW-22-2 MW-22-1 MW-4-5 MW-19-3	Chromium	J (all detects) UJ (all non-detects)	A	Matrix spike/Matrix spike duplicates (%R)

**NASA JPL**

**Chromium - Laboratory Blank Data Qualification Summary - SDG 03-5876/065632**

SDG	Sample	Analyte	Modified Final Concentration	A or P
03-5876/ 065632	MW-20-5	Chromium	1.3U ug/L	A
03-5876/ 065632	MW-20-2	Chromium	1.3U ug/L	A
03-5876/ 065632	Dupe-6-4-Q03**	Chromium	1.4U ug/L	A
03-5876/ 065632	MW-11-4	Chromium	0.8U ug/L	A
03-5876/ 065632	MW-11-2	Chromium	1U ug/L	A
03-5876/ 065632	MW-22-5	Chromium	0.7U ug/L	A
03-5876/ 065632	MW-22-2	Chromium	0.9U ug/L	A

SDG	Sample	Analyte	Modified Final Concentration	A or P
03-5876/ 065632	MW-20-4	Chromium	1.6U ug/L	A
03-5876/ 065632	MW-11-5	Chromium	1.5U ug/L	A
03-5876/ 065632	MW-24-4	Chromium	1.2U ug/L	A
03-5876/ 065632	Dupe-1-4-Q03**	Chromium	1.1U ug/L	A
03-5876/ 065632	MW-24-3	Chromium	1.7U ug/L	A
03-5876/ 065632	MW-18-5	Chromium	1.0U ug/L	A
03-5876/ 065632	MW-18-2	Chromium	1.9U ug/L	A
03-5876/ 065632	MW-18-1	Chromium	1.5U ug/L	A
03-5876/ 065632	MW-19-5	Chromium	1.8U ug/L	A
03-5876/ 065632	MW-19-1	Chromium	1.2U ug/L	A
03-5876/ 065632	MW-20-4	Chromium	2.4U ug/L	A
03-5876/ 065632	MW-4-1	Chromium	2.6U ug/L	A
03-5876/ 065632	MW-24-2	Chromium	2.7U ug/L	A
03-5876/ 065632	MW-18-4	Chromium	2.6U ug/L	A
03-5876/ 065632	MW-19-4	Chromium	2.4U ug/L	A

SDG	Sample	Analyte	Modified Final Concentration	A or P
03-5876/ 065632	MW-4-3	Chromium	0.4U ug/L	A

**NASA JPL  
Chromium - Field Blank Data Qualification Summary - SDG 03-5876/065632**

No Sample Data Qualified in this SDG

E

# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-001A

**Client Sample ID:** MW-20-5  
**Collection Date:** 10/24/2003 7:33:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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### ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4_031104B	QC Batch: R32245			PrepDate:			
Chromium	1.3	UT	0.11	1.0	µg/L	1	11/4/2003 3:18:39 PM

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**Qualifiers:**

ND - Not Detected at the Reporting Limit	S - Spike Recovery outside accepted recovery limits
J - Analyte detected below quantitation limits	R - RPD outside accepted recovery limits
B - Analyte detected in the associated Method Blank	E - Value above quantitation range
* - Value exceeds Maximum Contaminant Level	H-Samples exceed holding time

Results are wet unless otherwise specified

0010



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-002A

**Client Sample ID:** MW-20-4  
**Collection Date:** 10/24/2003 8:00:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

1.6

✓

0.11

1.0 µg/L

1

11/4/2003 3:21:08 PM

### Qualifiers:

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

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0011



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-003A

**Client Sample ID:** MW-20-3  
**Collection Date:** 10/24/2003 8:49:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

2.9



0.11

1.0 µg/L

1

11/4/2003 3:23:38 PM

*Handwritten signature*

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

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Results are wet unless otherwise specified

0012



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-004A

**Client Sample ID:** MW-20-2  
**Collection Date:** 10/24/2003 9:11:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	1.3	UJ	0.11	1.0	µg/L	1	11/4/2003 3:26:08 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

*Handwritten signature/initials*





# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-005A

**Client Sample ID:** MW-20-1  
**Collection Date:** 10/24/2003 9:49:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	1.9	J	0.11	1.0	µg/L	1	11/4/2003 3:28:38 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H - Samples exceed holding time

Results are wet unless otherwise specified

*Handwritten signature*



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-006A

**Client Sample ID:** EB-3-10-24-03  
**Collection Date:** 10/24/2003 8:41:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>		<b>EPA 200.8</b>			<b>Analyst: NS</b>		
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	ND	UJ	0.11	1.0	µg/L	1	11/4/2003 3:31:09 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

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0015



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-007A

**Client Sample ID:** Dupe-6-4-Q03  
**Collection Date:** 10/24/2003 9:29:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	1.4	UT	0.11	1.0	µg/L	1	11/4/2003 3:33:41 PM

*WDR*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

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Results are wet unless otherwise specified

0016



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-008A

**Client Sample ID:** MW-11-5  
**Collection Date:** 10/22/2003 7:48:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>				<b>Analyst: NS</b>
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	1.5	J	0.11	1.0	µg/L	1	11/4/2003 3:36:12 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

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0017



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-009A

**Client Sample ID:** MW-11-4  
**Collection Date:** 10/22/2003 8:10:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031104B	QC Batch: R32245			PrepDate:			
Chromium	0.8	J <i>WJ</i>	0.11	1.0	µg/L	1	11/4/2003 3:43:54 PM

*11/10/03*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0018



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-010A

**Client Sample ID:** MW-11-3  
**Collection Date:** 10/22/2003 8:47:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	3.4	J	0.11	1.0	µg/L	1	11/4/2003 3:46:27 PM

*11/18/03*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

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Results are wet unless otherwise specified

0019



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-011A

**Client Sample ID:** MW-11-2  
**Collection Date:** 10/22/2003 9:10:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>				<b>Analyst: NS</b>
RunID: ICP4_031104B	QC Batch: R32245			PrepDate:			
Chromium	1	J UJ	0.22	2.0	µg/L	2	11/4/2003 4:38:59 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

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**Advanced Technology Laboratories**

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-012A

**Client Sample ID:** MW-11-1  
**Collection Date:** 10/22/2003 9:32:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>		<b>EPA 200.8</b>		<b>Analyst: NS</b>			
RunID: ICP4_031104B	QC Batch: R32245			PrepDate:			
Chromium	2	J	0.22	2.0	µg/L	2	11/4/2003 4:41:32 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified





# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-013A

**Client Sample ID:** EB-1-10-22-03  
**Collection Date:** 10/22/2003 9:00:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	ND	UJ	0.11	1.0	µg/L	1	11/4/2003 3:56:39 PM

**Qualifiers:**  
ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Page 1 of 1

Results are wet unless otherwise specified

*Handwritten signature*



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-014A

**Client Sample ID:** MW-22-5  
**Collection Date:** 10/22/2003 10:21:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

0.7

J

*WJ*

0.11

1.0 µg/L

1

11/4/2003 3:59:06 PM

### Qualifiers:

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H-Samples exceed holding time

Page 1 of 1

Results are wet unless otherwise specified

*NOTE*



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-015A

**Client Sample ID:** MW-22-4  
**Collection Date:** 10/22/2003 10:41:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031104B	QC Batch: R32245			PrepDate:			
Chromium	3.1	J	0.22	2.0	µg/L	2	11/4/2003 4:44:05 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0024



**Advanced Technology Laboratories**

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-016A

**Client Sample ID:** MW-22-3  
**Collection Date:** 10/22/2003 11:00:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031104B	QC Batch: R32245		PrepDate:				
Chromium	3.2	J	0.22	2.0	µg/L	2	11/4/2003 4:46:39 PM

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0025



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-017A

**Client Sample ID:** MW-22-2  
**Collection Date:** 10/22/2003 11:20:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

0.9

J

*WJ*

0.56

5.0 µg/L

5

11/4/2003 5:09:44 PM

### Qualifiers:

ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H-Samples exceed holding time

Results are wet unless otherwise specified

Page 1 of 1

*WJ*



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-018A

**Client Sample ID:** MW-22-1  
**Collection Date:** 10/22/2003 11:40:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

3

J

J

0.56

5.0 µg/L

5

11/4/2003 5:12:14 PM

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

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Results are wet unless otherwise specified

0027



# Advanced Technology Laboratories

Date: 17-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-019A

**Client Sample ID:** MW-4-5  
**Collection Date:** 10/27/2003 7:44:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031104B

QC Batch: R32245

PrepDate:

Chromium

3.5

J

0.11

1.0 µg/L

1 11/4/2003 4:16:39 PM

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

Page 1 of 1

Results are wet unless otherwise specified

0028



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-020A

**Client Sample ID:** MW-4-4  
**Collection Date:** 10/27/2003 8:39:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105C	QC Batch: R32300			PrepDate:			
Chromium	2.4		0.11	1.0	µg/L	1	11/5/2003 11:58:43 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

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0029





# Advanced Technology Laboratories

Date: 10-Nov-03

CLIENT: Applied P & Ch Laboratories  
Lab Order: 065632  
Project: JPL GW Mon, 04-4428.10  
Lab ID: 065632-021A

Client Sample ID: MW-4-3  
Collection Date: 10/27/2003 9:01:00 AM  
Matrix: WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>	<b>(EPA 200.8)</b>			<b>Analyst: NS</b>
RunID: ICP4_031105D	QC Batch: 15669			PrepDate: 11/5/2003			
Chromium	0.4	J U	0.11	1.0	µg/L	1	11/5/2003 2:28:26 PM

Qualifiers: ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-022A

**Client Sample ID:** MW-4-2  
**Collection Date:** 10/27/2003 9:20:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	3.7		0.11	1.0	µg/L	1	11/5/2003 10:38:18 A

*Handwritten signature*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0031



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-023A

**Client Sample ID:** MW-4-1  
**Collection Date:** 10/27/2003 9:46:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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**ICP-MS METALS**

**EPA 200.8**

**Analyst: NS**

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

2.6

0.11

1.0 µg/L

1

11/5/2003 10:40:47 A

**Qualifiers:**

ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

*Handwritten signature*



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-024A

**Client Sample ID:** Dupe-3-4-Q03  
**Collection Date:** 10/27/2003 8:12:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031105B	QC Batch: R32298			PrepDate:			
Chromium	5.6		0.11	1.0	µg/L	1	11/5/2003 10:43:17 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

*Handwritten signature*



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-025A

**Client Sample ID:** EB-4-10-27-03  
**Collection Date:** 10/27/2003 9:59:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298			PrepDate:			
Chromium	ND		0.11	1.0	µg/L	1	11/5/2003 10:45:47 A

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

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# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-026A

**Client Sample ID:** MW-24-5  
**Collection Date:** 10/28/2003 7:58:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	3.7		0.11	1.0	µg/L	1	11/5/2003 10:48:17 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

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0035



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-027A

**Client Sample ID:** MW-24-4  
**Collection Date:** 10/28/2003 8:29:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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**ICP-MS METALS**

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.2

U

0.11

1.0 µg/L

1

11/5/2003 10:50:48 A

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

*Handwritten signature*



Advanced Technology Laboratories

Date: 10-Nov-03

CLIENT: Applied P & Ch Laboratories  
Lab Order: 065632  
Project: JPL GW Mon, 04-4428.10  
Lab ID: 065632-028A

Client Sample ID: Dupe-1-4Q03  
Collection Date: 10/28/2003 9:21:00 AM  
Matrix: WATER

Analyte Result Qual MDL PQL Units DF Date Analyzed

ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.1 U

0.11

1.0 µg/L

1

11/5/2003 10:53:19 A

Qualifiers: ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

0037



*Handwritten signature*



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-029A

**Client Sample ID:** MW-24-3  
**Collection Date:** 10/28/2003 9:46:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.7

U

0.11

1.0 µg/L

1

11/5/2003 10:55:51 A

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

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Results are wet unless otherwise specified

0038



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-030A

**Client Sample ID:** MW-24-2  
**Collection Date:** 10/28/2003 10:11:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	2.7	U	0.11	1.0	µg/L	1	11/5/2003 11:03:33 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

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11/2/03



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-031A

**Client Sample ID:** MW-24-1  
**Collection Date:** 10/28/2003 10:41:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	4.0		0.11	1.0	µg/L	1	11/5/2003 11:08:39 A

*Handwritten signature*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0040



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-032A

**Client Sample ID:** EB-5-10-28-03  
**Collection Date:** 10/28/2003 10:30:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	ND		0.11	1.0	µg/L	1	11/5/2003 11:11:13 A

*WJL*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0041



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-033A

**Client Sample ID:** MW-18-5  
**Collection Date:** 10/29/2003 7:46:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.0

U

0.11

1.0 µg/L

1

11/5/2003 11:13:47 A

**Qualifiers:** ND - Not Detected at the Reporting Limit

J - Analyte detected below quantitation limits

B - Analyte detected in the associated Method Blank

\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits

R - RPD outside accepted recovery limits

E - Value above quantitation range

H-Samples exceed holding time

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Results are wet unless otherwise specified



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-034A

**Client Sample ID:** MW-18-4  
**Collection Date:** 10/29/2003 8:14:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	2.6	U	0.11	1.0	µg/L	1	11/5/2003 11:16:18 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified



*M 12/10*

# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-035A

**Client Sample ID:** MW-18-3  
**Collection Date:** 10/29/2003 8:41:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105B	QC Batch: R32298			PrepDate:			
Chromium	5.9		0.11	1.0	µg/L	1	11/5/2003 11:18:45 A

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

*Handwritten signature*

Results are wet unless otherwise specified

0044



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-036A

**Client Sample ID:** MW-18-2  
**Collection Date:** 10/29/2003 9:06:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031105B	QC Batch: R32298		PrepDate:				
Chromium	1.9	U	0.11	1.0	µg/L	1	11/5/2003 11:21:13 A

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level  
S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

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*MW-18-2*

0045





# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-037A

**Client Sample ID:** MW-18-1  
**Collection Date:** 10/29/2003 9:55:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.5

U

0.11

1.0 µg/L

1

11/5/2003 11:26:10 A

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

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Results are wet unless otherwise specified

0046



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-038A

**Client Sample ID:** EB-6-10-29-03  
**Collection Date:** 10/29/2003 9:45:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>		<b>Analyst: NS</b>		
RunID: ICP4_031105B	QC Batch: R32298			PrepDate:			
Chromium	ND		0.11	1.0	µg/L	1	11/5/2003 11:33:48 A

*W/2K*

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
\* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified

0047



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-039A

**Client Sample ID:** MW-19-5  
**Collection Date:** 10/29/2003 11:17:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

1.8

U

0.11

1.0 µg/L

1

11/5/2003 11:36:18 A

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

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Results are wet unless otherwise specified

*Handwritten signature*



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-040A

**Client Sample ID:** MW-19-4  
**Collection Date:** 10/29/2003 11:40:00 AM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
---------	--------	------	-----	-----	-------	----	---------------

## ICP-MS METALS

EPA 200.8

Analyst: NS

RunID: ICP4\_031105B

QC Batch: R32298

PrepDate:

Chromium

2.4

U

0.11

1.0 µg/L

1

11/5/2003 11:38:48 A

### Qualifiers:

ND - Not Detected at the Reporting Limit

S - Spike Recovery outside accepted recovery limits

J - Analyte detected below quantitation limits

R - RPD outside accepted recovery limits

B - Analyte detected in the associated Method Blank

E - Value above quantitation range

\* - Value exceeds Maximum Contaminant Level

H-Samples exceed holding time

Results are wet unless otherwise specified

Page 40 of 43

*Handwritten signature*



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-041A

**Client Sample ID:** MW-19-3  
**Collection Date:** 10/29/2003 12:10:00 PM  
**Matrix:** WATER

**Analyte**                      **Result**   **Qual**      **MDL**      **PQL**   **Units**                      **DF**   **Date Analyzed**

**ICP-MS METALS**

**EPA 200.8**

**Analyst: NS**

RunID: ICP4\_031103A

QC Batch: R32245

PrepDate:

Chromium

4.3

J

0.22

2.0 µg/L

2

11/3/2003 4:54:04 PM

*Handwritten signature*

**Qualifiers:**      ND - Not Detected at the Reporting Limit                      S - Spike Recovery outside accepted recovery limits  
                         J - Analyte detected below quantitation limits                      R - RPD outside accepted recovery limits  
                         B - Analyte detected in the associated Method Blank                      E - Value above quantitation range  
                         \* - Value exceeds Maximum Contaminant Level                      H-Samples exceed holding time

Results are wet unless otherwise specified



# Advanced Technology Laboratories

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-042A

**Client Sample ID:** MW-19-2  
**Collection Date:** 10/29/2003 12:50:00 PM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
<b>ICP-MS METALS</b>			<b>EPA 200.8</b>			<b>Analyst: NS</b>	
RunID: ICP4_031105C	QC Batch: R32300			PrepDate:			
Chromium	4.0		0.11	1.0	µg/L	1	11/5/2003 12:01:16 P

*Handwritten:* 11/2/03

**Qualifiers:** ND - Not Detected at the Reporting Limit      S - Spike Recovery outside accepted recovery limits  
 J - Analyte detected below quantitation limits      R - RPD outside accepted recovery limits  
 B - Analyte detected in the associated Method Blank      E - Value above quantitation range  
 \* - Value exceeds Maximum Contaminant Level      H-Samples exceed holding time

Results are wet unless otherwise specified.

0051



**Advanced Technology Laboratories**

Date: 10-Nov-03

**CLIENT:** Applied P & Ch Laboratories  
**Lab Order:** 065632  
**Project:** JPL GW Mon, 04-4428.10  
**Lab ID:** 065632-043A

**Client Sample ID:** MW-19-1  
**Collection Date:** 10/29/2003 1:13:00 PM  
**Matrix:** WATER

Analyte	Result	Qual	MDL	PQL	Units	DF	Date Analyzed
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**ICP-MS METALS**

**EPA 200.8**

**Analyst: NS**

RunID: ICP4\_031105C

QC Batch: R32300

PrepDate:

Chromium

1.2

U

0.11

1.0 µg/L

1

11/5/2003 12:03:50 P

*1/2/04*

**Qualifiers:** ND - Not Detected at the Reporting Limit  
J - Analyte detected below quantitation limits  
B - Analyte detected in the associated Method Blank  
\* - Value exceeds Maximum Contaminant Level

S - Spike Recovery outside accepted recovery limits  
R - RPD outside accepted recovery limits  
E - Value above quantitation range  
H-Samples exceed holding time

Results are wet unless otherwise specified

0052



LDC #: 11272E4

**VALIDATION COMPLETENESS WORKSHEET**

Date: 12/15/03

SDG #: 03-5876/065632

Level III/IV

Page: 1 of 2

Laboratory: Advanced Technology Laboratories/Applied P & Ch Laboratory

Reviewer: MW

2nd Reviewer: [Signature]

**METHOD:** Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times	A	Sampling dates: 10/22 - 29/03
II.	Calibration	A	
III.	Blanks	SW	
IV.	ICP Interference Check Sample (ICS) Analysis	N	Not required
V.	Matrix Spike Analysis	SW	3 MS/MSD/MSR
VI.	Duplicate Sample Analysis	A	
VII.	Laboratory Control Samples (LCS)	A	LCS
VIII.	Internal Standard (ICP-MS)	A	
IX.	Furnace Atomic Absorption QC	N	Not Utilized
X.	ICP Serial Dilution	N	Not required
XI.	Sample Result Verification	A	Not reviewed for Level III validation.
XII.	Overall Assessment of Data	A	
XIII.	Field Duplicates	SW	(4,7), (19,24), (29,28)
XIV.	Field Blanks	MB	EB=6, 13, 25, 32, 38.

Note: A = Acceptable      ND = No compounds detected      D = Duplicate  
 N = Not provided/applicable      R = Rinsate      TB = Trip blank  
 SW = See worksheet      FB = Field blank      EB = Equipment blank

Validated Samples: \*\* Indicates sample underwent Level IV validation

1	MW-20-5	11	MW-11-2	21	MW-4-3	31	MW-24-1
2	MW-20-4	12	MW-11-1	22	MW-4-2	32	EB-5-10-28-03
3	MW-20-3	13	EB-1-10-22-03	23	MW-4-1	33	MW-18-5
4	MW-20-2	14	MW-22-5	24	Dupe-3-4-Q03**	34	MW-18-4
5	MW-20-1	15	MW-22-4	25	EB-4-10-27-03	35	MW-18-3
6	EB-3-10-24-03	16	MW-22-3	26	MW-24-5	36	MW-18-2
7	Dupe-6-4-Q03**	17	MW-22-2	27	MW-24-4	37	MW-18-1
8	MW-11-5	18	MW-22-1	28	Dupe-1-4-Q03**	38	EB-6-10-29-03
9	MW-11-4	19	MW-4-5	29	MW-24-3	39	MW-19-5
10	MW-11-3	20	MW-4-4	30	MW-24-2	40	MW-19-4

Notes: \_\_\_\_\_  
 \_\_\_\_\_  
 \_\_\_\_\_



LDC #: 11272E4

### VALIDATION COMPLETENESS WORKSHEET

Date: 12/15/03

SDG #: 03-5876/065632

Level III/IV

Page: 2 of 2

Laboratory: Advanced Technology Laboratories/Applied P & Ch Laboratory

Reviewer: MW

2nd Reviewer: [Signature]

METHOD: Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Technical holding times		Sampling dates:
II.	Calibration		
III.	Blanks		
IV.	ICP Interference Check Sample (ICS) Analysis		
V.	Matrix Spike Analysis	See page 1	
VI.	Duplicate Sample Analysis		
VII.	Laboratory Control Samples (LCS)		
VIII.	Internal Standard (ICP-MS)		
IX.	Furnace Atomic Absorption QC		
X.	ICP Serial Dilution		
XI.	Sample Result Verification		Not reviewed for Level III validation.
XII.	Overall Assessment of Data		
XIII.	Field Duplicates		
XIV.	Field Blanks		

Note: A = Acceptable  
N = Not provided/applicable  
SW = See worksheet

ND = No compounds detected  
R = Rinsate  
FB = Field blank

D = Duplicate  
TB = Trip blank  
EB = Equipment blank

Validated Samples: \*\* Indicates sample underwent Level IV validation

41	MW-19-3	51	MW-19-1MS				
42	MW-19-2	52	MW-19-1MSD				
43	MW-19-1	53	MW-19-1DUP				
44	MW-4-3MS	54	MW-19-3MS				
45	MW-4-3MSD	55	MW-19-3MSD				
46	MW-4-3DUP	56	MW-19-3DUP				
47	MW-11-3DUP	57	MW-24-2DUP				
48	MW-18-2MS	58	PB				
49	MW-18-2MSD	59					
50	MW-18-2DUP	60					

Notes: \_\_\_\_\_  
\_\_\_\_\_  
\_\_\_\_\_

LDC #: 11272 E4  
 SDG #: See color

VALIDATION FINDINGS CHECKLIST

Page: 1 of 1  
 Reviewer: WJ  
 2nd Reviewer: [Signature]

Method: Metals (EPA SW 826 Method 6010/7000/6020)

Validation Area	Yes	No	NA	Findings/Comments
<b>I. Technical holding times</b>				
All technical holding times were met.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Cooler temperature criteria was met.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>II. Calibration</b>				
Were all instruments calibrated daily, each set-up time?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the proper number of standards used?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all initial and continuing calibration verification %Rs within the 90-110% (80-120% for mercury and 85-115% for cyanide) QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all initial calibration correlation coefficients $\geq 0.995$ ?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was a midrange cyanide standard distilled?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>III. Blanks</b>				
Was a method blank associated with every sample in this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>IV. ICP Interference Check Sample</b>				
Were ICP interference check samples performed daily?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
Were the AB solution percent recoveries (%R) with the 80-120% QC limits?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
<b>IV. Matrix spike/Matrix spike duplicates</b>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq 20\%$ for waters and $\leq 35\%$ for soil samples? A control limit of $\pm 1 \text{ RL}$ ( $\pm 2 \text{ RL}$ for soil) was used for samples that were $\leq 5 \text{ X}$ the RL, including when only one of the duplicate sample values were $\leq 5 \text{ X}$ the RL.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>V. Laboratory control samples</b>				
Was an LCS analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was an LCS analyzed per extraction batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% QC limits for water samples and laboratory established QC limits for soils?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
<b>VI. Furnace Atomic Absorption QC</b>				
If MSA was performed, was the correlation coefficients $\geq 0.995$ ?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
Do all applicable analyses have duplicate injections?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	

LDC #: 1121284  
 SDG #: see cover

VALIDATION FINDINGS CHECKLIST

Page: 1 of 1  
 Reviewer: VJZ  
 2nd Reviewer: [Signature]

Validation Area	Yes	No	NA	Findings/Comments
For sample concentrations > RL, are applicable duplicate injection RSD values < 20%?			<input checked="" type="checkbox"/>	
Were analytical spike recoveries within the 85-115% QC limits?			<input checked="" type="checkbox"/>	
VII. ICP Serial Dilution				
Was an ICP serial dilution analyzed if analyte concentrations were > 50X the IDL?			<input checked="" type="checkbox"/>	
Were all percent differences (%Ds) ≤ 10%?			<input checked="" type="checkbox"/>	
Was there evidence of negative interference? If yes, professional judgement will be used to qualify the data.			<input checked="" type="checkbox"/>	
VIII. Internal Standards (EPA SW 846 Method 6020)				
Were all the percent recoveries (%R) within the <del>90-120</del> <sup>80-120</sup> % of the intensity of the internal standard in the associated initial calibration?	<input checked="" type="checkbox"/>			
If the %Rs were outside the criteria, was a reanalysis performed?			<input checked="" type="checkbox"/>	
IX. Regional Quality Assurance and Quality Control				
Were performance evaluation (PE) samples performed?			<input checked="" type="checkbox"/>	
Were the performance evaluation (PE) samples within the acceptance limits?			<input checked="" type="checkbox"/>	
X. Sample Result Verification				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	<input checked="" type="checkbox"/>			
XI. Overall assessment of data				
Overall assessment of data was found to be acceptable.	<input checked="" type="checkbox"/>			
XII. Field duplicates				
Field duplicate pairs were identified in this SDG.	<input checked="" type="checkbox"/>			
Target analytes were detected in the field duplicates.	<input checked="" type="checkbox"/>			
XIII. Field blanks				
Field blanks were identified in this SDG.	<input checked="" type="checkbox"/>			
Target analytes were detected in the field blanks.			<input checked="" type="checkbox"/>	

LDC #: 1127264  
 SDG #: Use conc

VALIDATION FINDINGS WORKSHEET  
 PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace Metals (EPA SW 846 Method 6010/7000) Soil preparation factor applied: 1.5X from 750mg  
 Sample Concentration units, unless otherwise noted: ug/L Associated Samples: 1-19, 4

Page: 1 of 4  
 Reviewer: MK  
 2nd Reviewer: [Signature]

Analyte	Maximum PB* (mg/Kg)	Maximum PB* (ug/L)	Maximum ICB/CCB* (ug/L)	Blank Action Limit	1	4	7	9	11	14	17	2	8	
Al														Al
Sb														Sb
As														As
Ba														Ba
Be														Be
Cd														Cd
Ca														Ca
Cr			0.284	1.995 1.495 1.16	1.3	1.3	1.4	0.8	1	0.7	0.9	1.6	1.5	Cr
Co														Co
Cu														Cu
Fe														Fe
Pb														Pb
Mg														Mg
Mn														Mn
Hg														Hg
Ni														Ni
K														K
Se														Se
Ag														Ag
Na														Na
Tl														Tl
V														V
Zn														Zn
B														B
Mo														Mo
Sr														Sr

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".  
 Note: a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

LDC #: 11272 E4  
 SDG #: See cover

VALIDATION FINDINGS WORKSHEET

METHOD: Trace Metals (EPA SW 846 Method 6010/7000) Soil preparation factor applied: 1.25  
 Sample Concentration units, unless otherwise noted: ug/kg

Associated Samples: PB: 20, 43, 43, ICB/CCB: 20, 22-40, 43, 43

Page: 2 of 84  
 Reviewer: MW  
 2nd Reviewer: J

Analyte	Maximum PB* (mg/Kg)	Maximum PB* (ug/L)	Maximum ICB/CCB* (ug/L)	Blank Action Limit	27	28	29	33	36	37	39	43	20	23
Al														
Sb														
As														
Ba														
Be														
Cd														
Ca														
Cr		0.46	0.442	2.00 2.00 2.00	1.2	1.1	1.7	1.0	1.9	1.5	1.8	1.2	2.4	2.6
Cu														
Fe														
Pb														
Mg														
Mn														
Hg														
Ni														
K														
Se														
Ag														
Na														
Tl														
V														
Zn														
B														
Mo														
Sr														

Sample Identification:

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected. "U".  
 Note: a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

LDC #: 112224  
 SDG #: SAC  
 VALIDATION FINDINGS WORKSHEET  
 PB/ICB/CCB QUALIFIED SAMPLES  
 METHOD: Trace Metals (EPA SW 846 Method 6010/7000) Soil preparation factor applied:  
 Sample Concentration units, unless otherwise noted: ug/L Associated Samples: PB: 20, 43, 43, ICB/CCB: 20, 22-40, 45, 43  
 Page: 3 of 4  
 Reviewer: MP  
 2nd Reviewer:

Analyte	Maximum PB* (mg/Kg)	Maximum PB* (ug/L)	Maximum ICB/CCB* (ug/L)	Blank Action Limit	30	34	40	Sample Identification
Al								
Sb								
As								
Ba								
Be								
Cd								
Ca				2.76				
Cr		0.461	0.442	2.765 Limit	2.7	2.6	2.4	
Cc								
Cu								
Fe								
Pb								
Mg								
Mn								
Hg								
Ni								
K								
Se								
Ag								
Na								
Tl								
V								
Zn								
B								
Mo								
Sr								

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".  
 Note: a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

VALIDATION FINDINGS WORKSHEET  
 PB/CB/CCB QUALIFIED SAMPLES

Soil preparation factor applied: 1.25X Associated Samples: 2

Analyte	Maximum PB* (mg/Kg)	Maximum PB* (ug/L)	Maximum CB/CCB* (ug/L)	Blank Action Limit	Sample Identification
Al					
Sb					
As					
Ba					
Be					
Cd					
Ca					
Cr	0.1625 0.446	0.420	2.0	0.4	21
Cc					
Cu					
Fe					
Pb					
Mg					
Mn					
Hg					
Ni					
K					
Se					
Ag					
Na					
Tl					
V					
Zn					
B					
Mo					
Sr					

Samples with analyte concentrations within five times the associated [CB, CCB or PB] concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".  
 Note: a - The listed analyte concentration is the highest [CB, CCB, or PB] detected in the analysis of each element.

LDC #: 11272 E4  
SDG #: See com

Page: 1 of 1  
Reviewer: MD  
2nd Reviewer: A

**VALIDATION FINDINGS WORKSHEET**  
**Matrix Spike/Matrix Spike Duplicates**

METHOD: Trace metals (EPA SW 846 Method 6010/7000)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

N N/A Was a matrix spike analyzed for each matrix in this SDG?  
 N N/A Were matrix spike percent recoveries (%R) within the control limits of 75-125? *Lab limit*  
if the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.

N N/A Were all duplicate sample relative percent differences (RPD)  $\leq 20\%$  for water samples and  $\leq 35\%$  for soil samples?

N N/A Were recalculated results acceptable? See Level IV Recalculation Worksheet for recalculations.

#	MS/MSD ID	Matrix	Analyte	MS %Recovery	MSD %Recovery	RPD (Limits)	Associated Samples	Qualifications
1	C4155	A2	CY	68.4 (80-120)	72.8 (80-120)		1-19, 41	J/UT/A

Comments:



LDC #: 11272E4  
 SDG #: See com

**VALIDATION FINDINGS WORKSHEET**  
**Field Duplicates**

Page: 1 of 1  
 Reviewer: ML  
 2nd reviewer: A

METHOD: Cr ~~Inorganics~~, Method 200.8

N N/A Were field duplicate pairs identified in this SDG?  
 N N/A Were target analytes detected in the field duplicate pairs?

Analyte	Concentration (ug/L)		RPD (Limits)	Qualifier
	4	7		
Cr	1.3	1.4	7	

Analyte	Concentration (ug/L)		RPD (Limits)	Qualifier
	19	24		
Cr	3.5	5.6	46	

Analyte	Concentration (ug/L)		RPD (Limits)	Qualifier
	27	28		
Cr	1.2	1.1	9	

Analyte	Concentration ( )		RPD (Limits)	Qualifier

LDC #: 112p264  
 SDG #: See cover

**VALIDATION FINDINGS WORKSHEET**  
**Initial and Continuing Calibration Calculation Verification**

Page: 1 of 1  
 Reviewer: M10  
 2nd Reviewer: A

METHOD: Trace Metals (EPA SW 846 Method 6010/7000)

An initial and continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$\%R = \frac{\text{Found} \times 100}{\text{True}}$  Where Found = concentration (in ug/L) of each analyte measured in the analysis of the ICV or CCV solution  
 True = concentration (in ug/L) of each analyte in the ICV or CCV source

Standard ID	Type of Analysis	Element	Found (ug/L)	True (ug/L)	Recalculated		Reported		Acceptable (Y/N)
					%R	%R	%R	%R	
<u>ICV</u>	<u>ICP/Initial calibration</u>	<u>Cr</u>	<u>10.19</u>	<u>10</u>	<u>102</u>	<u>102</u>	<u>102</u>	<u>Y</u>	
	GFAA (Initial calibration)								
	CVAA (Initial calibration)								
<u>CCV</u>	<u>ICP/Continuing calibration</u>	<u>Cr</u>	<u>9.958</u>	<u>10</u>	<u>100</u>	<u>100</u>	<u>100</u>	<u>Y</u>	
	GFAA (Continuing calibration)								
	CVAA (Continuing calibration)								
	Cyanide (Initial calibration)								
	Cyanide (Continuing calibration)								

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 11272-24  
 SDG #: See Cover

**VALIDATION FINDINGS WORKSHEET**  
**Level IV Recalculation Worksheet**

Page: 1 of 1  
 Reviewer: juv  
 2nd Reviewer: [Signature]

METHOD: Trace Metals (EPA SW 846 Method 6010/7000)

Percent recoveries (%R) for an ICP interference check sample, a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found} \times 100}{\text{True}}$$
 Where, Found = Concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation,  
 Found = SSR (spiked sample result) - SR (sample result).  
 True = Concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$RPD = \frac{|S-D|}{(S+D)/2} \times 100$$
 Where, S = Original sample concentration  
 D = Duplicate sample concentration

An ICP serial dilution percent difference (%D) was recalculated using the following formula:

$$\%D = \frac{|I-SDRI|}{I} \times 100$$
 Where, I = Initial Sample Result (mg/L)  
 SDR = Serial Dilution Result (mg/L) (Instrument Reading x 5)

Sample ID	Type of Analysis	Element	Found / S / I (units)	True / D / SDR (units)	Recalculated		Reported		Acceptable (Y/N)
					%R / RPD / %D	%R / RPD / %D			
NA	ICP interference check	I							
LC5	Laboratory control sample	Cr	10.153	10	102	102	102		
54	Matrix spike	Cr	(SSR-SR) 6.838	10	68.4	68.4	68.4		Y
56	Duplicate	Cr	4.362	4.286	1.76	1.76	1.76		Y
NA	ICP serial dilution								

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 1127224  
 SDG #: See work

**VALIDATION FINDINGS WORKSHEET**  
Sample Calculation Verification

Page: 1 of 1  
 Reviewer: MJS  
 2nd reviewer: [Signature]

METHOD: Trace Metals (EPA SW 846 Method 6010/7000)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- Y / N / N/A Have results been reported and calculated correctly?
- Y / N / N/A Are results within the calibrated range of the instruments and within the linear range of the ICP?
- Y / N / N/A Are all detection limits below the CRDL?

Detected analyte results for 7 were recalculated and verified using the following equation:

$$\text{Concentration} = \frac{(RD)(FV)(Dil)}{(In. Vol.)(\%S)}$$

Recalculation:

*From the row 10th*  
 $Cr = 1.370 \mu g/L$

- RD = Raw data concentration
- FV = Final volume (ml)
- In. Vol. = Initial volume (ml) or weight (G)
- Dil = Dilution factor
- %S = Decimal percent solids

Sample ID	Analyte	Reported Concentration ( $\mu g/L$ )	Calculated Concentration ( $\mu g/L$ )	Acceptable (Y/N)
7	Cr	1.4	1.4	Y