

```

=====
Sample Name: 2809-1 F=2.5                      Date: 04/22/2003 11:58:39
Data File  : C:\DX\DATA\03W2479\2809-101.D07
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 7           Detector:COND
Analyst    : David                            Column: Dionex AS4A-SC
=====

```

```

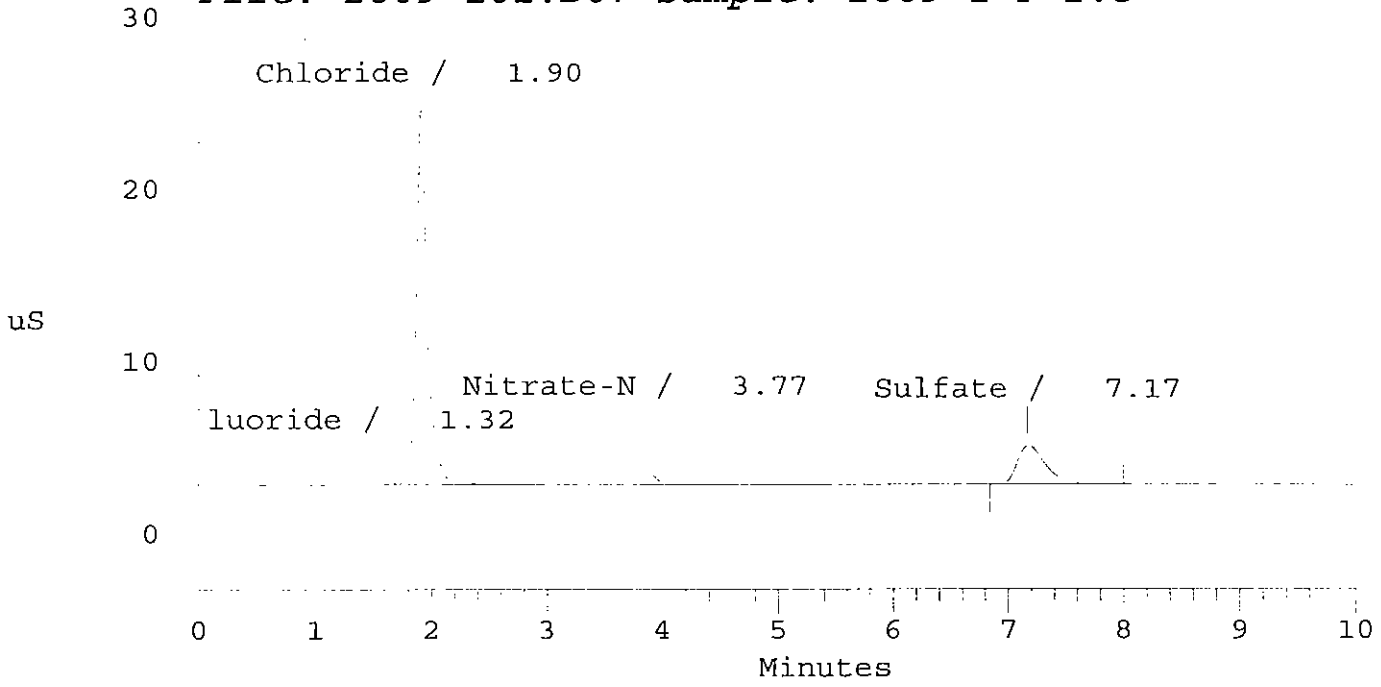
-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External          1          2.5    3000  5Hz   0.00 10.00    1000
-----

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	0.728	16910	212174	2	0.00
2	1.90	Chloride	24.191	686962	4889400	2	0.00
0	0.00	Nitrite-N	0.000	0	0	0	0.00
0	0.00	Bromide	0.000	0	0	0	0.00
3	3.77	Nitrate-N	1.746	80734	814055	1	0.00
0	0.00	Phosphate-P	0.000	0	0	0	0.00
4	7.17	Sulfate	9.537	75701	1296160	1	0.00
Totals			36.202	860307	7211789		

File: 2809-101.D07 Sample: 2809-1 F=2.5



```

=====
Sample Name: MB RW1408                               Date: 04/22/2003 10:25:06
Data File  : C:\DX\DATA\03W2479\W2479K01.D02
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 2                 Detector: COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====
    
```

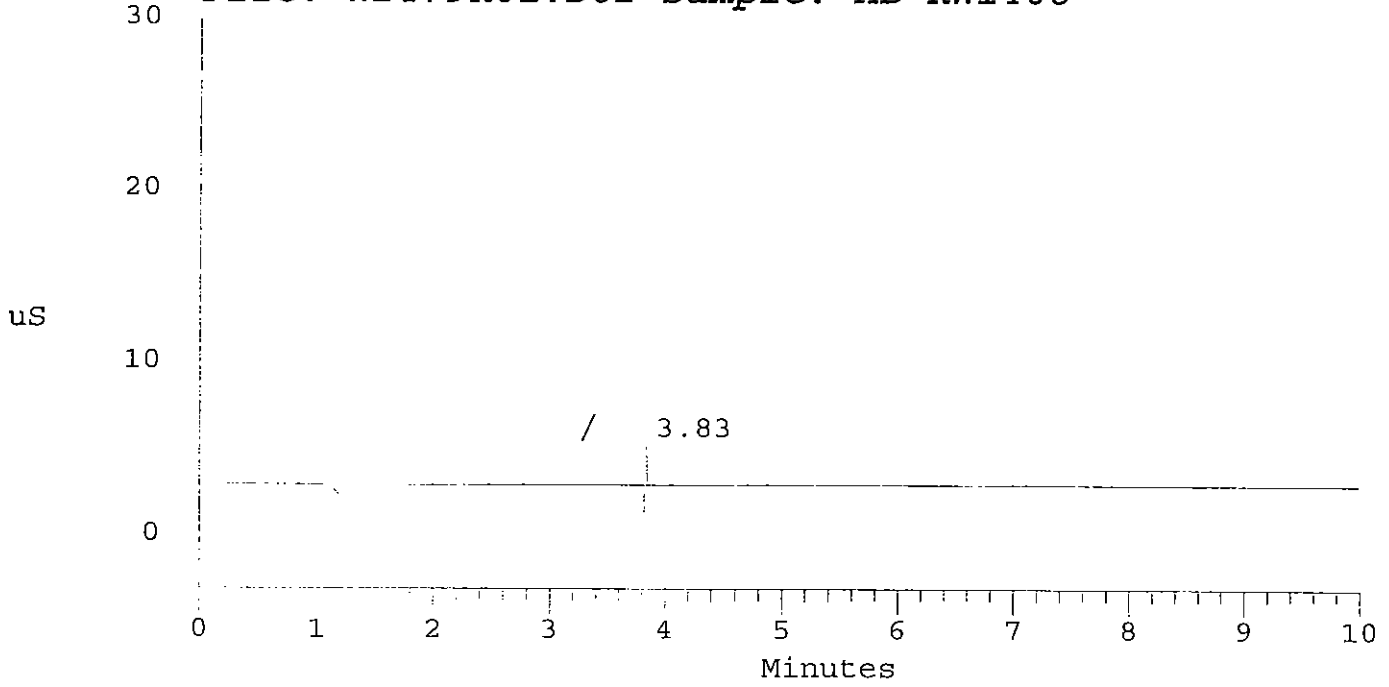
```

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

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

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

File: W2479K01.D02 Sample: MB RW1408



After
 4/22/03 DV
 reason: 3.5
 4578

```

=====
Sample Name: MB RW1408                               Date: 04/22/2003 10:25:06
Data File  : C:\DX\DATA\03W2479\W2479K01.D02
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 2                 Detector: COND
Analyst    : David                                   Column: Dionex AS4A-SC
=====

```

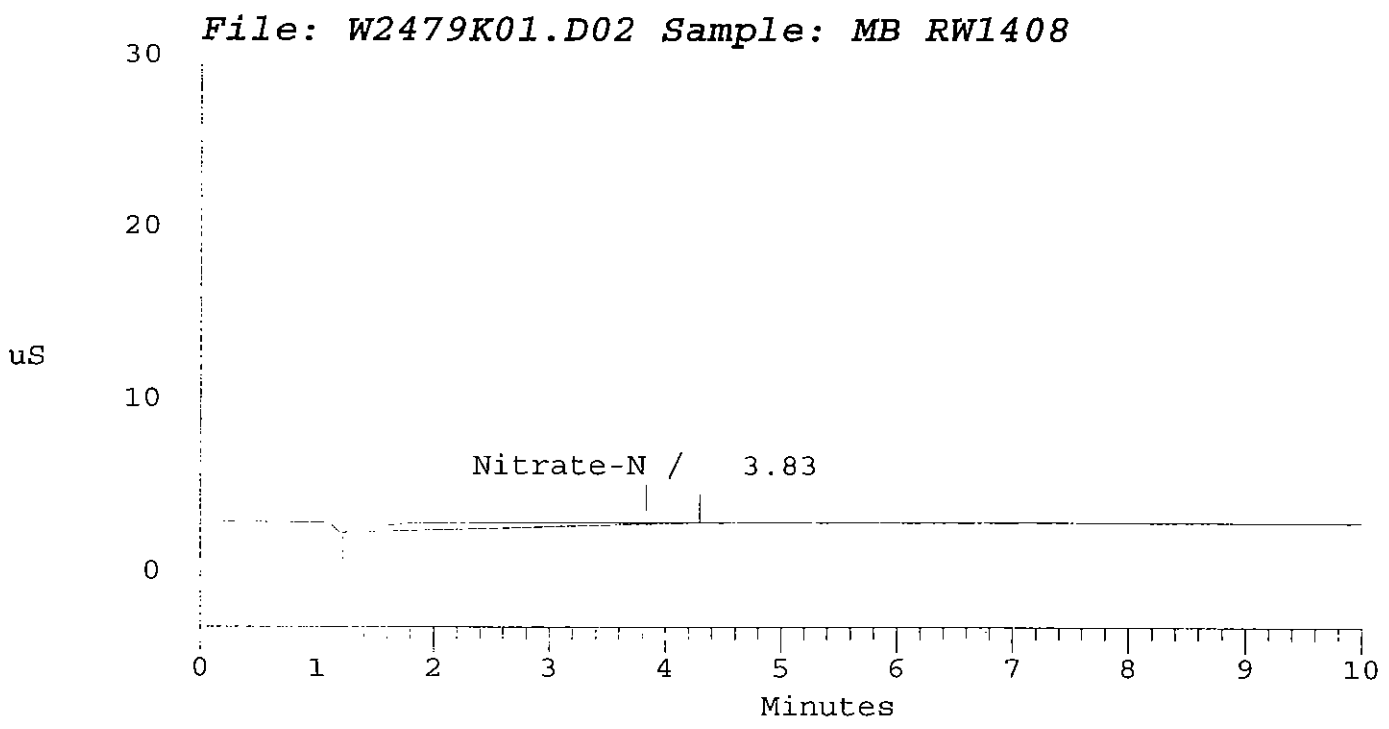
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
0	0.00	Fluoride	0.000	0	0	0	0.00
0	0.00	Chloride	0.000	0	0	0	0.00
0	0.00	Nitrite-N	0.000	0	0	0	0.00
0	0.00	Bromide	0.000	0	0	0	0.00
1	3.83	Nitrate-N	1.307	3121	1570052	1	0.00
0	0.00	Phosphate-P	0.000	0	0	0	0.00
0	0.00	Sulfate	0.000	0	0	0	0.00
Totals			1.307	3121	1570052		



before

```

=====
Sample Name: MB RW1408                               Date: 04/22/2003 14:06:15
Data File  : C:\DX\DATA\03W2479\W2479K01.D13
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 13                 Detector:COND
Analyst    : David                                   Column: Dionex AS4A-SC
=====

```

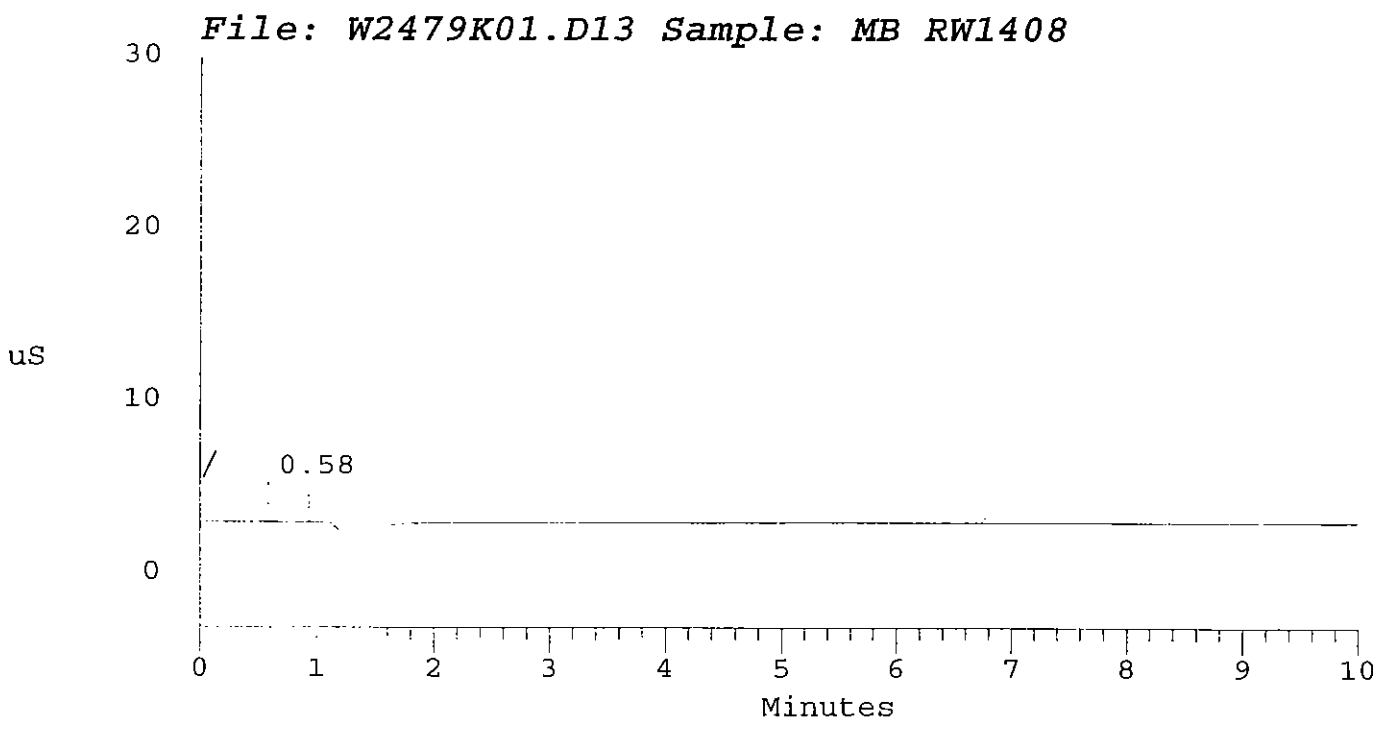
```

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

```

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

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



```

=====
Sample Name: MB RW1408                               Date: 04/22/2003 16:46:10
Data File  : C:\DX\DATA\03W2479\W2479K11.D19
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 19                 Detector:COND
Analyst    : David      Column: Dionex AS4A-SC
=====

```

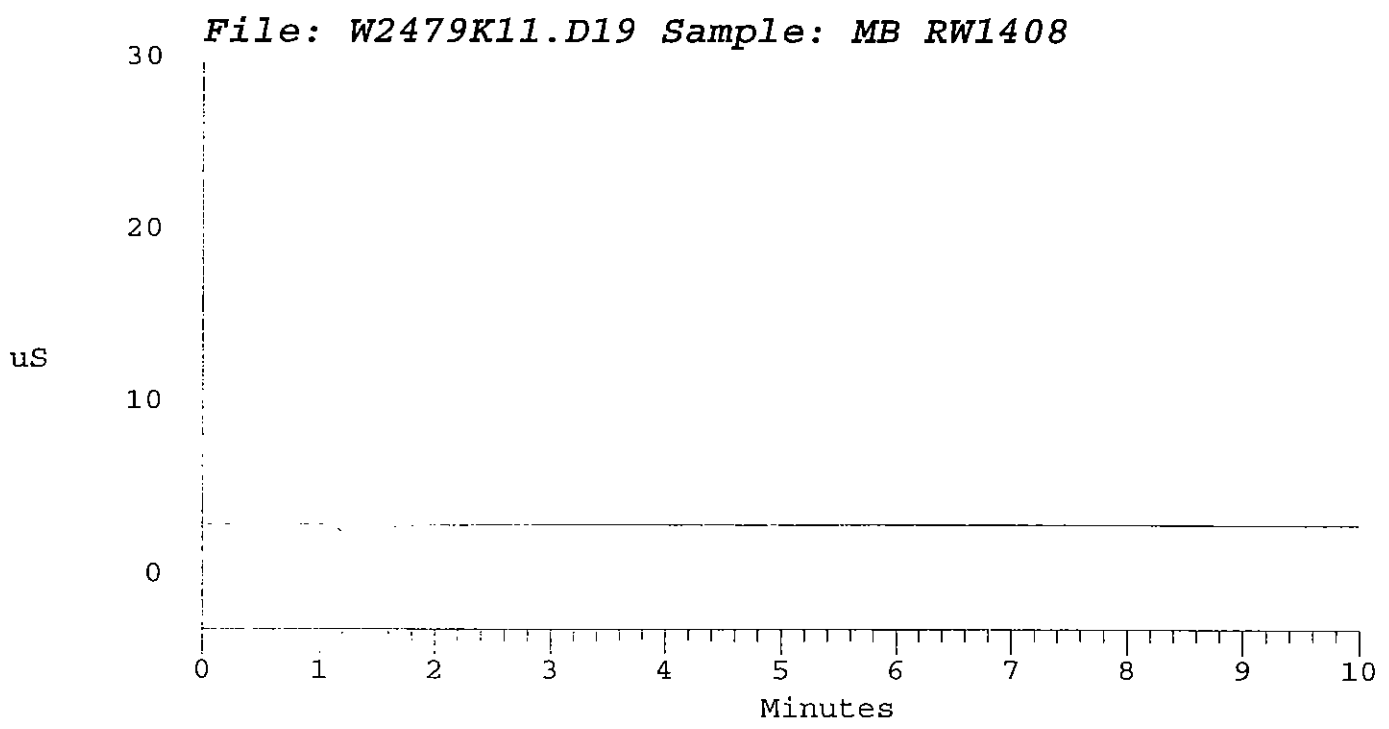
```

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

```

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

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



FORM-3

Applied P & Ch Laboratory

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

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2479	
LCS Filename: -	Date Analyzed: 042203	Time Analyzed: 10:43
LCSD Filename: -	Date Analyzed: 042203	Time Analyzed: 10:56

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
CHLORIDE CL ⁻	mg/L	4.0	0	3.96	99	80-120
NITRATE AS N	mg/L	1.5	0	1.51	101	80-120
SULFATE SO ₄ ⁻	mg/L	15	0	14.8	99	80-120
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHLORIDE CL ⁻	mg/L	4.0	3.93	98	1	20	80-120
NITRATE AS N	mg/L	1.5	1.50	100	1	20	80-120
SULFATE SO ₄ ⁻	mg/L	15	14.7	98	1	25	80-120
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

```

=====
Sample Name: LCS W7768-100X                               Date: 04/22/2003 10:43:49
Data File  : C:\DX\DATA\03W2479\W2479L01.D03
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 3                       Detector:COND
Analyst    : David                                         Column: Dionex AS4A-SC
=====

```

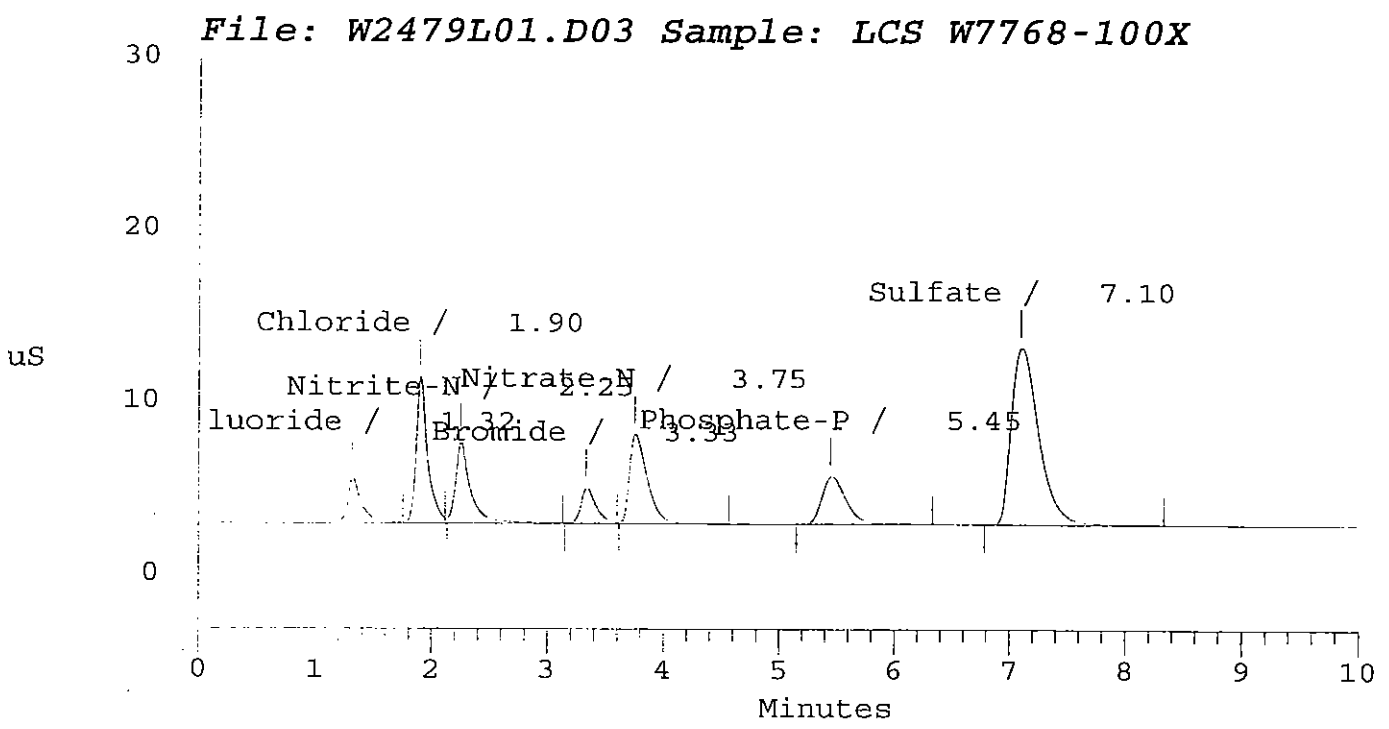
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	1.025	87568	745302	2	0.00
2	1.90	Chloride	3.956	277396	1960299	2	0.00
3	2.25	Nitrite-N	1.438	154386	1448414	2	0.00
4	3.33	Bromide	3.034	66690	618380	2	0.37
5	3.75	Nitrate-N	1.514	170870	1826336	2	0.00
6	5.45	Phosphate-P	2.928	90756	1338642	1	0.00
7	7.10	Sulfate	14.753	338834	5615256	1	0.00
Totals			28.648	1186501	13552630		



```

=====
Sample Name: LCSD W7768-100X           Date: 04/22/2003 10:56:32
Data File  : C:\DX\DATA\03W2479\W2479J01.D04
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 4      Detector: COND
Analyst    : David                     Column: Dionex AS4A-SC
=====

```

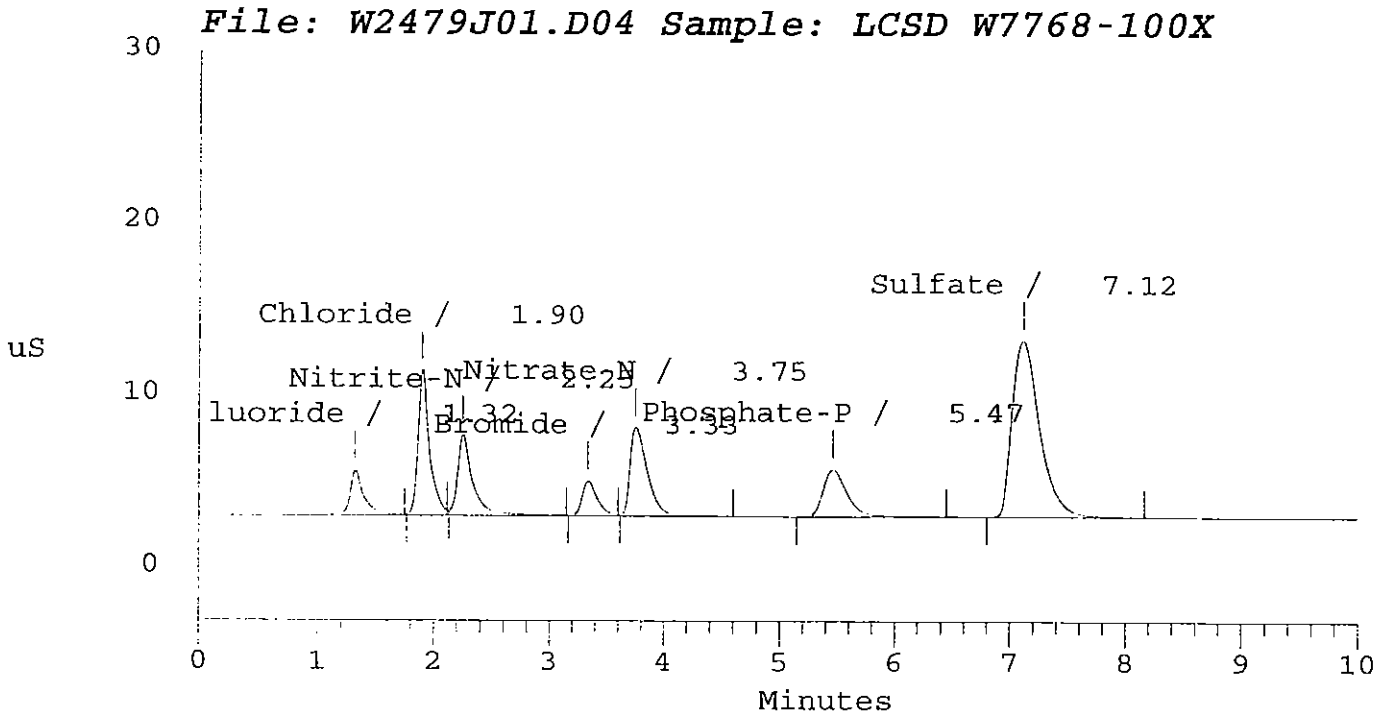
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	1.020	86862	741871	2	0.00
2	1.90	Chloride	3.930	277914	1946867	2	0.00
3	2.25	Nitrite-N	1.442	154917	1453082	2	0.00
4	3.33	Bromide	3.022	67290	615743	2	0.37
5	3.75	Nitrate-N	1.502	170692	1812154	2	0.00
6	5.47	Phosphate-P	2.919	91574	1334304	1	0.00
7	7.12	Sulfate	14.659	339223	5578007	1	0.00
Totals			28.494	1188472	13482030		



FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 300.0

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2479	
MS Filename: -	Date Analyzed: 042203	Time Analyzed: 14:22
MSD Filename: -	Date Analyzed: 042203	Time Analyzed: 14:35
MS Sample No: MW-4-2	Sample Lab ID: 03-2809-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
CHLORIDE CL ⁻	mg/L	160	95.4	257	101	75-125
NITRATE AS N	mg/L	60.0	10.0	69.3	99	75-125
SULFATE SO ₄ ⁻	mg/L	600	124	719	99	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHLORIDE CL ⁻	mg/L	160	256	100	1	20	75-125
NITRATE AS N	mg/L	60.0	70.1	100	1	20	75-125
SULFATE SO ₄ ⁻	mg/L	600	713	98	1	25	75-125
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

```

=====
Sample Name: $2809-4 MS F=40                               Date: 04/22/2003 14:22:48
Data File  : C:\DX\DATA\03W2479\W2479M01.D14
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 14                       Detector: COND
Analyst    : David                                         Column: Dionex AS4A-SC
=====

```

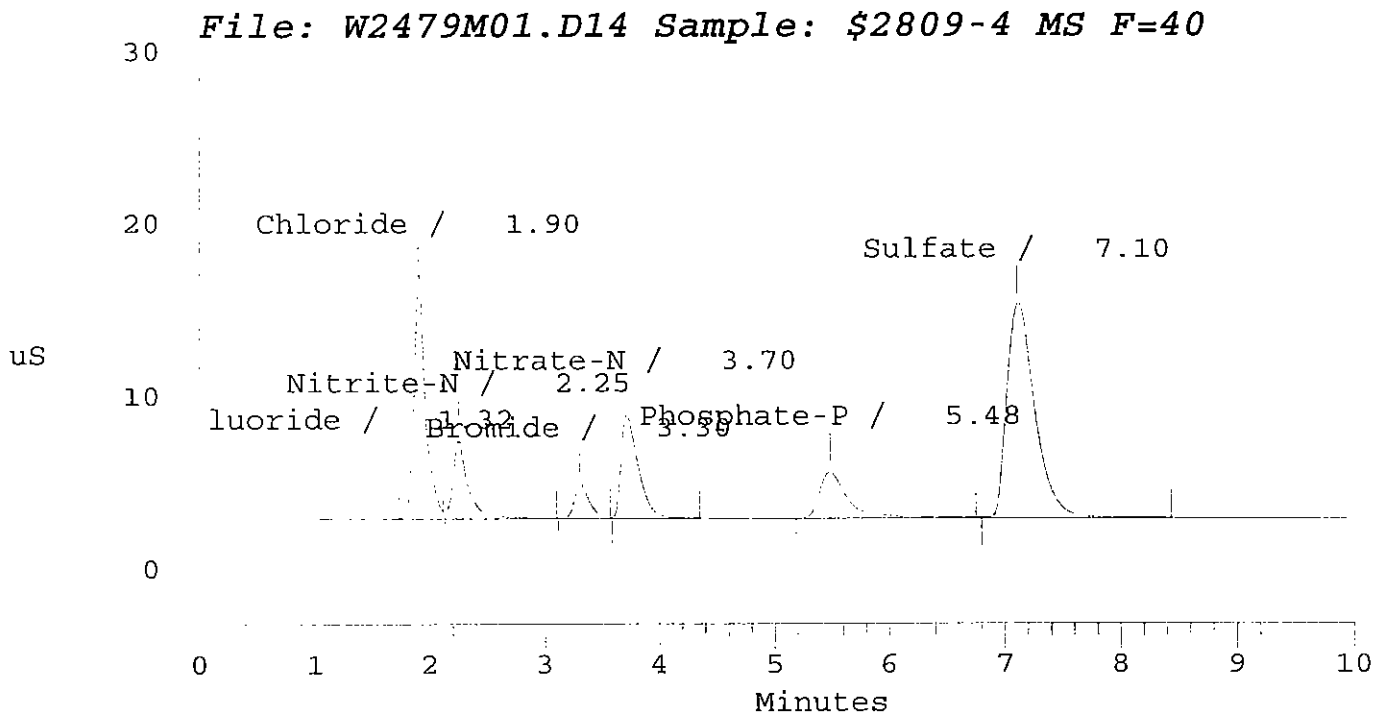
```

-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External          1          40    3000   5Hz   0.00 10.00      1000
-----

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	41.825	82701	760523	2	0.00
2	1.90	Chloride	256.662	459912	3220066	2	0.00
3	2.25	Nitrite-N	56.934	152454	1433599	2	0.00
4	3.30	Bromide	120.216	66386	612379	2	0.71
5	3.70	Nitrate-N	69.326	195629	2098781	2	0.00
6	5.48	Phosphate-P	136.572	88842	1567853	2	0.00
7	7.10	Sulfate	719.313	410163	6890583	2	0.00
Totals			1400.848	1456087	16583784		



```

=====
Sample Name: $2809-4 MSD F=40                               Date: 04/22/2003 14:35:30
Data File  : C:\DX\DATA\03W2479\W2479N01.D15
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 15                       Detector:COND
Analyst    : David                                         Column: Dionex AS4A-SC
=====

```

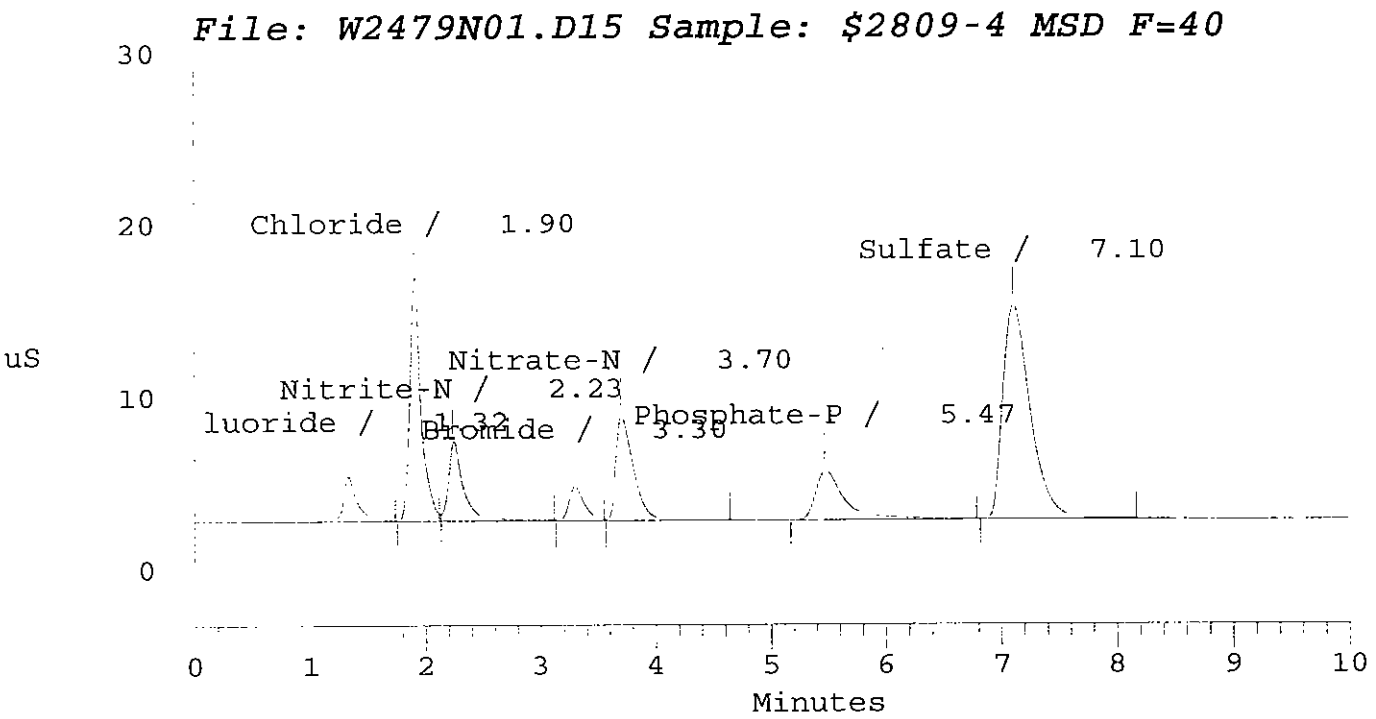
```

-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External          1          40    3000  5Hz   0.00 10.00      1000

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	41.597	82508	756379	2	0.00
2	1.90	Chloride	255.602	464903	3206498	2	0.00
3	2.23	Nitrite-N	57.448	145928	1446758	2	-0.74
4	3.30	Bromide	120.209	66738	612341	2	0.71
5	3.70	Nitrate-N	70.064	199598	2121675	2	0.00
6	5.47	Phosphate-P	149.479	93590	1719897	2	0.00
7	7.10	Sulfate	713.046	410183	6828718	2	0.00
Totals			1407.444	1463449	16692266		



FORM-3

Applied P & Ch Laboratory

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

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2496	
LCS Filename: -	Date Analyzed: 042203	Time Analyzed: 18:49
LCSD Filename: -	Date Analyzed: 042203	Time Analyzed: 19:07

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
PERCHLORATE	µg/L	25	0	20.8	83	80-120
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
PERCHLORATE	µg/L	25	20.4	82	1	20	80-120
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

APCL Perchlorate Analysis Report

Sample Name : lcs 25ppb w7827d

Data File Name : C:\DATA\03W2496K\W2496K L01_004.DXD

Method File Name : c:\peaknet\method\314-011.met

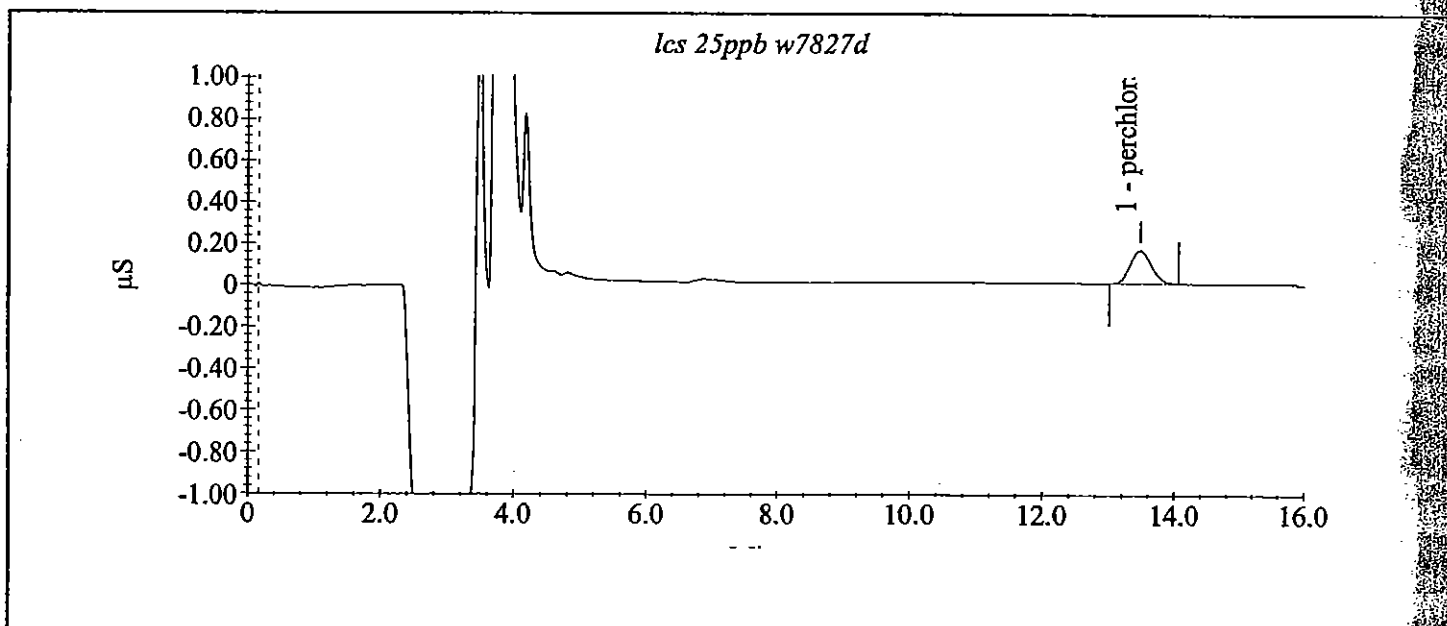
Date Time Collected : 04/22/2003 6:49:00 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.48	20.81	35317.40	1591.25



APCL Perchlorate Analysis Report

Sample Name : LCS 18PPB W7685D

Data File Name : C:\DATA\03W2496K\W2496K J01_005.DXD

Method File Name : c:\peaknet\method\314-011.met

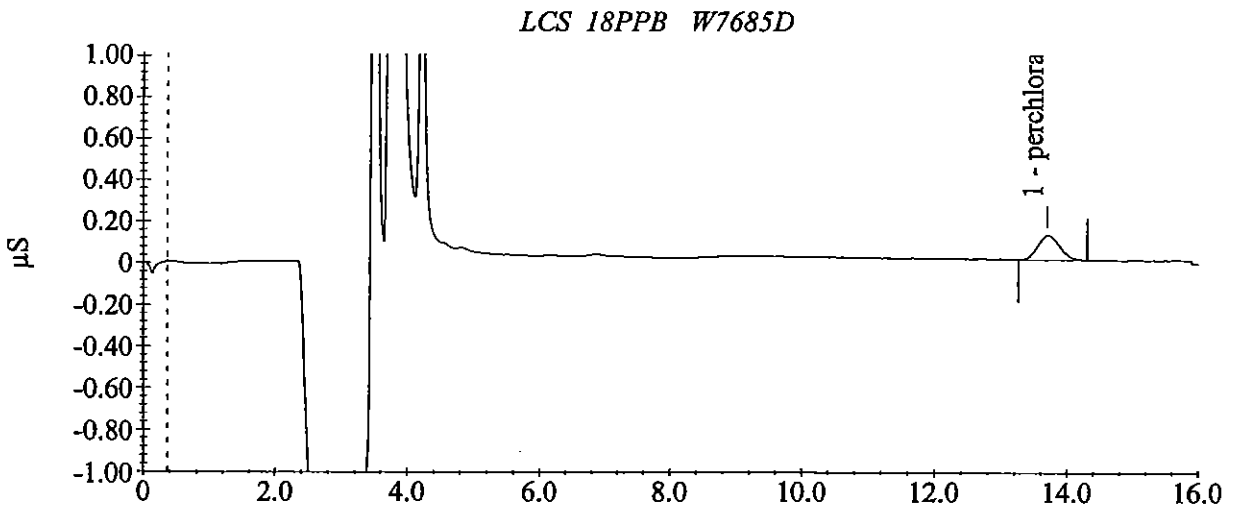
Date Time Collected : 04/22/2003 7:07:40 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.70	16.23	27547.60	1168.31



FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 314.0

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2496	
MS Filename: -	Date Analyzed: 042303	Time Analyzed: 01:57
MSD Filename: -	Date Analyzed: 042303	Time Analyzed: 02:15
MS Sample No: MW-4-2	Sample Lab ID: 03-2809-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
PERCHLORATE	µg/L	50.0	6.6	55.4	98	75-125
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
PERCHLORATE	µg/L	50.0	56.7	100	2	20	75-125
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

APCL Perchlorate Analysis Report

Sample Name : 2809-04 MS 50PPB F=1

Data File Name : C:\DATA\03W2496K\W2496K M01_027.DXD

Method File Name : c:\peaknet\method\314-011.met

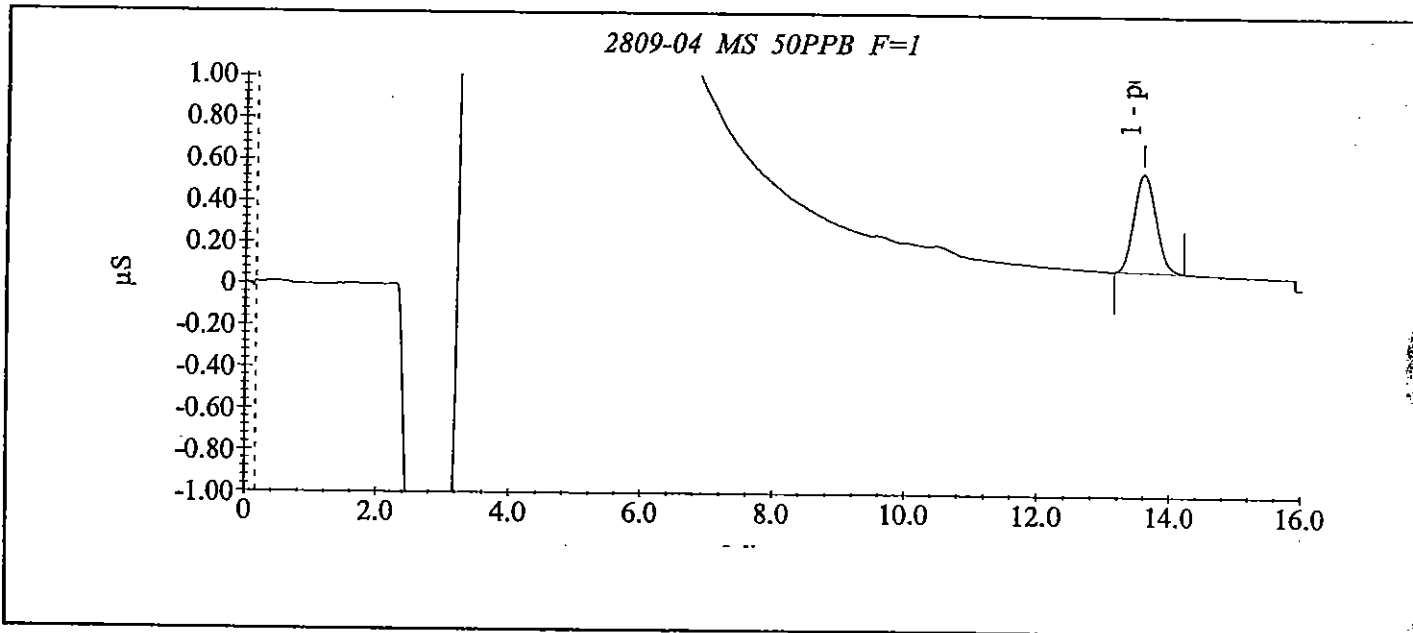
Date Time Collected : 04/23/2003 1:57:22 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.58	64.12	108809.80	4728.45



Rec 115.14%



APCL Perchlorate Analysis Report

Sample Name : 2809-04 MSD 50PPB F=1

Data File Name : C:\DATA\03W2496K\W2496K N01_028.DXD

Method File Name : c:\peaknet\method\c314-011.met

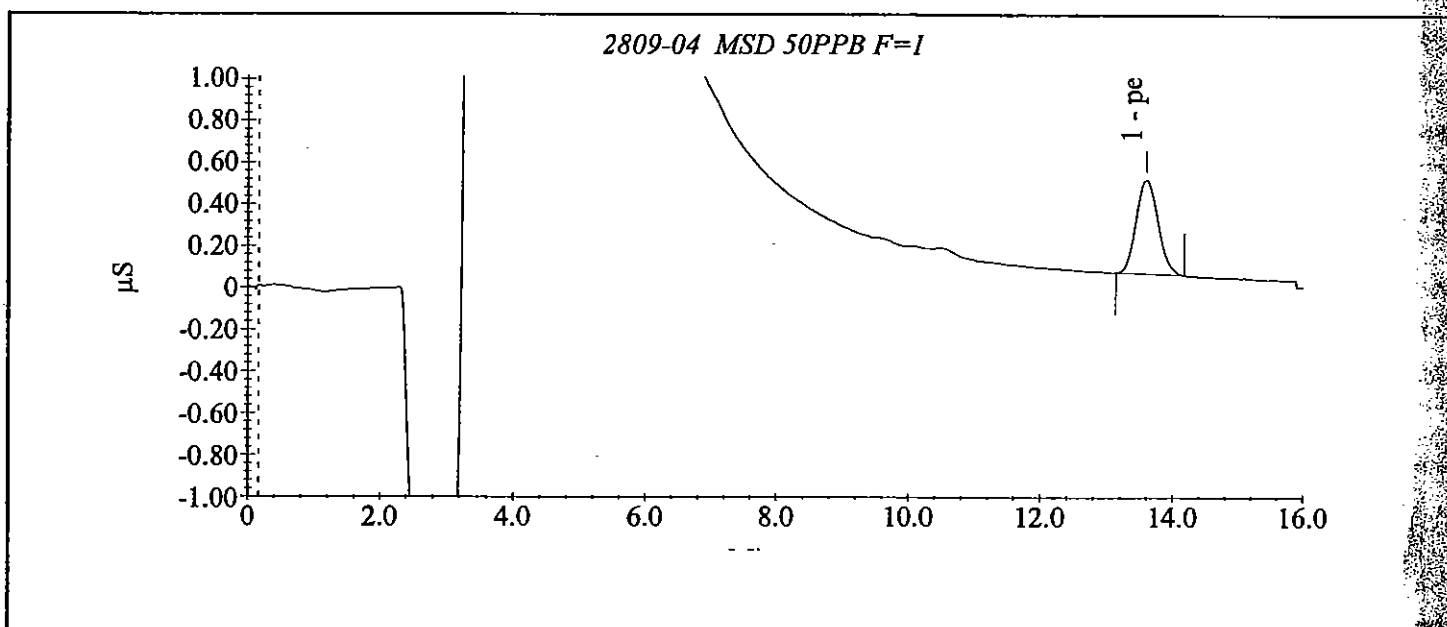
Date Time Collected : 04/23/2003 2:15:52 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.60	60.77	103133.40	4475.04



Rec 108.44%



FORM-3

Applied P & Ch Laboratory

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

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2500	
LCS Filename: -	Date Analyzed: 042203	Time Analyzed: 16:56
LCSD Filename: -	Date Analyzed: 042203	Time Analyzed: 16:56

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	0	421	105	88-108
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	428	107	2	20	88-108
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: _____

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 160.1

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2500	
MS Filename: -	Date Analyzed: 042203	Time Analyzed: 16:56
MSD Filename: -	Date Analyzed: 042203	Time Analyzed: 16:56
MS Sample No: MW-19-2	Sample Lab ID: 03-2809-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	587	1010	106	80-119
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
SOLIDS, TOTAL DISSOLVED (TDS)	mg/L	400	971	96	10	20	80-119
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

* - Values outside of contract required QC Limits

D - Spiked components diluted out

Comments: _____

FORM-3

Applied P & Ch Laboratory

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

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2476	
LCS Filename: -	Date Analyzed: 042103	Time Analyzed: 17:22
LCSD Filename: -	Date Analyzed: 042103	Time Analyzed: 17:22

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
CHROMIUM (VI)	mg/L	0.25	0	0.225	90	80-115
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHROMIUM (VI)	mg/L	0.25	0.230	92	2	19	80-115
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

FORM-3

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 7196

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 32809
Project ID: JPL	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03W2476	
MS Filename: -	Date Analyzed: 042103	Time Analyzed: 17:22
MSD Filename: -	Date Analyzed: 042103	Time Analyzed: 17:22
MS Sample No: MW-4-2	Sample Lab ID: 03-2809-4	

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

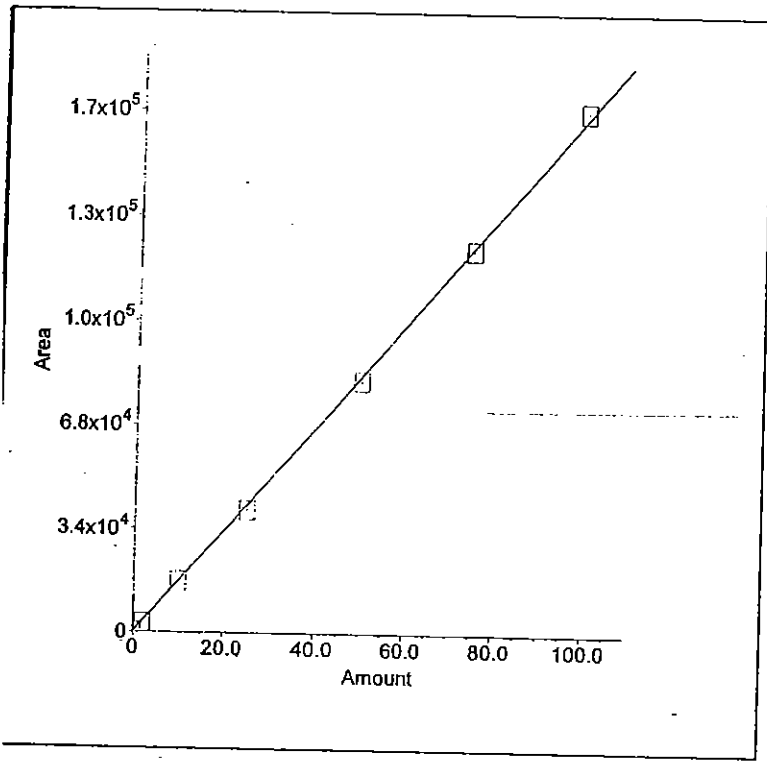
Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
CHROMIUM (VI)	mg/L	0.25	0.226	90	1	19	78-115
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

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



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

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

APCL Perchlorate Analysis Report

Sample Name : Cal blank

Data File Name : C:\data\E314-011\Mb_001.DXD

Method File Name : c:\peaknet\method\e314-011.met

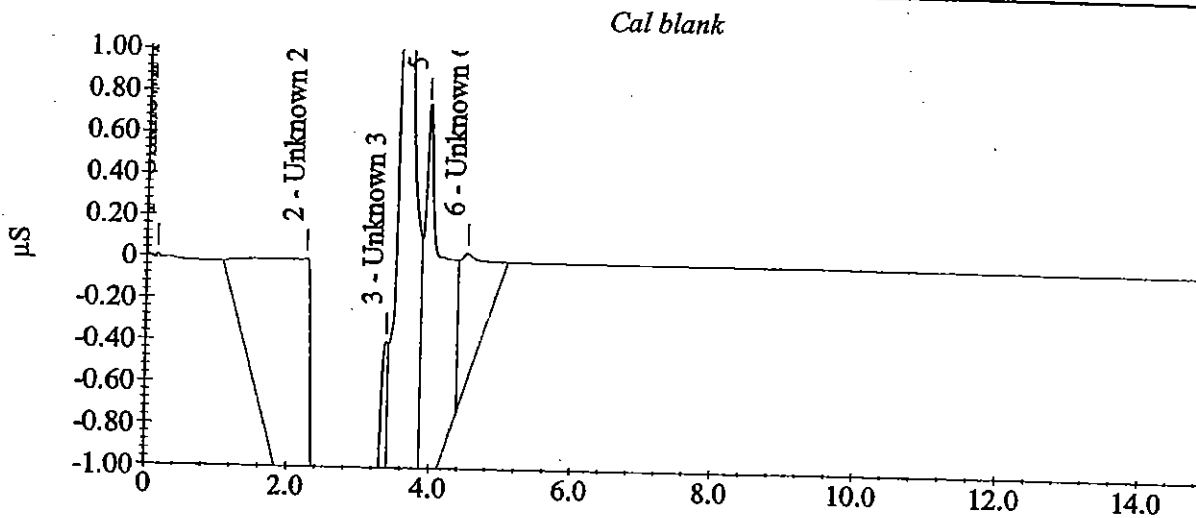
Date Time Collected : 03/12/2003 5:55:39 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
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APCL Perchlorate Analysis Report

Sample Name : cal standard 2ppb W7827a

Data File Name : C:\DATA\E314-011\std-2pb_002.DXD

Method File Name : C:\PEAKNET\METHOD\e314-011.met

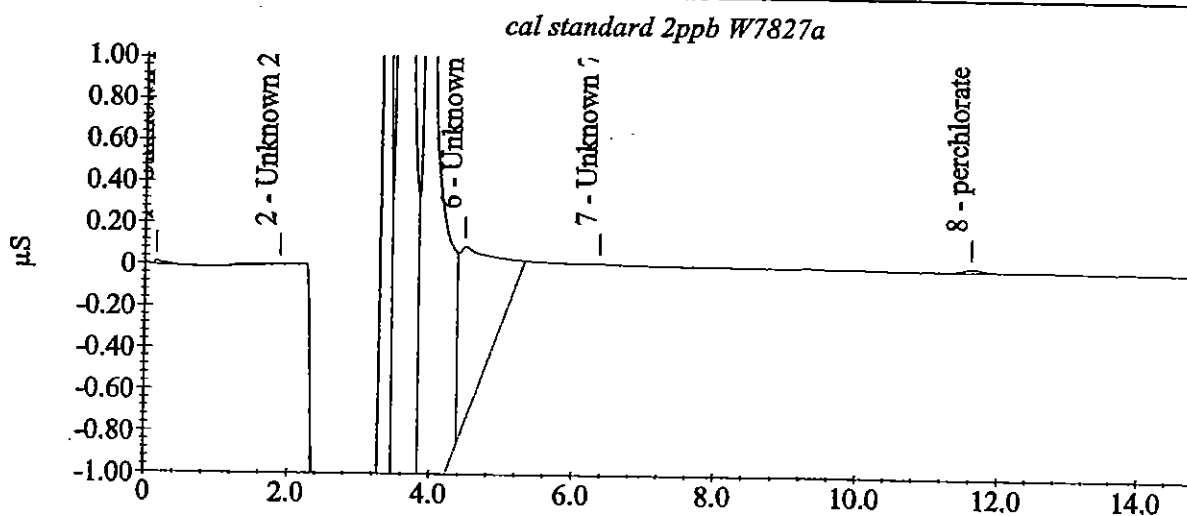
Date Time Collected : 03/12/2003 6:13:12 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
8	perchlorate	11.62	1.92	2910	164



APCL Perchlorate Analysis Report

Sample Name : cal standard 10ppb W7827c

Data File Name : C:\DATA\E314-011\std-10pb_004.DXD

Method File Name : C:\PEAKNET\METHOD\314-011.met

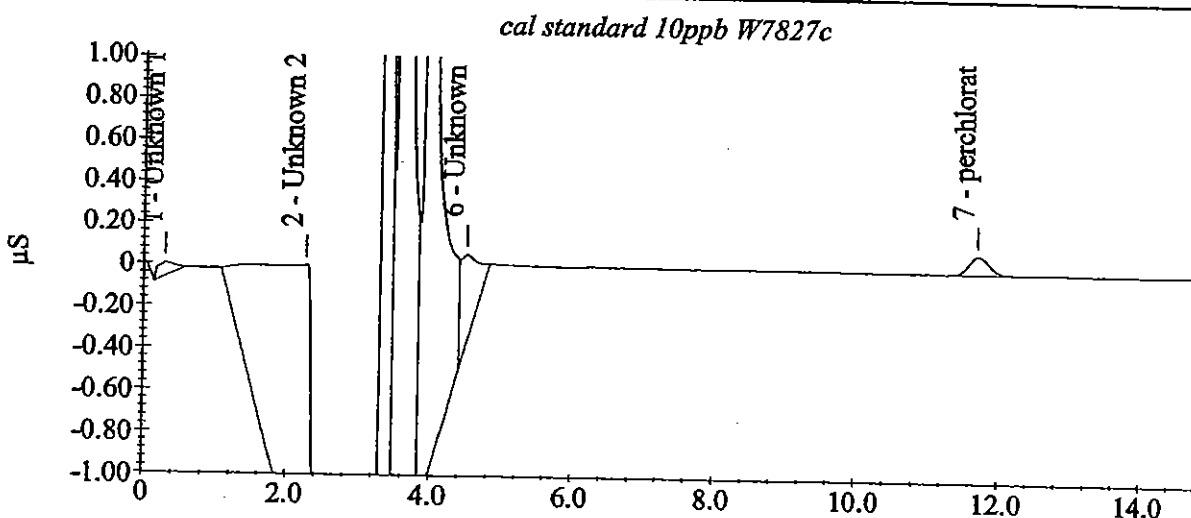
Date Time Collected : 03/12/2003 6:48:21 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
7	perchlorate	11.70	11.16	16917	879



APCL Perchlorate Analysis Report

Sample Name : cal standard 25ppb W7827d

Data File Name : C:\DATA\E314-011\std-25pb_005.DXD

Method File Name : C:\PEAKNET\METHOD\314-011.met

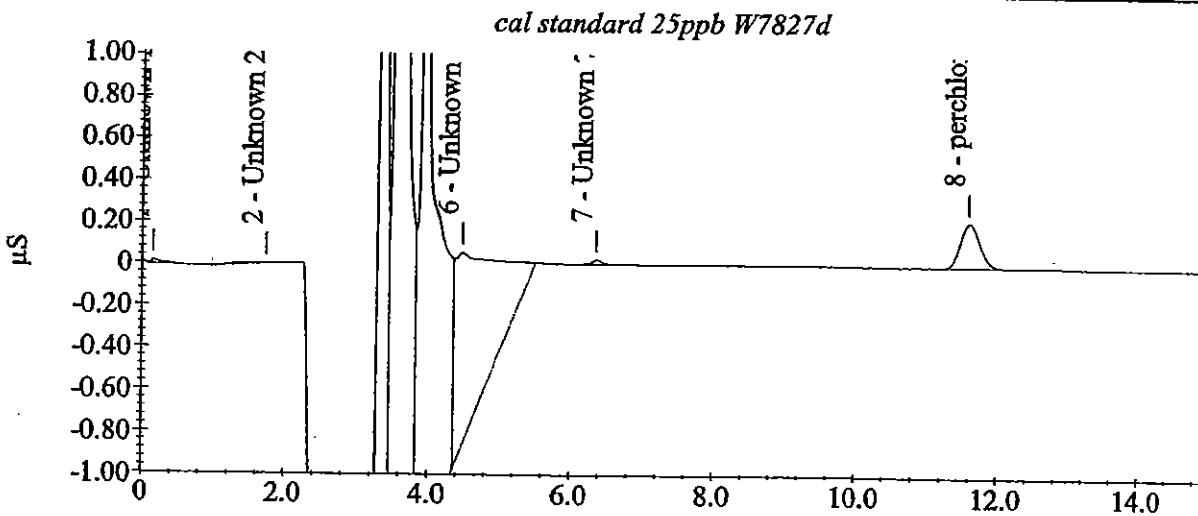
Date Time Collected : 03/12/2003 7:05:54 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
8	perchlorate	11.60	26.84	40702	2125



APCL Perchlorate Analysis Report

Sample Name : cal standard 50ppb W7827e

Data File Name : C:\DATA\E314-011\std-50pb_006.DXD

Method File Name : C:\PEAKNET\METHOD\314-011.met

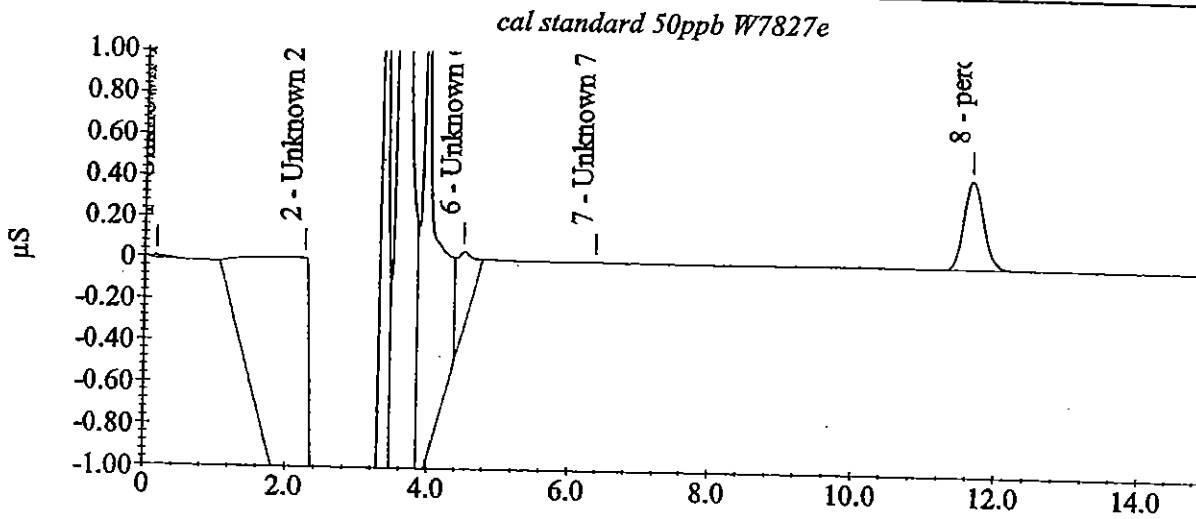
Date Time Collected : 03/12/2003 7:23:30 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
8	perchlorate	11.67	54.89	83240	4320



APCL Perchlorate Analysis Report

Sample Name : cal standard 75ppb W7827f

Data File Name : C:\DATA\E314-011\std-75pb_007.DXD

Method File Name : C:\PEAKNET\METHOD\314-011.met

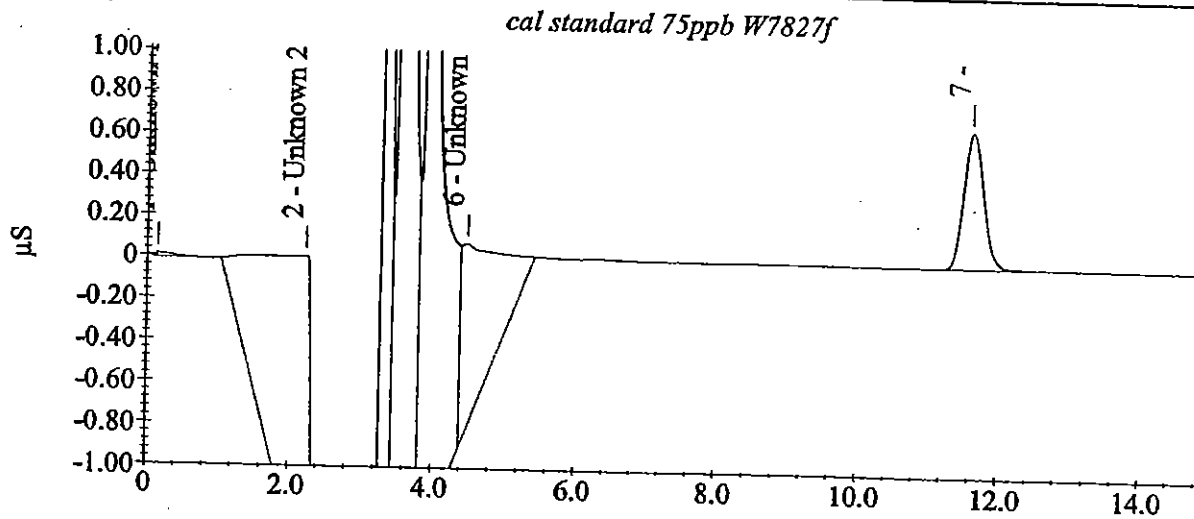
Date Time Collected : 03/12/2003 7:41:05 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
7	perchlorate	11.62	83.23	126224	6553



APCL Perchlorate Analysis Report

Sample Name : cal standard 100ppb W7827g

Data File Name : C:\DATA\E314-011\std-100pb_008.DXD

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

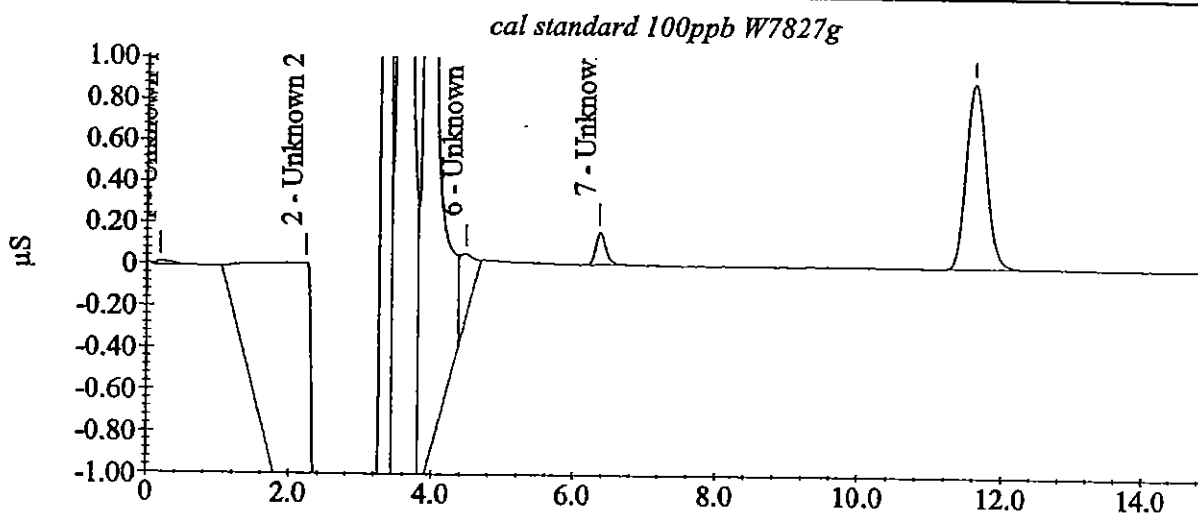
Date Time Collected : 03/12/2003 7:58:39 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
8	perchlorate	11.62	113.21	171686	8927



APCL Perchlorate Analysis Report

Sample Name : ICV 50 ppb w7828a

Data File Name : C:\DATA\E314-011\icv-50pb_009.DXD

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

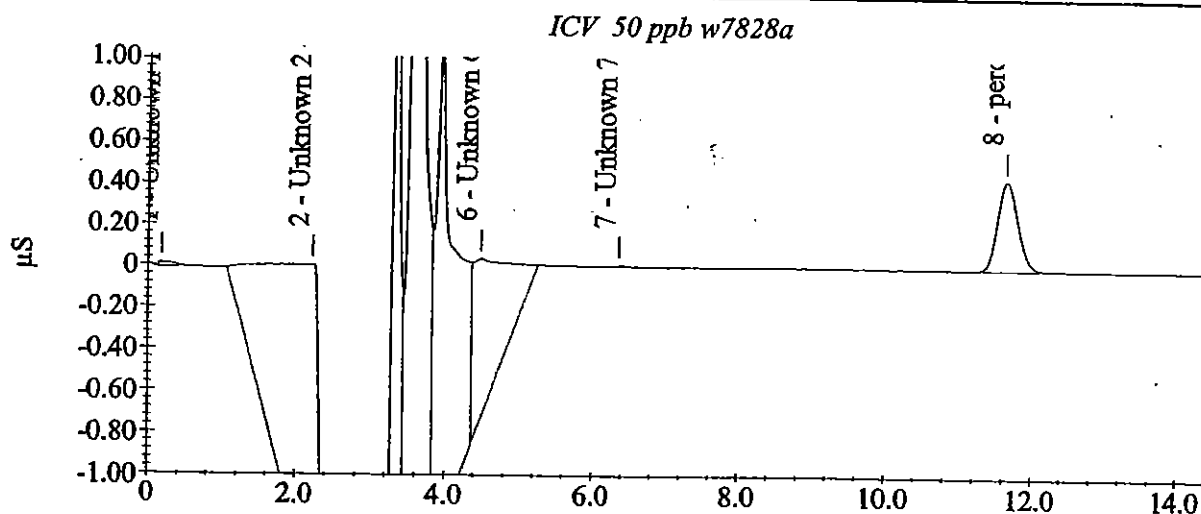
Date Time Collected : 03/12/2003 8:16:15 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

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



APCL Perchlorate Analysis Report

Sample Name : icb

Data File Name : C:\DATA\E314-011\ICB_010.DXD

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

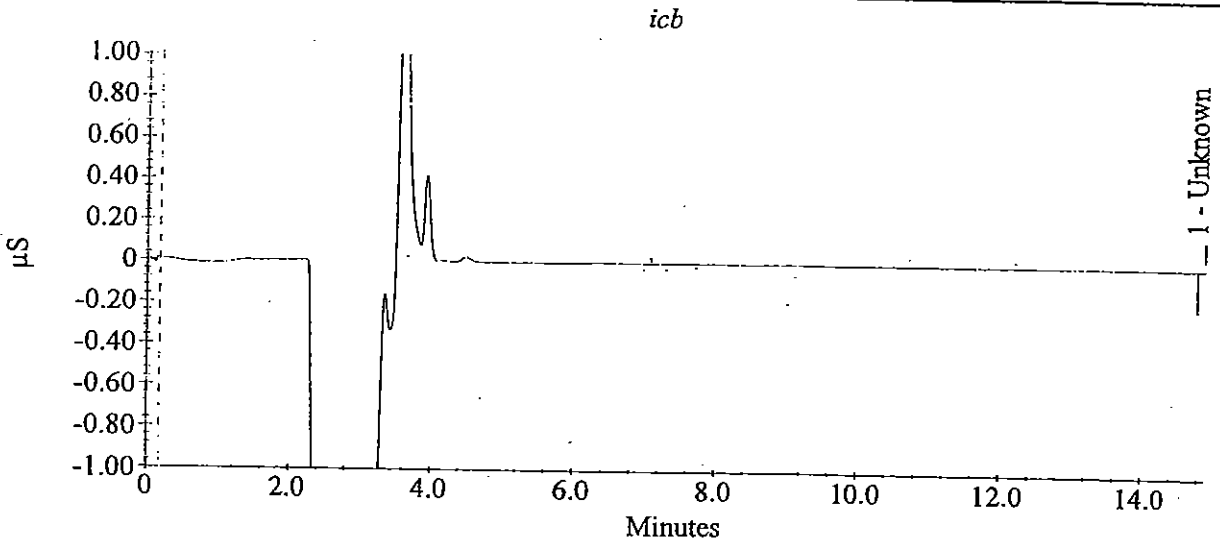
Date Time Collected : 03/12/2003 8:33:51 PM

System Operator : wei wang

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
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APCL Perchlorate Analysis Report

Sample Name : ICCS 4ppb w7827b

Data File Name : C:\DATA\03W2496K\W2496K ICCS 4PPB_006.DXD

Method File Name : c:\peaknet\method\314-011.met

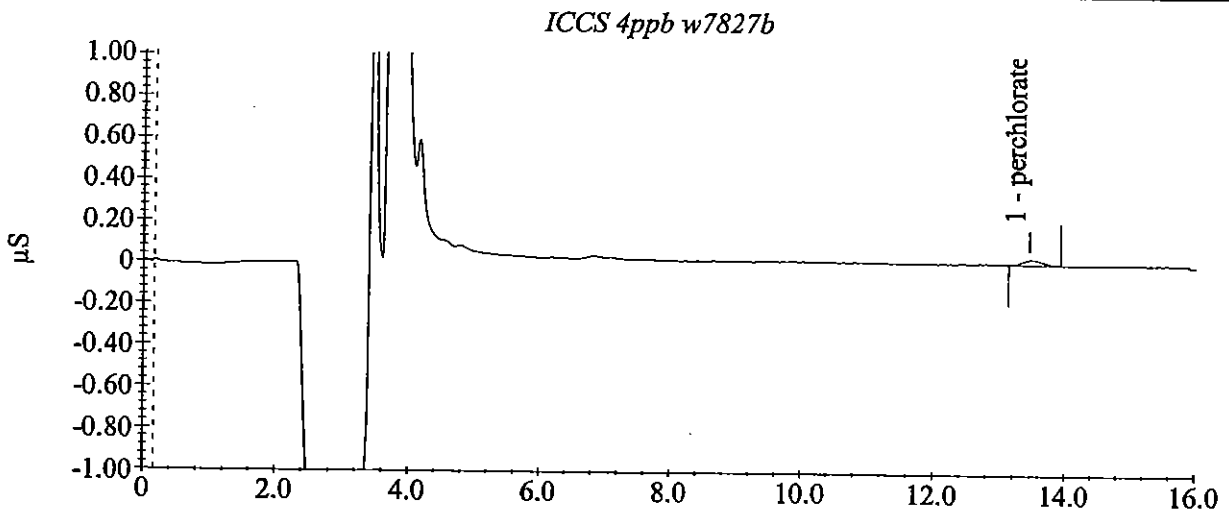
Date Time Collected : 04/22/2003 7:26:27 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.45	3.27	5550.15	242.10



APCL Perchlorate Analysis Report

Sample Name : ##03w2496kw ipc 25ppb w7759

Data File Name : C:\DATA\03W2496K\W2496K IPC 25PPB_001.DXD

Method File Name : c:\peaknet\method\ve314-011.met

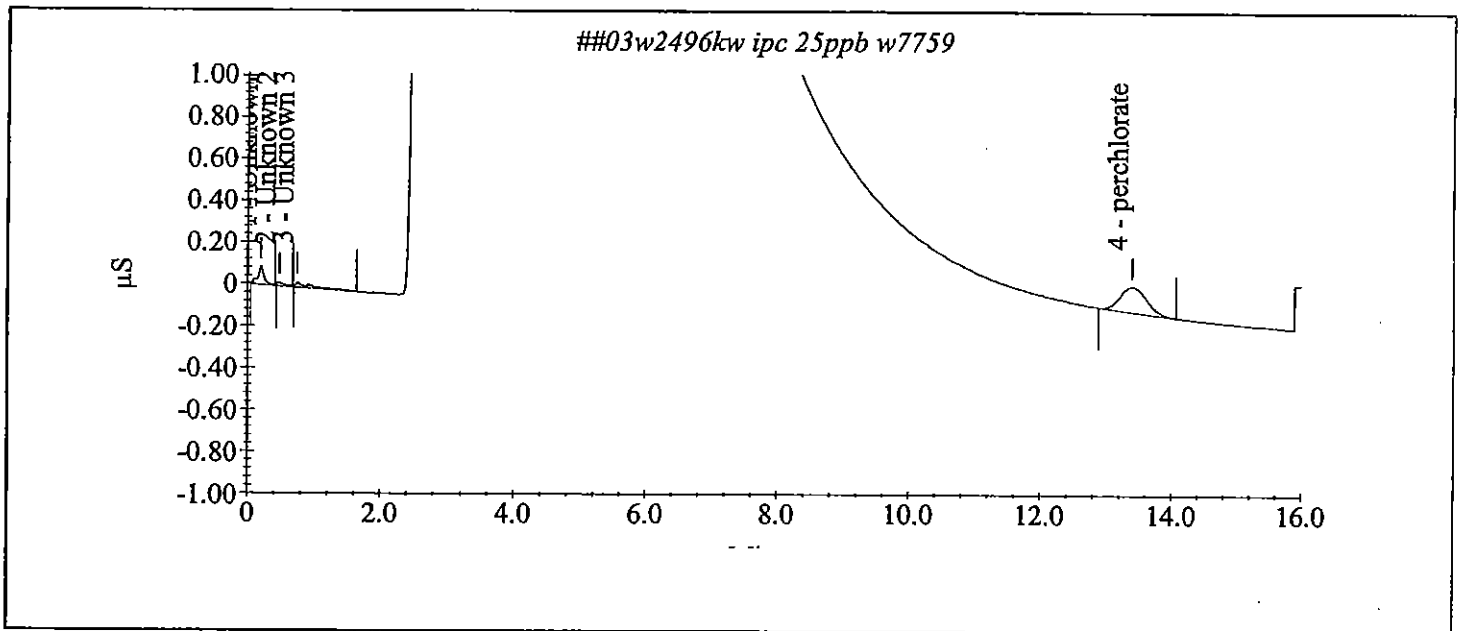
Date Time Collected : 04/22/2003 4:43:22 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
4	perchlorate	13.40	20.01	33963.15	1197.81



APCL Perchlorate Analysis Report

Sample Name : CCB

Data File Name : C:\DATA\03W2496K\W2496K K03_025.DXD

Method File Name : c:\peaknet\method\314-011.met

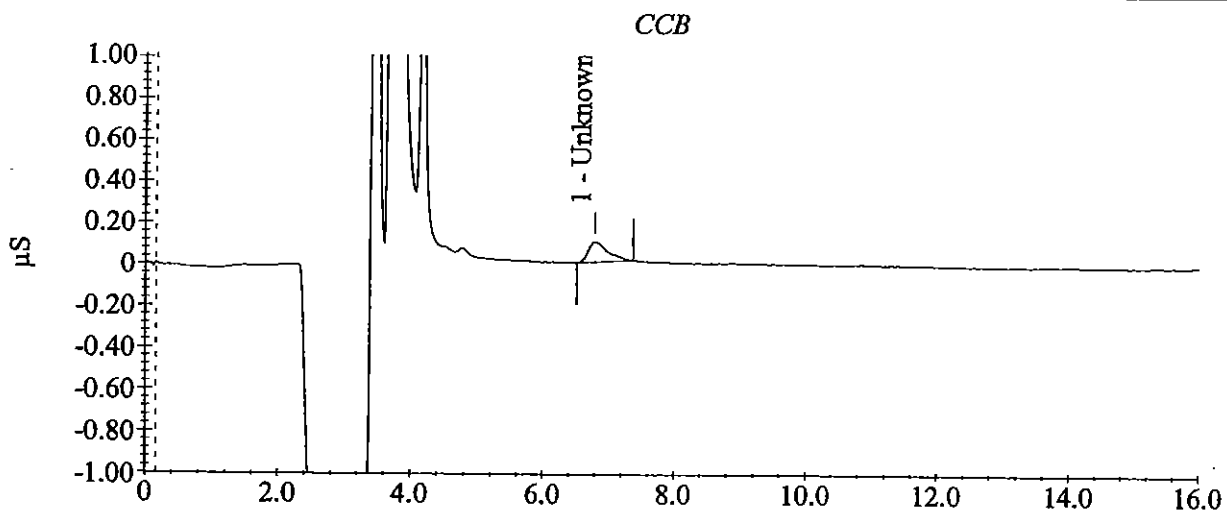
Date Time Collected : 04/23/2003 1:20:02 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
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APCL Perchlorate Analysis Report

Sample Name : CCB

Data File Name : C:\DATA\03W2496K\W2496K K04_033.DXD

Method File Name : c:\peaknet\method\314-011.met

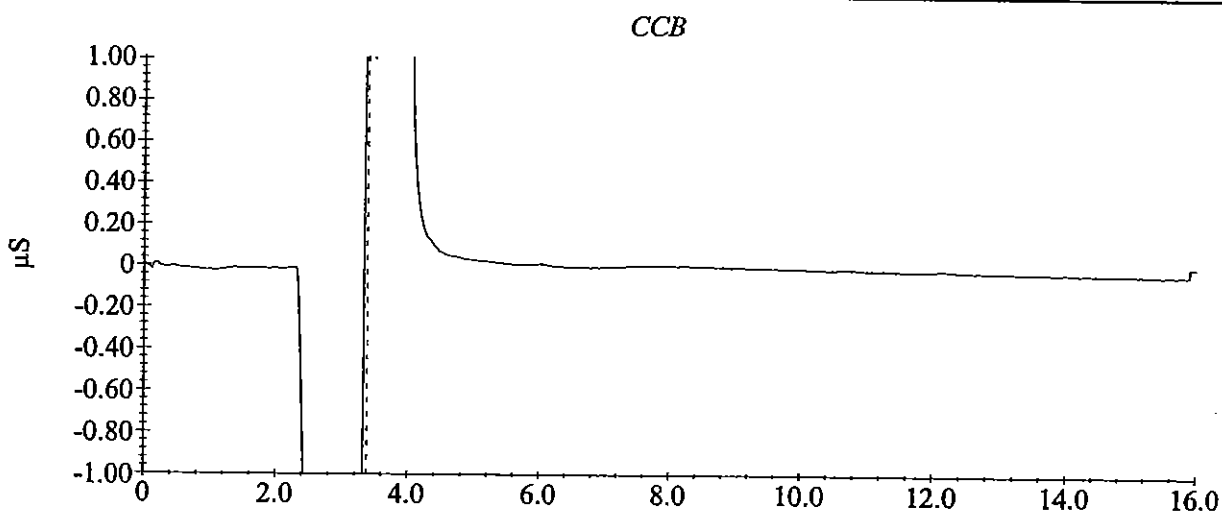
Date Time Collected : 04/23/2003 10:36:23 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
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APCL Perchlorate Analysis Report

Sample Name : ccb

Data File Name : C:\DATA\03W2496K\W2496K CCB01_003.DXD

Method File Name : c:\peaknet\method\314-011.met

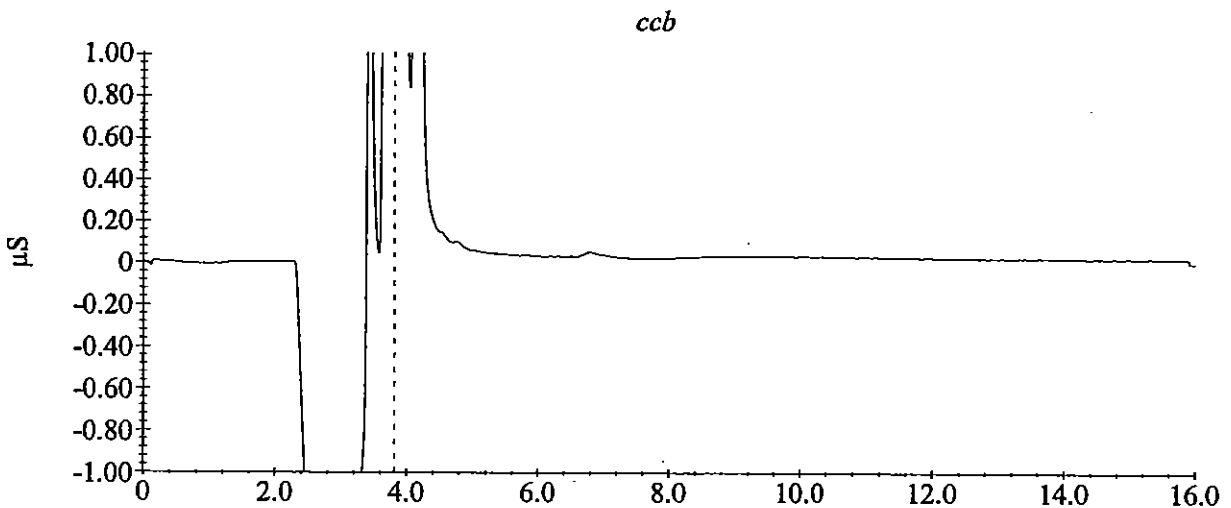
Date Time Collected : 04/22/2003 6:30:26 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
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APCL Perchlorate Analysis Report

Sample Name : ccb

Data File Name : C:\DATA\03W2496K\W2496K K02_014.DXD

Method File Name : c:\peaknet\method\314-011.met

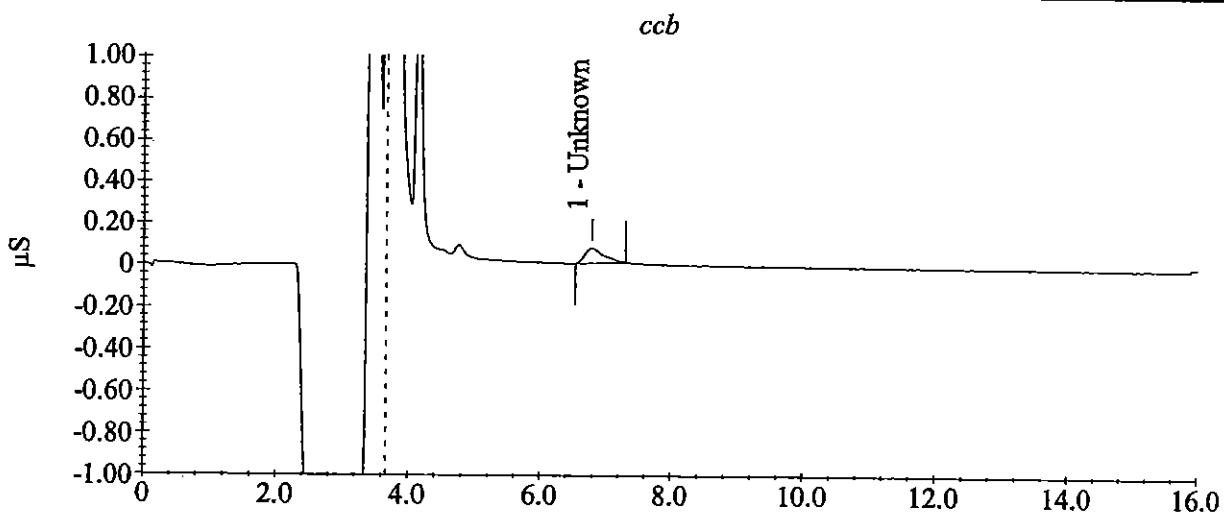
Date Time Collected : 04/22/2003 9:55:23 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

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

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

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

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

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

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

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

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

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

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

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

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

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

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

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

Component: Fluoride

Fit Type: Linear

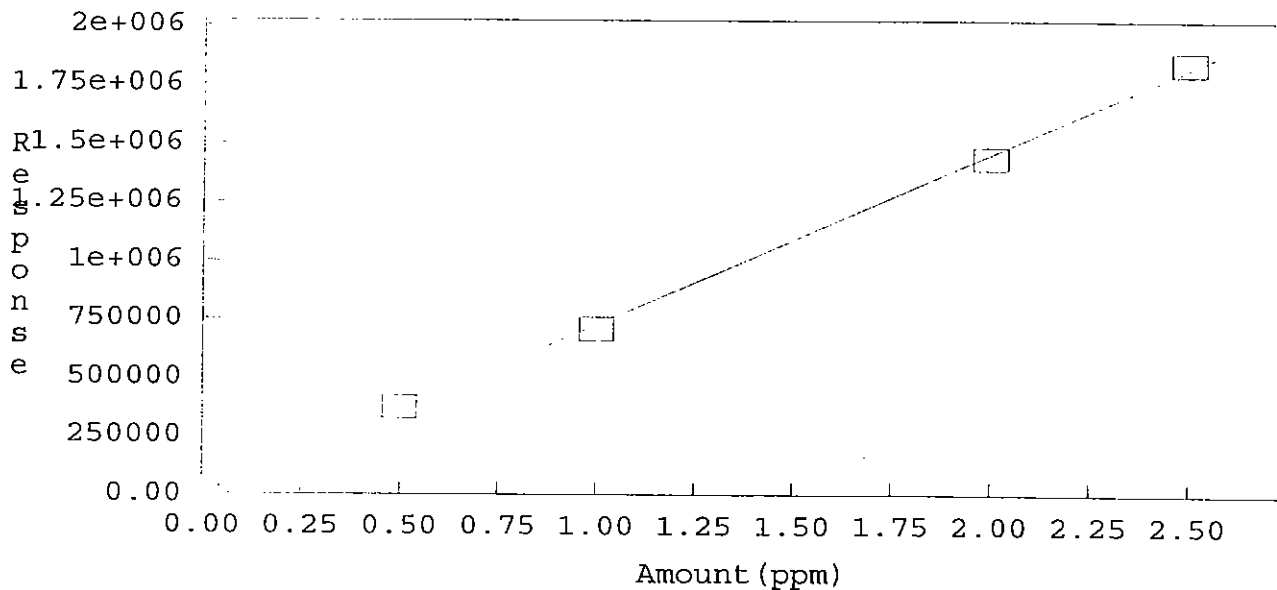
$r^2 = 0.999552$

$Amt = Resp * 1.376e-006 + -0.000962$

$Resp = Amt * 7.267e+005 + 699.7$

Standardization: External

Calibration: Area



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

Component: Chloride

Fit Type: Linear

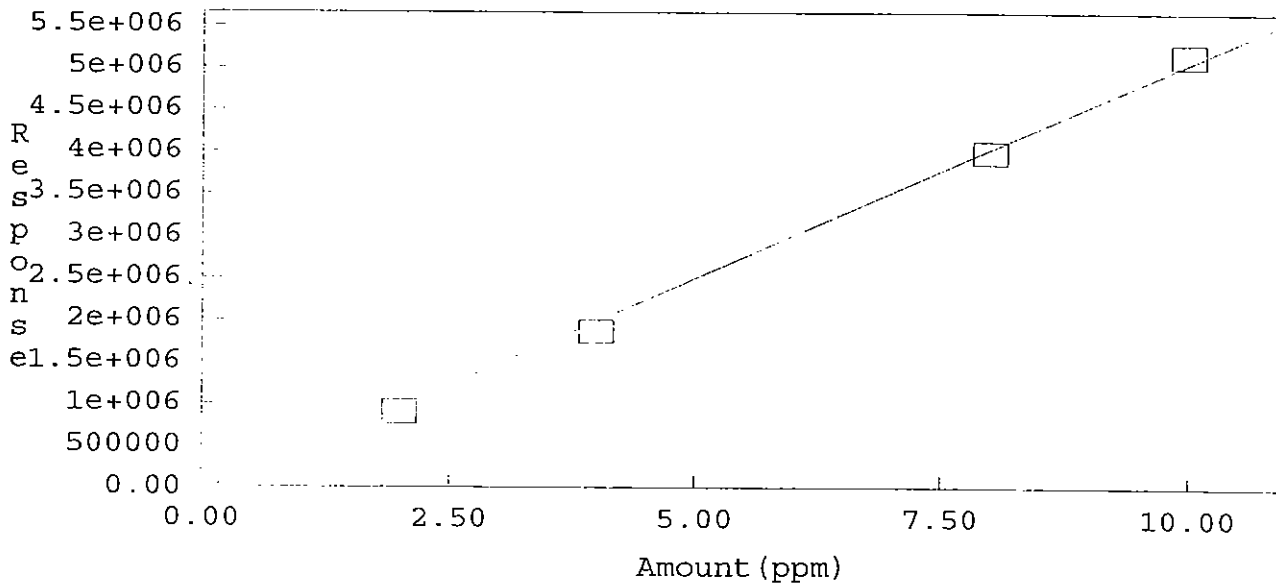
$r^2 = 0.998409$

$Amt = Resp * 1.953e-006 + 0.1282$

$Resp = Amt * 5.121e+005 + -6.564e+00$

Standardization: External

Calibration: Area



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

Component: Nitrite-N

Fit Type: Linear

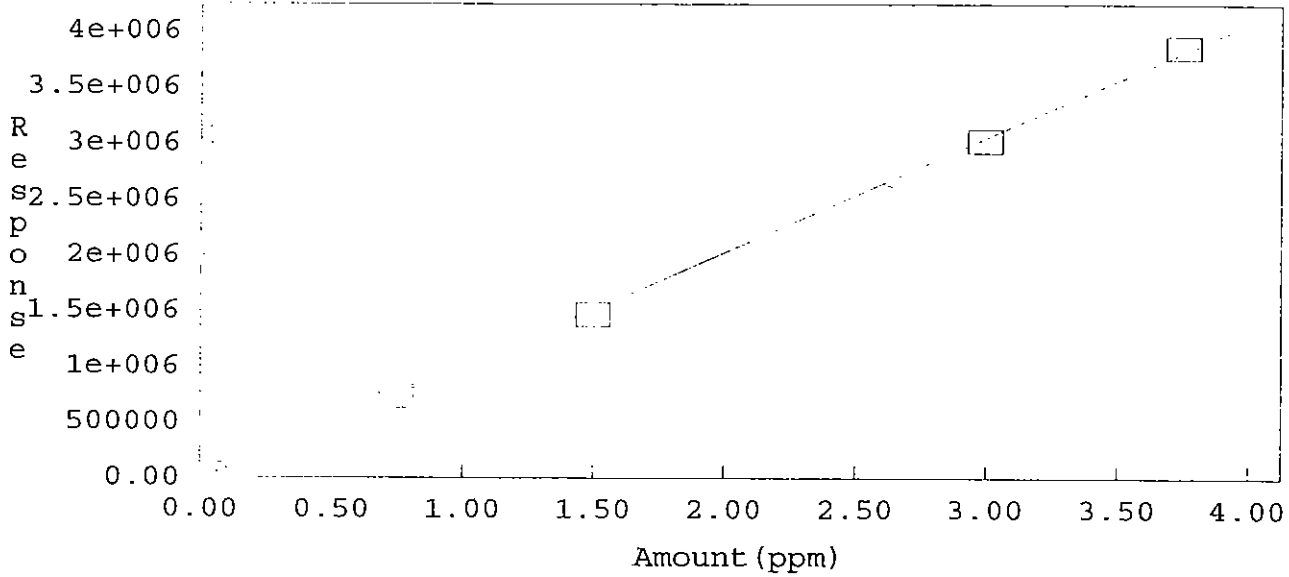
$r^2 = 0.999594$

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

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

Standardization: External

Calibration: Area



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

Component: Bromide

Fit Type: Linear

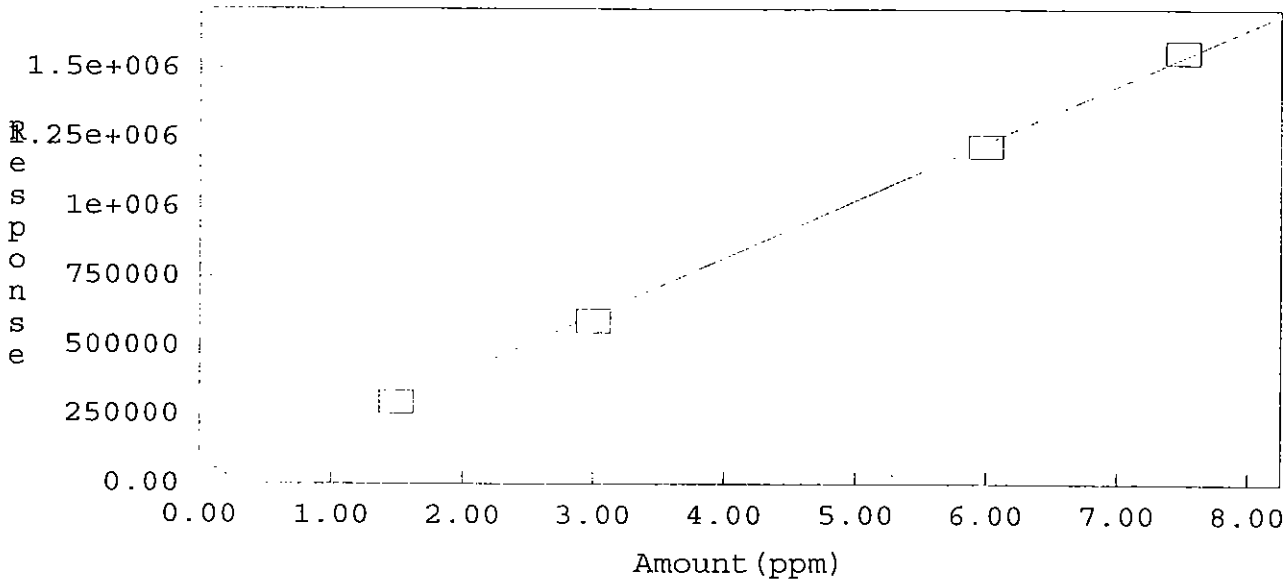
$r^2 = 0.999518$

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

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

Standardization: External

Calibration: Area



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

Component: Nitrate-N

Fit Type: Linear

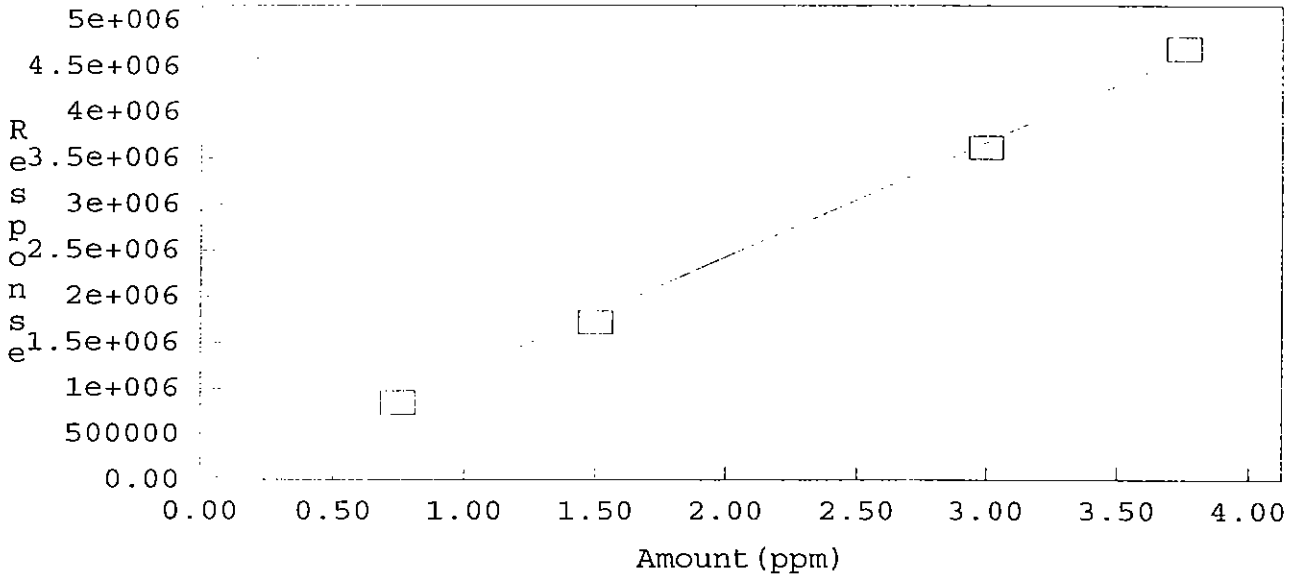
$r^2 = 0.998618$

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

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

Standardization: External

Calibration: Area



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

Component: Phosphate-P

Fit Type: Linear

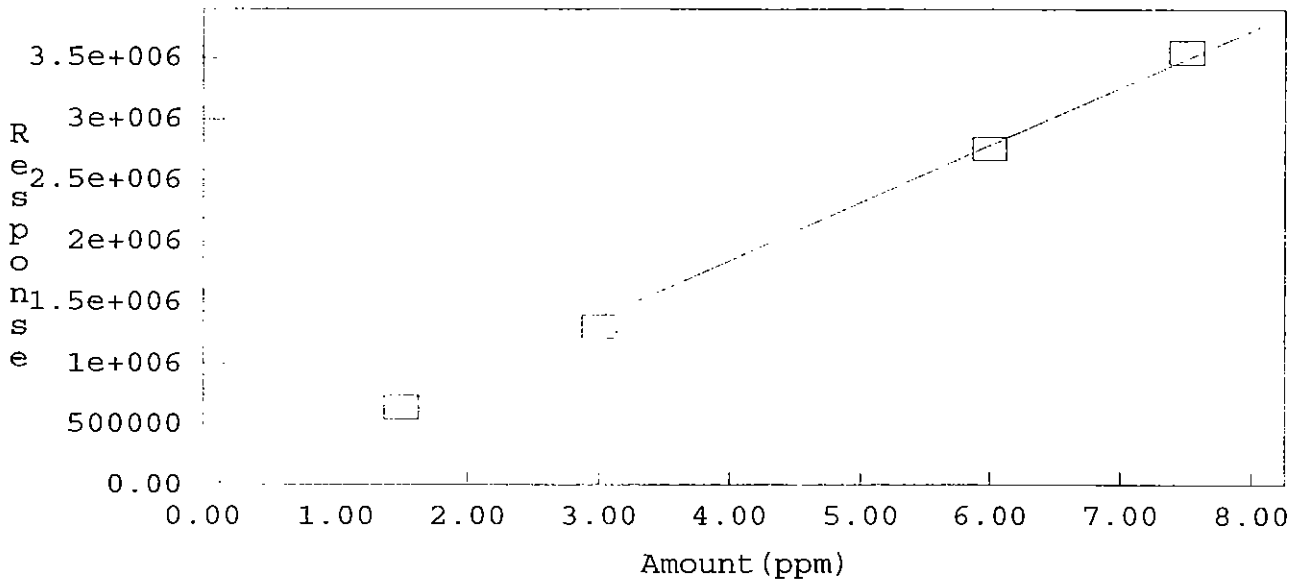
$r^2 = 0.998898$

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

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

Standardization: External

Calibration: Area



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

Component: Sulfate

Fit Type: Linear

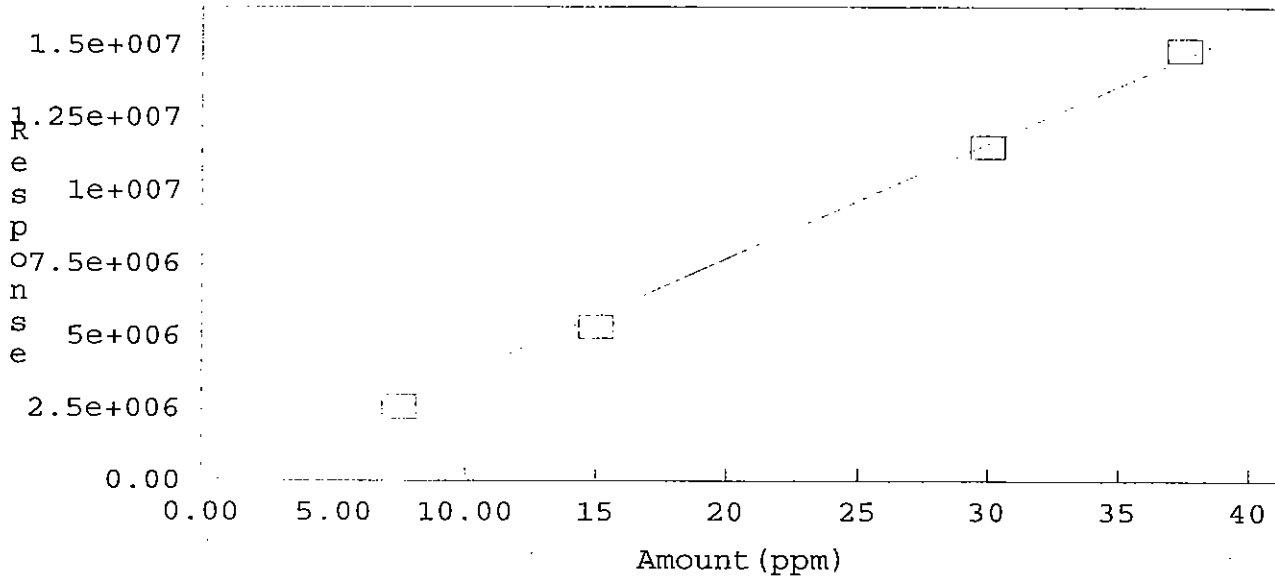
$r^2 = 0.998245$

$Amt = Resp * 2.533e-006 + 0.5323$

$Resp = Amt * 3.949e+005 + -2.102e+00$

Standardization: External

Calibration: Area



```

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

```

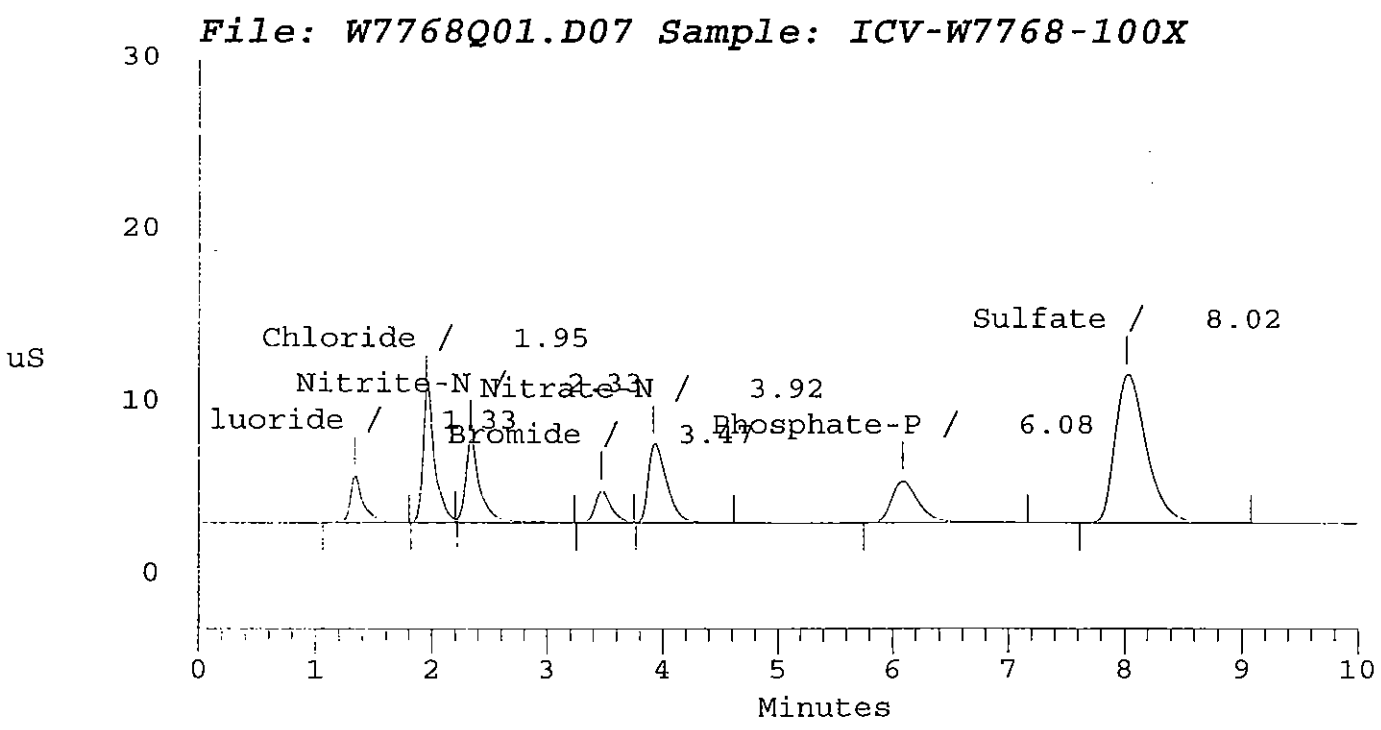
```

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

```

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

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



```

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

```

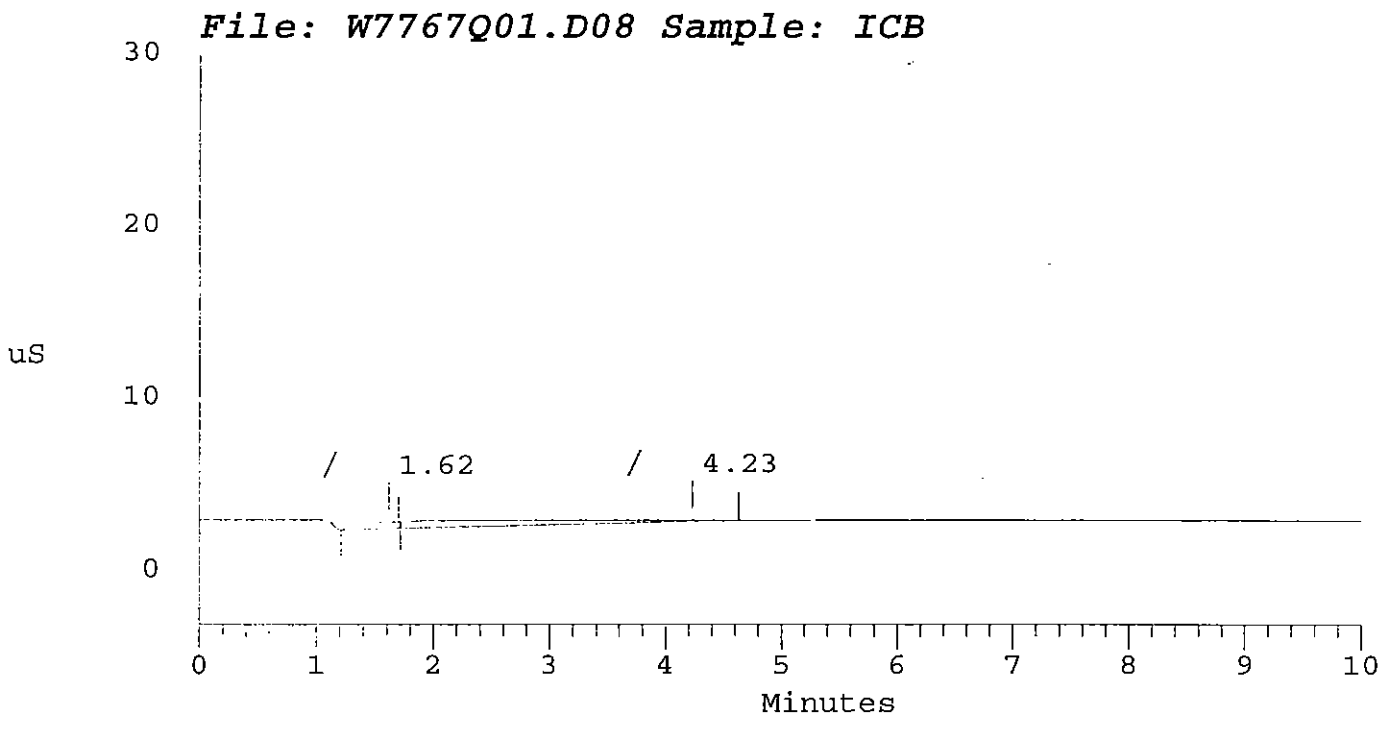
```

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

```

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

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




```

=====
Sample Name: AUTOCAL1R                               Date: 03/21/2003 16:23:08
Data File  : C:\DX\DATA\E300-063\W7767Q01.D01
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 1                 Detector: COND
Analyst    : David                                   Column: Dionex AS4A-SC
=====

```

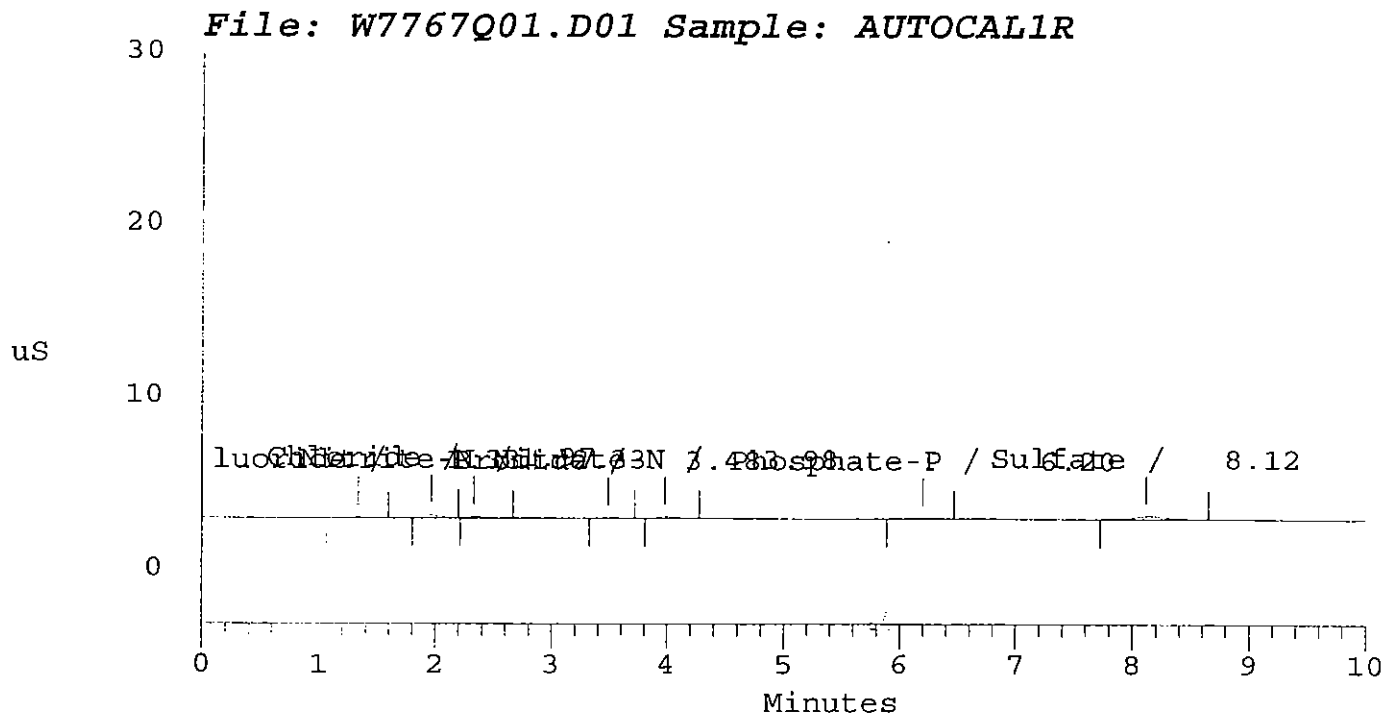
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.33	Fluoride	0.025	2732	28534	1	0.00
2	1.97	Chloride	0.100	7044	51206	2	0.00
3	2.33	Nitrite-N	0.038	3582	30884	2	-3.33
4	3.48	Bromide	0.075	1488	13830	1	-1.47
5	3.98	Nitrate-N	0.038	3802	40157	1	0.00
6	6.20	Phosphate-P	0.075	1450	24783	1	0.00
7	8.12	Sulfate	0.376	6524	129999	1	0.00
Totals			0.726	26621	319393		



```

=====
Sample Name: AUTOCAL2R                               Date: 03/21/2003 16:35:53
Data File  : C:\DX\DATA\E300-063\W7767Q01.D02
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 2                 Detector:COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====

```

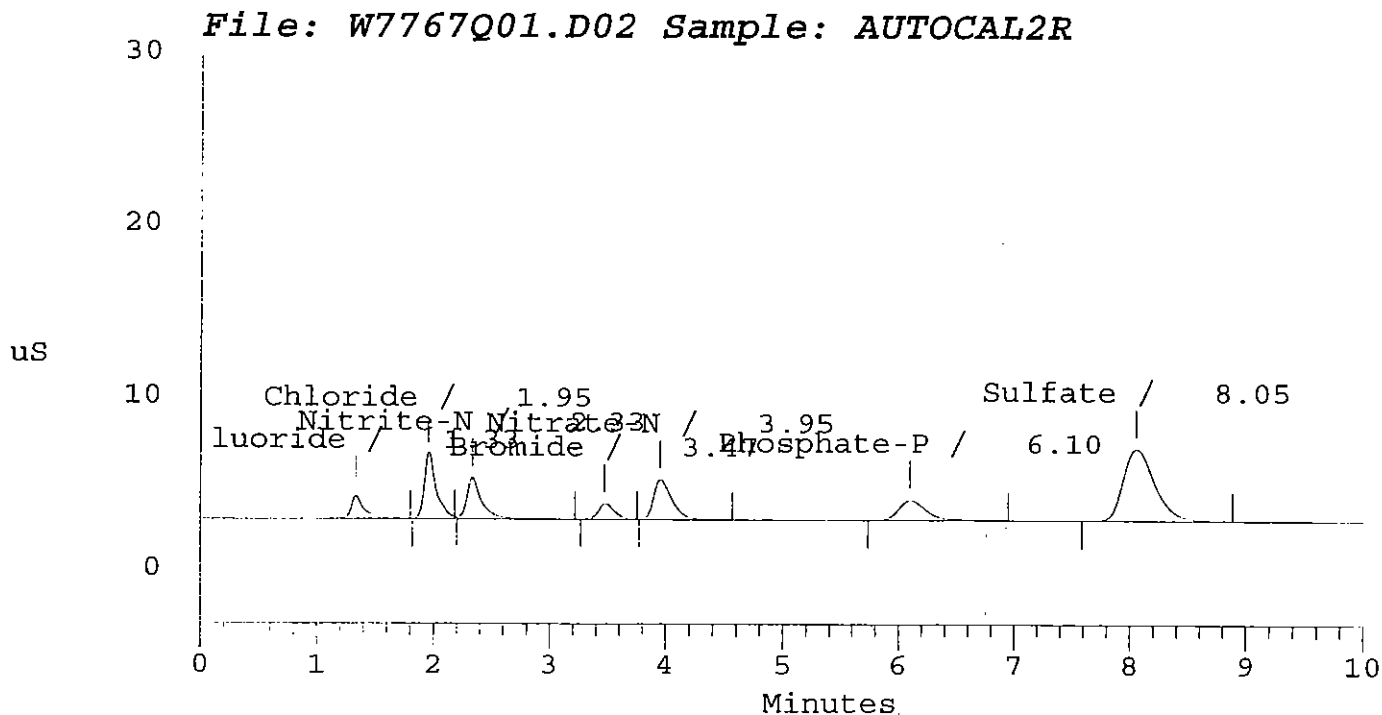
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.33	Fluoride	0.500	44629	373164	2	0.00
2	1.95	Chloride	2.000	126181	909455	2	0.00
3	2.33	Nitrite-N	0.750	79701	734006	2	0.85
4	3.47	Bromide	1.500	30100	298206	2	0.36
5	3.95	Nitrate-N	0.750	77129	849179	2	0.00
6	6.10	Phosphate-P	1.500	38579	642376	1	0.00
7	8.05	Sulfate	7.500	138579	2598757	1	0.00
Totals			14.500	534896	6405142		



```

=====
Sample Name: AUTOCAL3R                               Date: 03/21/2003 16:48:37
Data File  : C:\DX\DATA\E300-063\W7767Q01.D03
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 3                 Detector:COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====

```

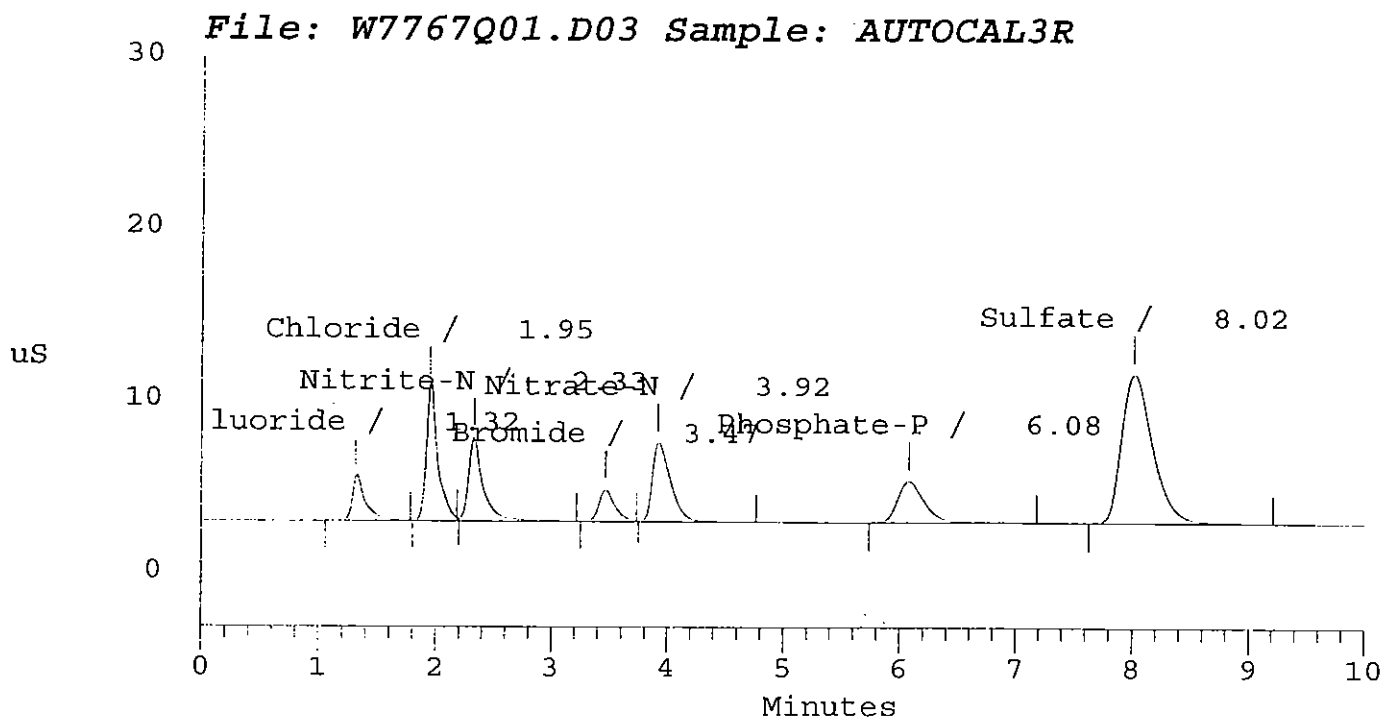
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	1.000	82595	707646	2	0.00
2	1.95	Chloride	4.000	261681	1856586	2	0.00
3	2.33	Nitrite-N	1.500	162005	1468106	2	0.00
4	3.47	Bromide	3.000	61776	591234	2	0.85
5	3.92	Nitrate-N	1.500	152776	1713421	2	0.00
6	6.08	Phosphate-P	3.000	79971	1301126	1	0.00
7	8.02	Sulfate	15.000	287851	5330209	1	0.00
Totals			29.000	1088657	12968327		



```

=====
Sample Name: AUTOCAL4R                               Date: 03/21/2003 17:01:21
Data File  : C:\DX\DATA\E300-063\W7767Q01.D04
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 4                 Detector: COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====

```

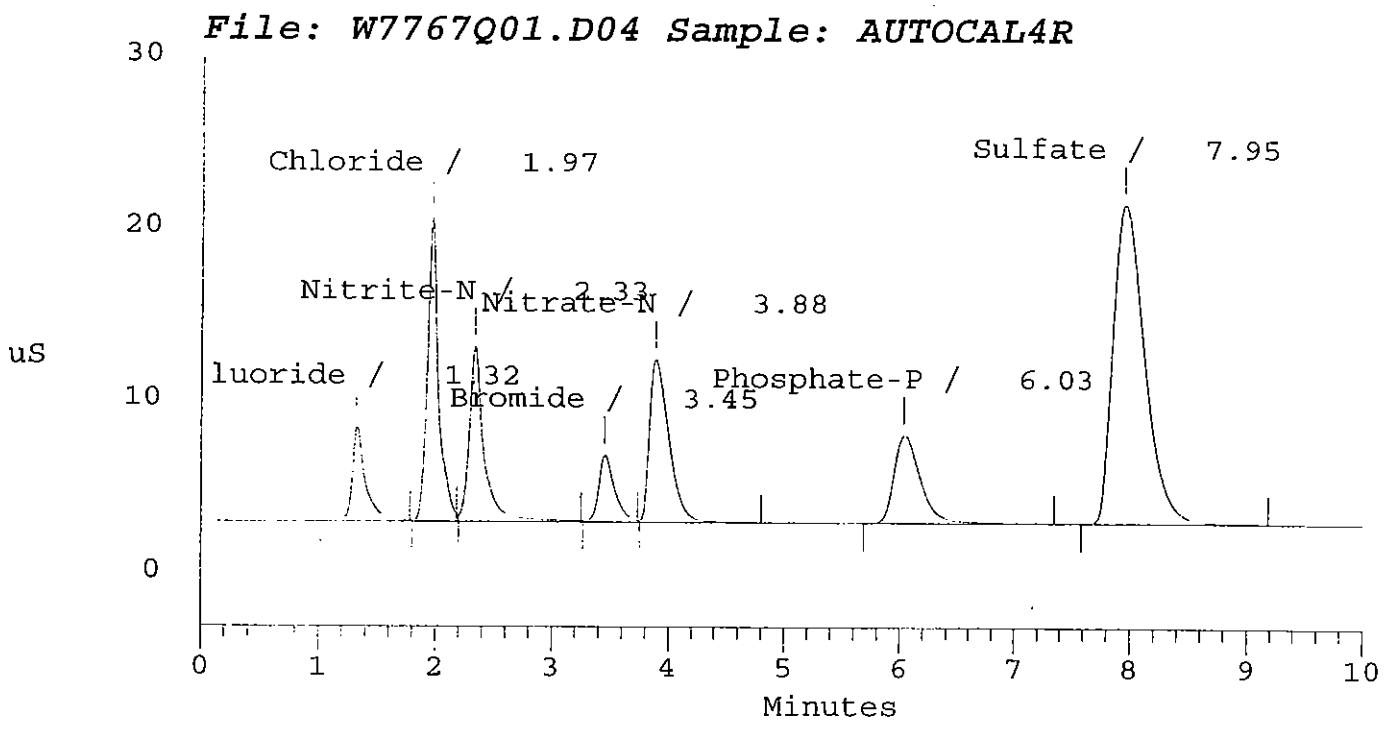
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	2.000	173007	1435865	2	0.00
2	1.97	Chloride	8.000	585791	3987563	2	0.00
3	2.33	Nitrite-N	3.000	336616	3021523	2	-0.85
4	3.45	Bromide	6.000	128845	1219933	2	0.37
5	3.88	Nitrate-N	3.000	313707	3610927	2	0.00
6	6.03	Phosphate-P	6.000	168994	2756481	2	0.00
7	7.95	Sulfate	30.000	615917	11507107	2	0.00
Totals			58.000	2322876	27539399		



```

=====
Sample Name: AUTOCAL5R                               Date: 03/21/2003 17:14:05
Data File  : C:\DX\DATA\E300-063\W7767Q01.D05
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 5                 Detector:COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====

```

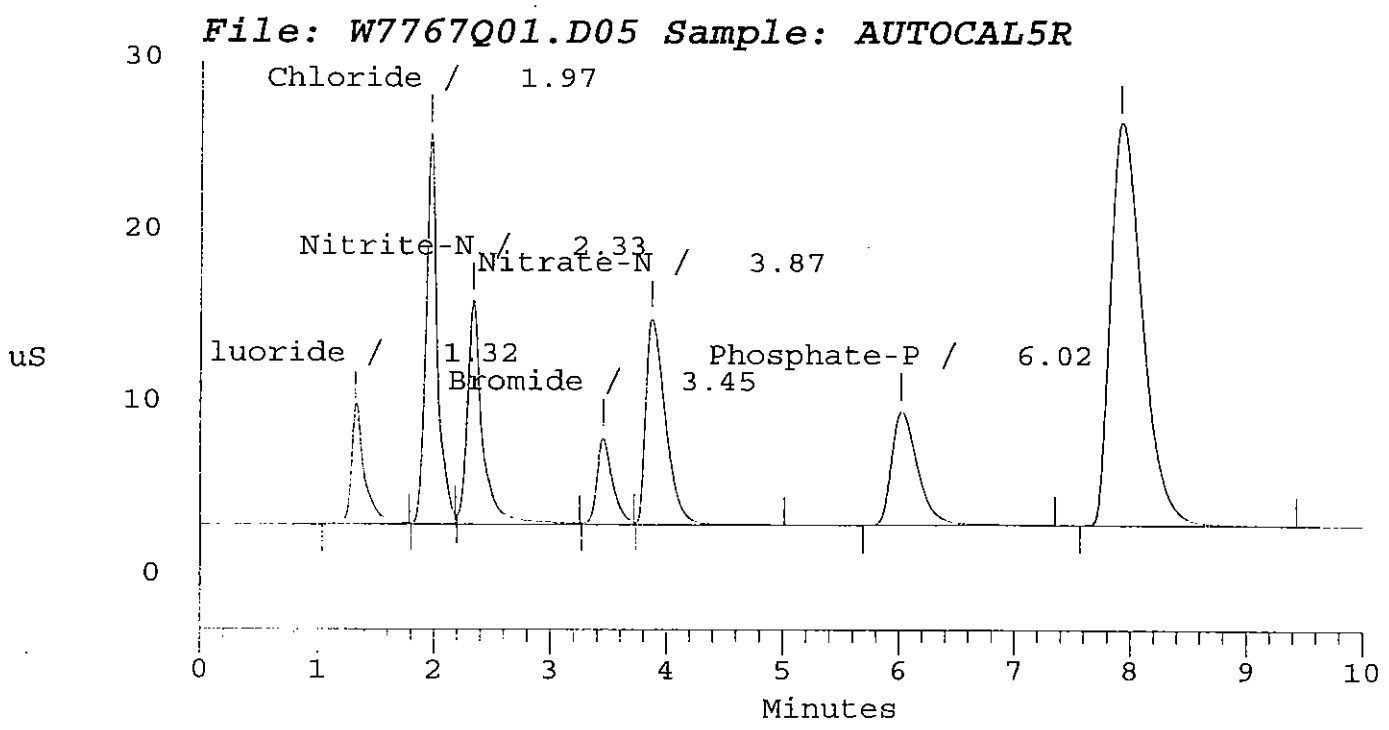
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	2.500	220914	1837162	2	0.00
2	1.97	Chloride	10.000	754321	5142155	2	0.00
3	2.33	Nitrite-N	3.750	429219	3856614	2	0.00
4	3.45	Bromide	7.500	166594	1559887	2	0.43
5	3.87	Nitrate-N	3.750	396441	4688990	2	0.00
6	6.02	Phosphate-P	7.500	217521	3546397	2	0.00
7	7.92	Sulfate	37.500	776426	14859049	2	0.00
Totals			72.500	2961435	35490255		



Data Reprocessed On 03/21/2003 17:39:26

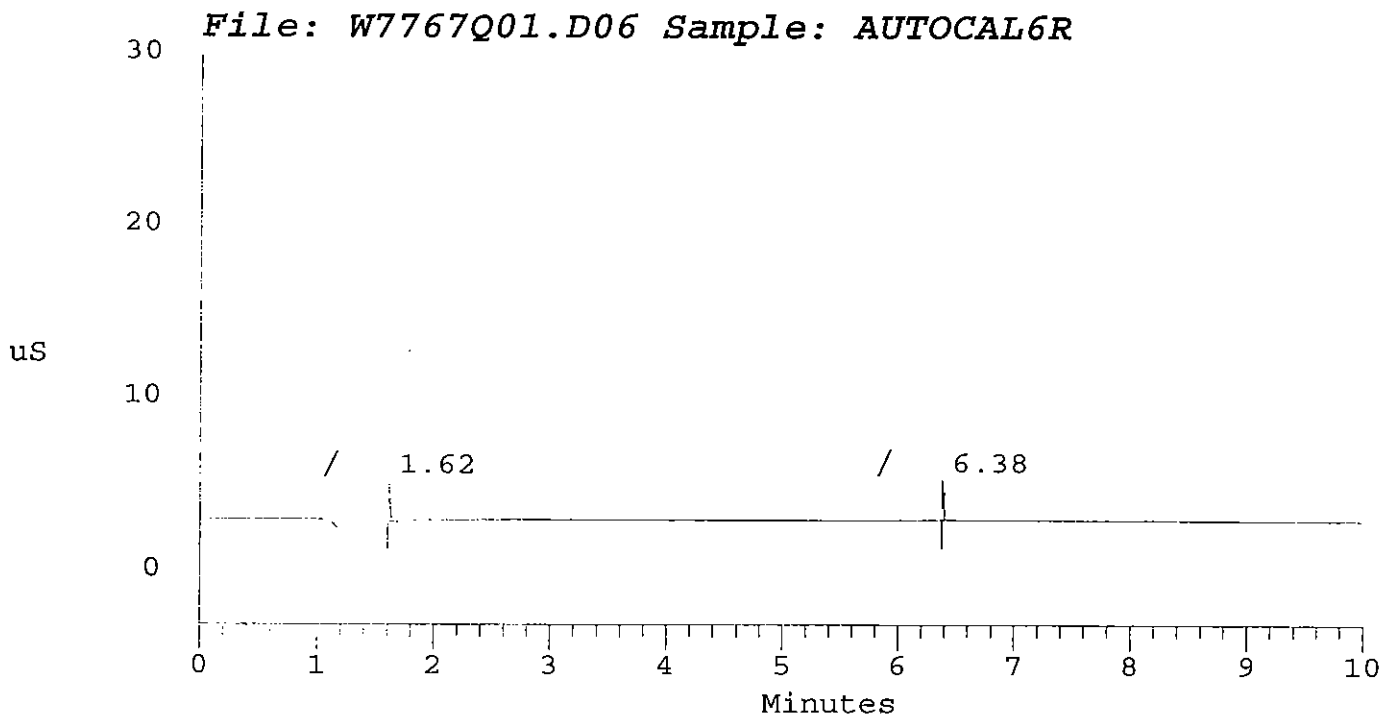
=====
Sample Name: AUTOCAL6R Date: 03/21/2003 17:37:58
Data File : C:\DX\DATA\E300-063\W7767Q01.D06
Method : C:\DX\METHOD\E300-063.MET
AGI Address: 1 System: 1 Inject#: 6 Detector: COND
Analyst : David Column: Dionex AS4A-SC
=====

Calibration Volume Dilution Points Rate Start Stop Area Reject

External 1 1 3000 5Hz 0.00 10.00 1000

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

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



After
3/21/03
V... 27

```

=====
Sample Name: AUTOCAL6R                               Date: 03/21/2003 17:37:58
Data File  : C:\DX\DATA\E300-063\W7767Q01.D06
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 6                 Detector:COND
Analyst    : David                                  Column: Dionex AS4A-SC
=====

```

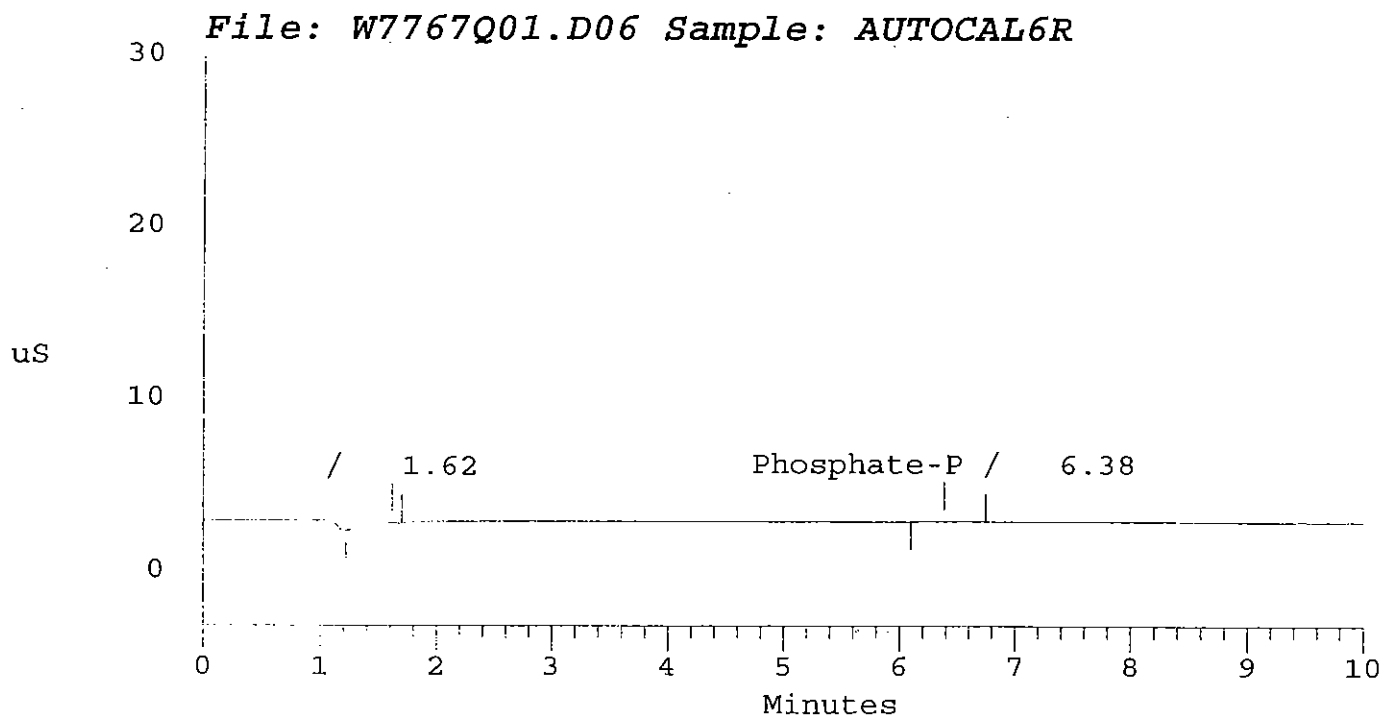
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
0	0.00	Fluoride	0.000	0	0	0	0.00
0	0.00	Chloride	0.000	0	0	0	0.00
0	0.00	Nitrite-N	0.000	0	0	0	0.00
0	0.00	Bromide	0.000	0	0	0	0.00
0	0.00	Nitrate-N	0.000	0	0	0	0.00
2	6.38	Phosphate-P	0.000	582	11752	1	0.00
0	0.00	Sulfate	0.000	0	0	0	0.00
Totals			0.000	582	11752		



Before

6A
INITIAL CALIBRATION DATA

Lab Name: Applied P & Ch Lab Contract: _____

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

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

A = 0.000 + 0.836C

A = Absorbance

C = Concentration (mg/L)

r = 0.9997

Wet Chemistry QC Report B
Duplicate Results

Matrix: Water

APCL Service ID: 03-2809

Analysis	Batch ID	Analysis Date	Sample Name	Unit	Result	Duplicate Result	RPD %	RPD Control limit
Alkalinity	03W2555	04/24/2003	DUPE-1-2Q03	mg/L	161.0	158.4	2	20
pH	03W2477	04/21/2003	MW-4-1	pH unit	7.05	7.08	0	20

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

FORM-7

Applied P & Ch Laboratory

CCV Recovery for Wet Analysis

Client Name: GEOPON, Inc.

Contract No.:

Lab Code:

APCL

Case No:

SAS No.:

Service ID:

32809

Project ID: JPL

Project No.: 04-4428.10

#	Component Name	Method	Batch No.	Unit	Expected	Test Result	Rec. %	Dev. %	Flag	Control Limit, %	Test Date
1	Chloride Cl ⁻	300.0	03W2479	mg/L	4.0	3.89	97	-3	✓	90-110	04/22/2003
	NITRATE as N-NO ₃ ⁻ , BY	300.0	03W2479	mg/L	1.5	1.50	100	0	✓	90-110	04/22/2003
	SULFATE SO ₄ ⁻ , BY I	300.0	03W2479	mg/L	15	14.6	97	-3	✓	90-110	04/22/2003
	Chloride Cl ⁻	300.0	03W2479	mg/L	4.0	3.92	98	-2	✓	90-110	04/22/2003
	NITRATE as N-NO ₃ ⁻ , BY	300.0	03W2479	mg/L	1.5	1.52	101	1	✓	90-110	04/22/2003
	SULFATE SO ₄ ⁻ , BY I	300.0	03W2479	mg/L	15	14.7	98	-2	✓	90-110	04/22/2003
	Chloride Cl ⁻	300.0	03W2479	mg/L	4.0	3.97	99	-1	✓	90-110	04/22/2003
	NITRATE as N-NO ₃ ⁻ , BY	300.0	03W2479	mg/L	1.5	1.53	102	2	✓	90-110	04/22/2003
	SULFATE SO ₄ ⁻ , BY I	300.0	03W2479	mg/L	15	14.8	99	-1	✓	90-110	04/22/2003
	Chloride Cl ⁻	300.0	03W2479	mg/L	4.0	3.92	98	-2	✓	90-110	04/22/2003
	NITRATE as N-NO ₃ ⁻ , BY	300.0	03W2479	mg/L	1.5	1.52	101	1	✓	90-110	04/22/2003
	SULFATE SO ₄ ⁻ , BY I	300.0	03W2479	mg/L	15	14.7	98	-2	✓	90-110	04/22/2003
2	Perchlorate	314.0	03W2496	μg/L	50	52.4	105	5	✓	85-115	04/22/2003
	Perchlorate	314.0	03W2496	μg/L	50	53.9	108	8	✓	85-115	04/23/2003
	Perchlorate	314.0	03W2496	μg/L	50	53.7	107	7	✓	85-115	04/23/2003
	Perchlorate	314.0	03W2496	μg/L	50	48.2	96	-4	✓	85-115	04/23/2003
	Perchlorate	314.0	03W2496	μg/L	50	51.8	104	4	✓	85-115	04/23/2003
3	Chromium (VI)	7196	03W2476	mg/L	0.25	0.252	101	1	✓	90-110	04/21/2003
	Chromium (VI)	7196	03W2476	mg/L	0.25	0.250	100	0	✓	90-110	04/21/2003

```

=====
Sample Name:  ##03W2479, W CCVW7767-100X      Date: 04/22/2003 09:08:43
Data File   :  C:\DX\DATA\03W2479\W2479Q01.D01
Method      :  C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 1           Detector: COND
Analyst     :  David                          Column: Dionex AS4A-SC
=====

```

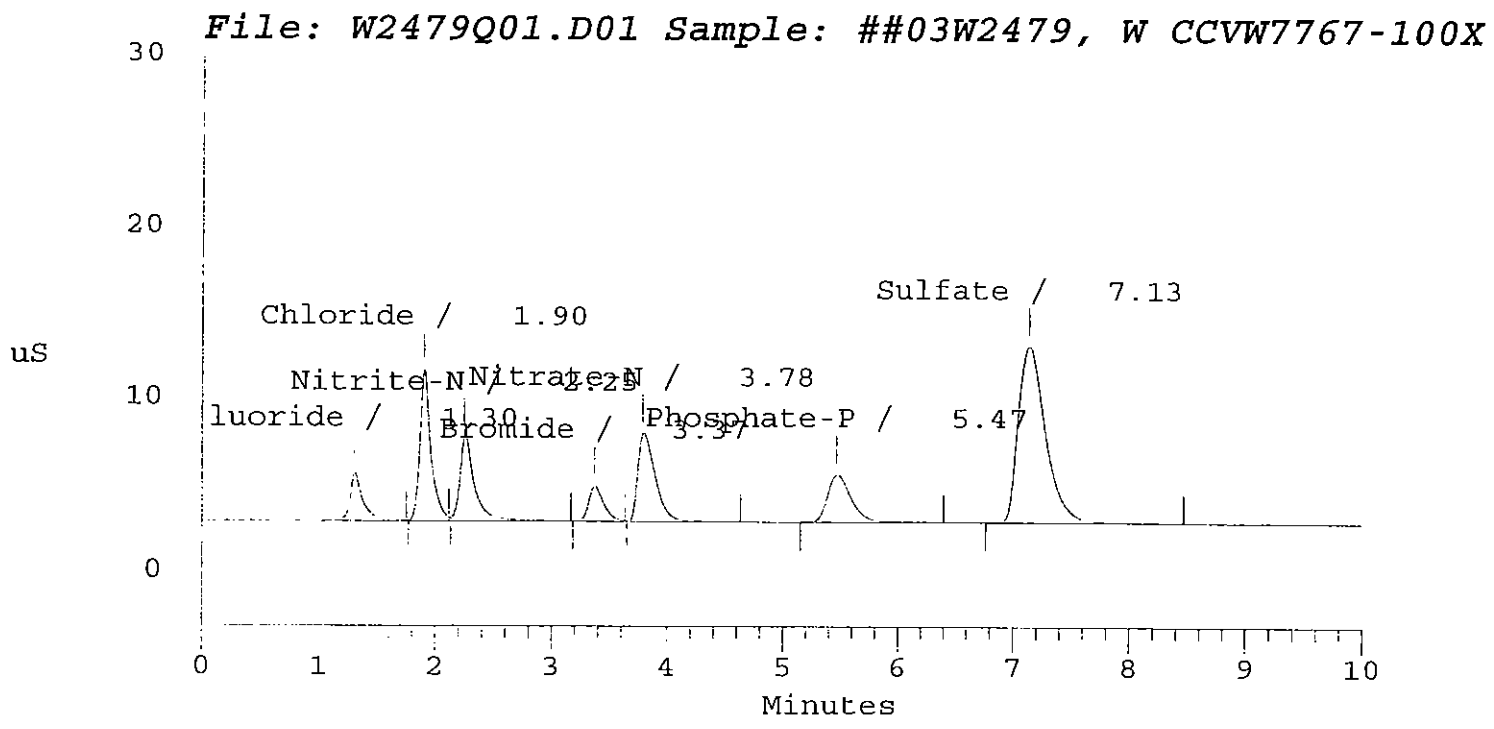
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.30	Fluoride	1.018	89676	740676	2	0.00
2	1.90	Chloride	3.891	286553	1926728	2	0.00
3	2.25	Nitrite-N	1.446	159608	1456884	2	0.00
4	3.37	Bromide	3.016	67928	614566	2	0.48
5	3.78	Nitrate-N	1.498	168870	1806348	2	0.00
6	5.47	Phosphate-P	2.886	91074	1318884	1	0.00
7	7.13	Sulfate	14.603	339530	5556185	1	0.00
Totals			28.358	1203240	13420271		



```

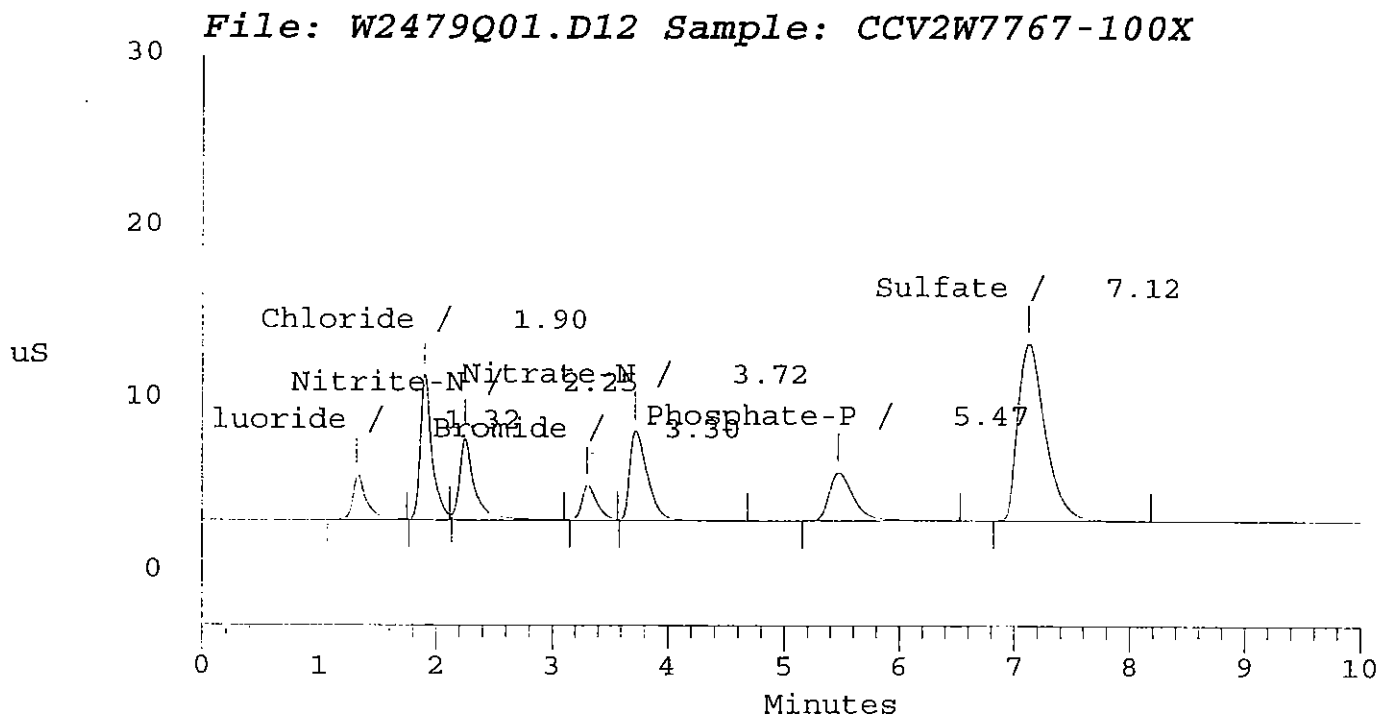
=====
Sample Name: CCV2W7767-100X           Date: 04/22/2003 13:41:40
Data File  : C:\DX\DATA\03W2479\W2479Q01.D12
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 12   Detector: COND
Analyst    : David                    Column: Dionex AS4A-SC
=====
    
```

```

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

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.32	Fluoride	1.015	83929	738138	2	0.00
2	1.90	Chloride	3.922	277552	1942703	2	0.00
3	2.25	Nitrite-N	1.431	156174	1441205	2	0.00
4	3.30	Bromide	3.044	67081	620425	2	0.25
5	3.72	Nitrate-N	1.518	171805	1831322	2	0.00
6	5.47	Phosphate-P	3.050	92318	1396394	1	0.00
7	7.12	Sulfate	14.702	339765	5594980	1	0.00
Totals			28.682	1188622	13565167		



```

=====
Sample Name: CCV3W7767-100X           Date: 04/22/2003 16:01:56
Data File  : C:\DX\DATA\03W2479\W2479Q01.D18
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 18   Detector: COND
Analyst    : David                    Column: Dionex AS4A-SC
=====

```

```

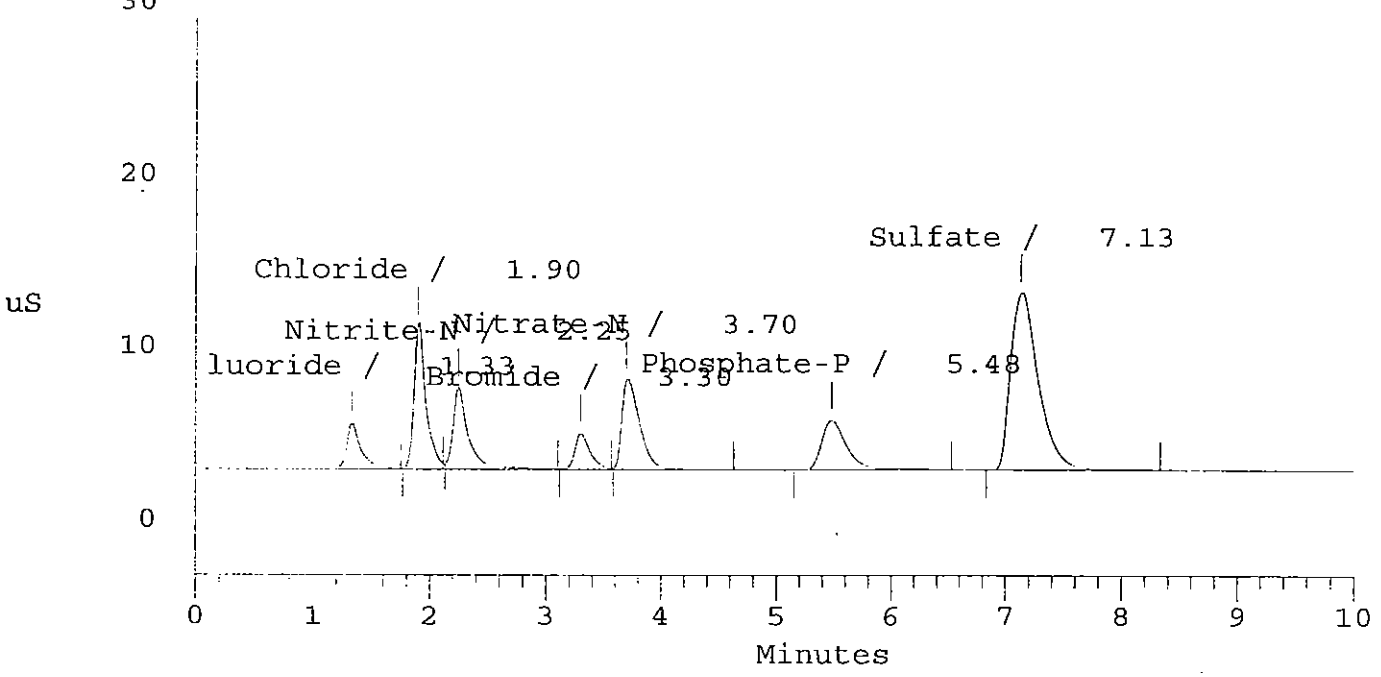
-----
Calibration Volume Dilution Points Rate Start Stop Area Reject
-----
External          1          1 3000 5Hz 0.00 10.00          1000

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.33	Fluoride	1.032	88200	750697	2	0.00
2	1.90	Chloride	3.969	274665	1966649	2	0.00
3	2.25	Nitrite-N	1.450	156672	1461106	2	0.00
4	3.30	Bromide	3.087	68327	629322	2	0.71
5	3.70	Nitrate-N	1.531	169909	1847431	2	0.00
6	5.48	Phosphate-P	3.142	94267	1439318	1	0.00
7	7.13	Sulfate	14.828	339078	5644979	1	0.00
Totals			29.039	1191118	13739501		

File: W2479Q01.D18 Sample: CCV3W7767-100X



```

=====
Sample Name: CCV4W7767-100X           Date: 04/22/2003 18:46:30
Data File  : C:\DX\DATA\03W2479\W2479Q01.D28
Method     : C:\DX\METHOD\E300-063.MET
ACI Address: 1 System: 1 Inject#: 28
Analyst    : David                    Column: Dionex AS4A-SC
Detector: COND
=====

```

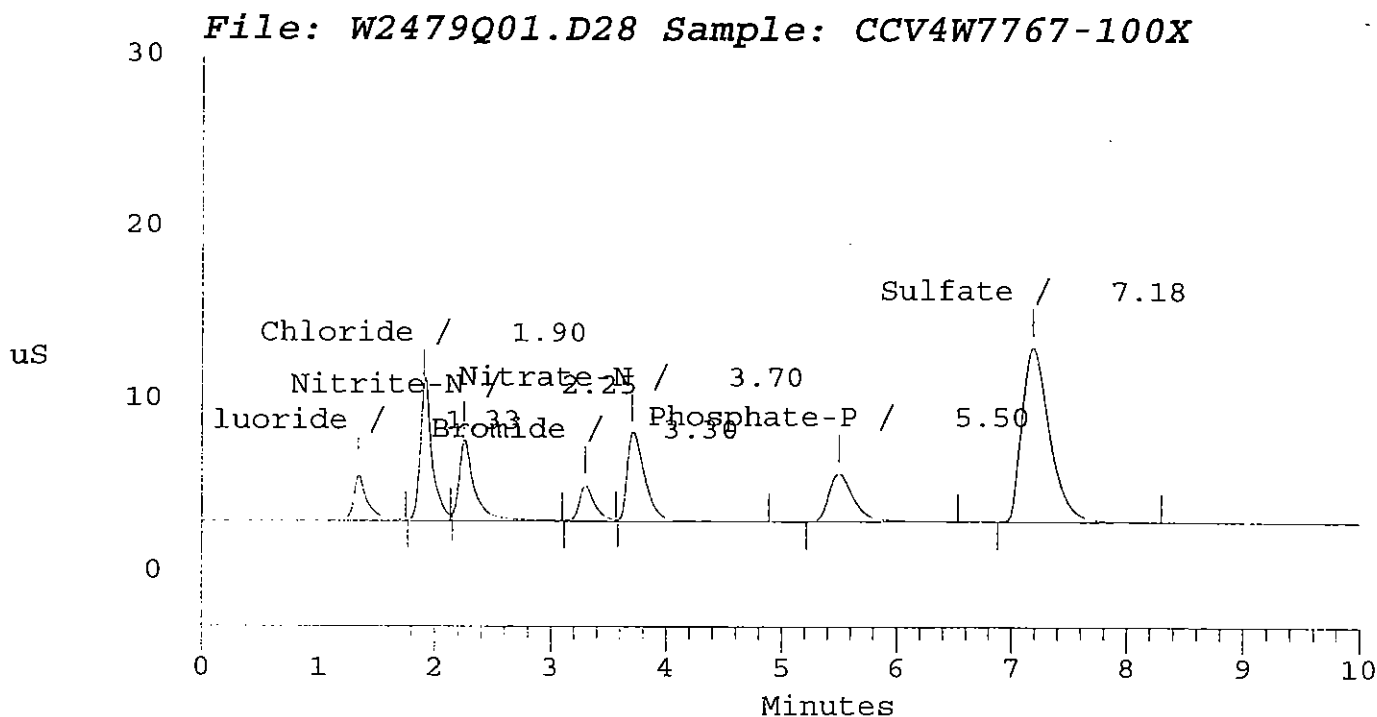
```

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

```

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

Pk. Num	Ret Time	Component Name	Concentration ppm	Height	Area	Bl. Code	%Delta
1	1.33	Fluoride	1.005	84714	730909	2	0.00
2	1.90	Chloride	3.915	253791	1938964	2	0.00
3	2.25	Nitrite-N	1.441	153836	1452033	2	0.00
4	3.30	Bromide	3.078	67663	627394	2	0.71
5	3.70	Nitrate-N	1.517	168718	1830118	2	0.00
6	5.50	Phosphate-P	3.078	92353	1409370	1	0.00
7	7.18	Sulfate	14.706	335912	5596807	1	0.00
Totals			28.740	1156988	13585594		



APCL Perchlorate Analysis Report

Sample Name : ccv 50ppb w7827e

Data File Name : C:\DATA\03W2496K\W2496K Q02_013.DXD

Method File Name : c:\peaknet\method\314-011.met

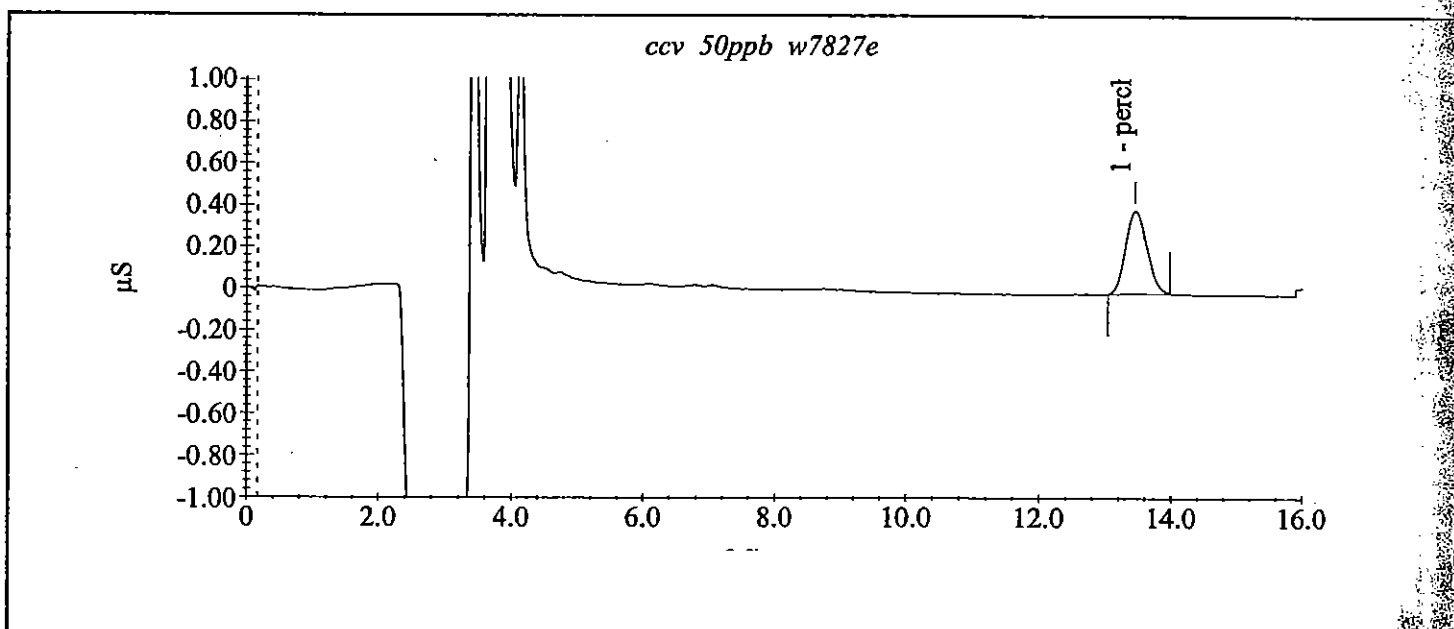
Date Time Collected : 04/22/2003 9:36:32 PM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.45	52.42	88964.40	3967.48



APCL Perchlorate Analysis Report

Sample Name : ccv 50ppb w7827e

Data File Name : C:\DATA\03W2496K\W2496K Q03_024.DXD

Method File Name : c:\peaknet\method\e314-011.met

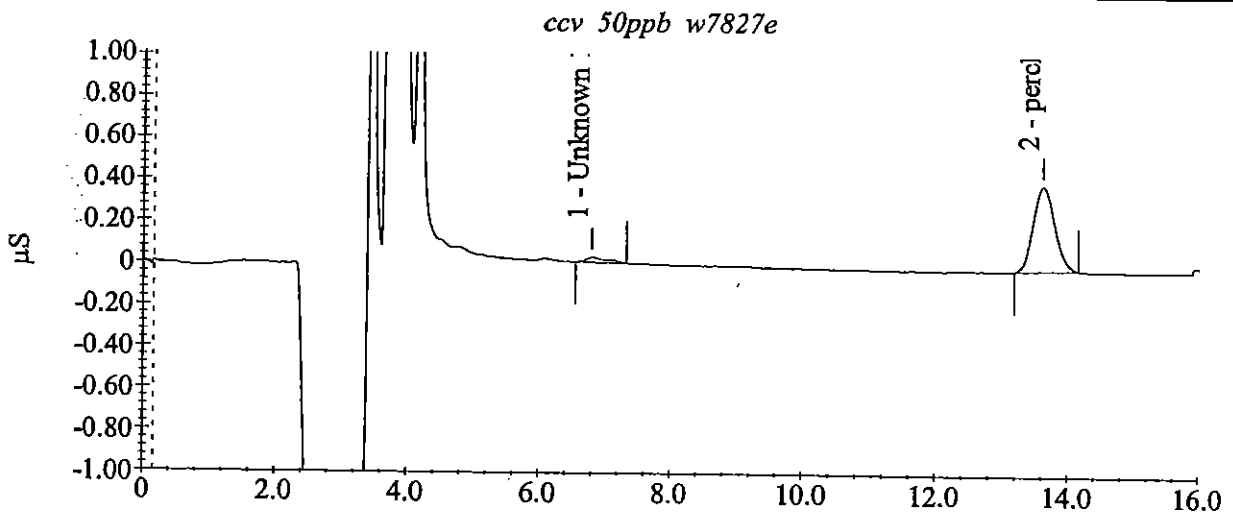
Date Time Collected : 04/23/2003 1:01:22 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
2	perchlorate	13.60	53.85	91392.50	4047.88



APCL Perchlorate Analysis Report

Sample Name : ccv 50ppb w7827e

Data File Name : C:\DATA\03W2496K\W2496K Q04_031.DXD

Method File Name : c:\peaknet\method\314-011.met

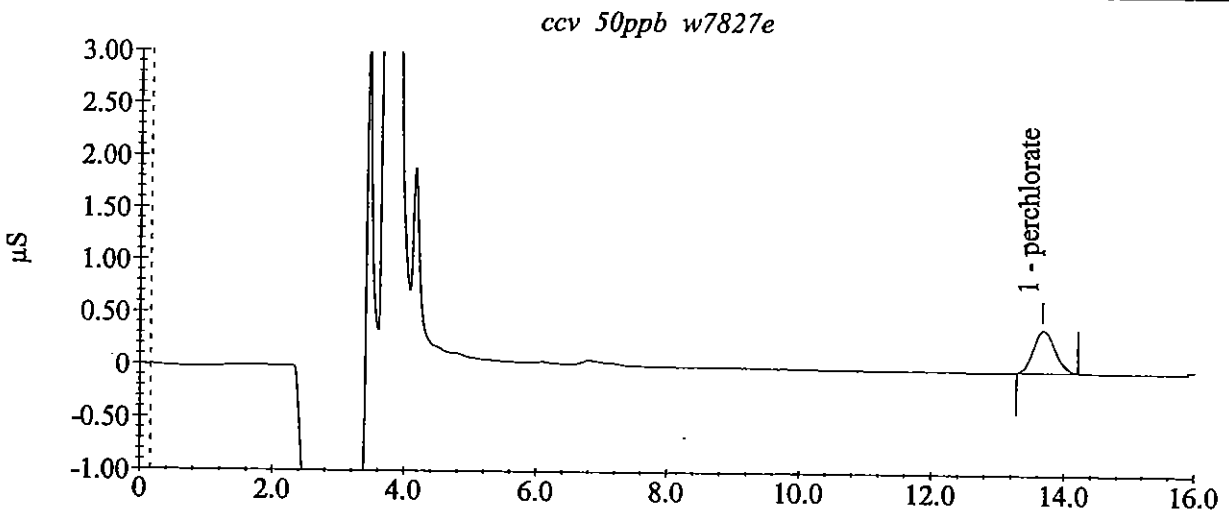
Date Time Collected : 04/23/2003 3:11:29 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	13.67	53.67	91073.15	3991.43



APCL Perchlorate Analysis Report

Sample Name : ccv 50ppb w7827e

Data File Name : C:\DATA\03W2496K\W2496K Q05_032.DXD

Method File Name : c:\peaknet\method\314-011.met

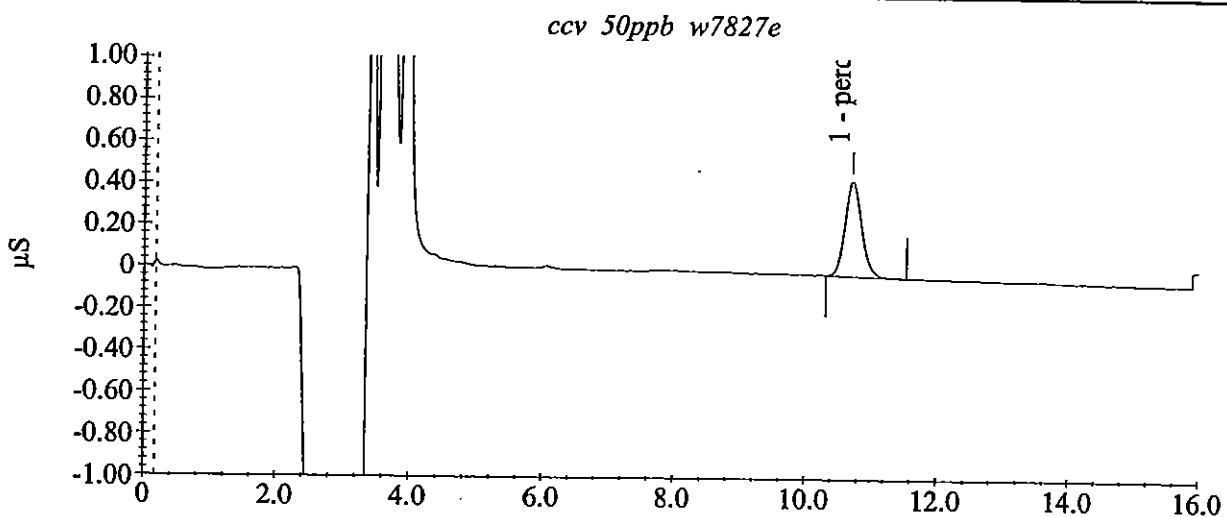
Date Time Collected : 04/23/2003 10:17:46 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	10.72	48.24	81872.40	4536



APCL Perchlorate Analysis Report

Sample Name : ccv 50ppb w7827e

Data File Name : C:\DATA\03W2496K\W2496K Q06_037.DXD

Method File Name : c:\peaknet\method\314-011.met

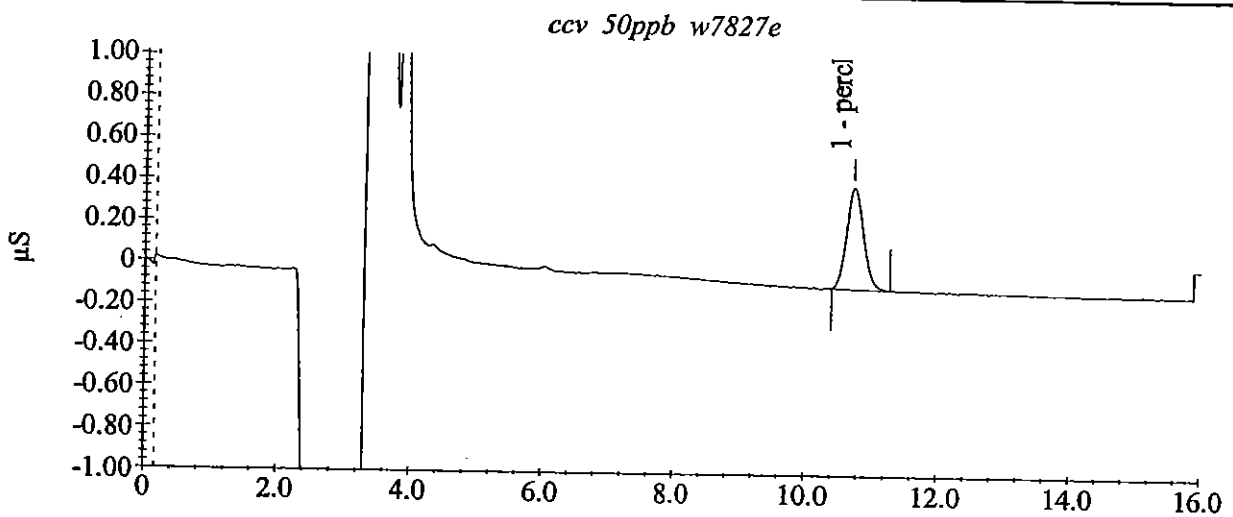
Date Time Collected : 04/23/2003 11:51:10 AM

System Operator : C.W and W.W

Dilution Factor : 1.00

Peak Information : All Components

Peak #	Component Name	Retention Time	Amount (ppb)	Peak Area	Peak Height
1	perchlorate	10.73	51.84	87969.80	4894.80



Alkalinity / OH / CO₃ / HCO₃ (310.1 / SM12320B) Worksheets

Batch # 02002539 Matrix: W Titrant: H₂SO₄ Lot # W2980 Concentration (C): 0.02505 N: Test Date: 4/24/03 Analyst: TR SOP: G-51

#	Sample ID	Dilution V _f /V _i =f ₁	Smp Amt V _i , mL	H ₂ SO ₄ (mL) by Phln S _A	H ₂ SO ₄ (mL) by Phln E _A	H ₂ SO ₄ (mL) by MR-BCG S _B	H ₂ SO ₄ (mL) by MR-BCG E _B	Phln-Alk., P	Tot. Alk., T (in unit of mgCaCO ₃ /L)	OH ⁻	CO ₃ ²⁻	HCO ₃ ⁻	Note & Anomaly
1	MB: <u>T111.1</u>	1 =	100	0	0	0	0	0	0	0	0	0	
2	<u>LCS</u>	1 =	100	7.80	0	7.80	0	99.6	0	0	0	0	
3	<u>LCS</u>	1 =	100	7.81	0	7.81	0	100.3	0	0	0	0	
4	<u>2809-1</u>	1 =	100	6.30	0	6.30	0	161.0	0	0	0	161.0	
5	<u>-2</u>	1 =	100	0	0	0	0	0	0	0	0	0	
6	<u>-3</u>	1 =	100	7.00	0	7.00	0	178.9	0	0	0	178.9	
7	<u>-4</u>	1 =	100	8.71	0	8.71	0	223.16	0	0	0	223.16	
8	<u>-5</u>	1 =	100	7.70	0	7.70	0	196.7	0	0	0	196.7	
9	<u>-6</u>	1 =	100	5.71	0	5.71	0	146.9	0	0	0	146.9	
10	<u>-7</u>	1 =	100	4.90	0	4.90	0	130.3	0	0	0	130.3	
11	<u>-8</u>	1 =	100	0	0	0	0	0	0	0	0	0	
12	<u>2815-1</u>	1 =	100	14.60	0	14.60	0	373.0	0	0	0	373.0	
13	<u>2819-6</u>	1 =	100	6.00	0	6.00	0	154.6	0	0	0	154.6	
14	<u>-5</u>	1 =	100	5.80	0	5.80	0	148.2	0	0	0	148.2	
15	<u>-4</u>	1 =	100	9.90	0	9.90	0	252.9	0	0	0	252.9	
16	<u>-3</u>	1 =	100	8.00	0	8.00	0	209.5	0	0	0	209.5	
17	<u>-2</u>	1 =	100	5.00	0	5.00	0	129.0	0	0	0	129.0	
18	<u>-1</u>	1 =	100	0	0	0	0	0	0	0	0	0	
19	<u>2843-6</u>	1 =	100	6.10	0	6.10	0	151.9	0	0	0	151.9	
20	<u>-7</u>	1 =	100	5.20	0	5.20	0	132.9	0	0	0	132.9	
Dup.	<u>2809-1</u>	1 =	100	6.20	0	6.20	0	158.4	0	0	0	158.4	

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

Calculations:
 A = S_A - E_A
 B = S_B - E_B
 P = 50,000 f₁ A C / V
 T = 50,000 f₁ (A+B) C / V

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

Temperature compensation must be performed by the instrument automatically.

Analyst PC

SOP: G-44

Batch # <u>03W2455</u>	Analysis Date: <u>4/18/03</u>	Batch # <u>03W2477</u>	Analysis Date: <u>4/21/03</u>				
Starting Time: <u>15:51</u>	Ending Time: _____	Starting Time: <u>17:23</u>	Ending Time: _____				
Matrix <input checked="" type="checkbox"/> Aqueous <input type="checkbox"/> Soil		Matrix <input checked="" type="checkbox"/> Aqueous <input type="checkbox"/> Soil					
Standard	4.00	7.00	10.00	Standard	4.00	7.00	10.00
Lot #		<u>2120</u>	<u>030659-24</u>	Lot #		<u>2120</u>	<u>030659-24</u>
Temperature °C		<u>22.8</u>	<u>22.8</u>	Temperature °C		<u>23.0</u>	<u>23.0</u>
pH Reading		<u>7.02</u>	<u>10.01</u>	pH Reading		<u>7.03</u>	<u>10.01</u>
T-corrected pH		<u>7.07</u>	<u>10.02</u>	T-corrected pH		<u>7.01</u>	<u>10.02</u>
Control Limit	±0.05 pH unit			Control Limit	±0.05 pH unit		

#	Sample ID	Pre-treat	pH	Note	#	Sample ID	Pre-treat	pH	Note
MB	<u>T115</u>		<u>6.89</u>		MB	<u>T115</u>		<u>6.86</u>	
1	<u>2789-2</u>		<u>6.79</u>		1	<u>2810-2</u>		<u>7.44</u>	
2					2	<u>2809-1</u>		<u>7.88</u>	
3					3	<u>-2</u>		<u>7.85</u>	
4					4	<u>-3</u>		<u>7.28</u>	
5					5	<u>-4</u>		<u>7.05</u>	
6					6	<u>-5</u>		<u>7.58</u>	
7					7	<u>-6</u>		<u>8.01</u>	
8					8	<u>-7</u>		<u>8.47</u>	
9					9	<u>-8</u>		<u>6.60</u>	
10					10				
11					11				
12					12				
13					13				
14					14				
15					15				
16					16				
17					17				
18					18				
19					19				
20					20				
Dup.	<u>2789-2</u>		<u>6.77</u>		Dup.	<u>2809-4</u>		<u>7.08</u>	

13760 Magnolia Ave. Chino CA 91710

T65
Solid Analysis (160.1, 160.2, 160.3) Worksheet

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

Batch # 2712500 Matrix W Method: 160.1 Balance No. _____

Date: 4/22/03 Analyst: Ru

EPA 160.1 TDS - Total Dissolved (filterable) Solids - Dry for 1hr. or more at 180 °C
 EPA 160.2 TSS - Total Suspended (nonfilterable) Solids - Dry for 1hr. or more at 103-105 °C
 EPA 160.3 TS - Total Solids - Dry for 1hr. or more at 103-105 °C
 Other method (specify): _____

Result = $10^6 \times \Delta W \times f_1 / V$ SOP: G-81

#	Analysis Type	Sample ID (STD Lot #)	Treatment Ratio $V_f/X=f_1$	Volume V_f , mL	W ₁ g	W ₂ 1st, g	W ₂ 2nd, g	$\Delta W = W_2 - W_1$, g	Results (ppm)	Note
1	Blank	T1115	1 =	100	115.5103	115.5104	115.5104		1	C
2	LCS	"	1 =	v	114.0245	114.0666	114.0666		421	D
3	Sample-1	2899-1	1 =	100	116.6483	116.6482	116.6483		0	Z5
4	MS on S-1	3	1 =	↓	115.3396	115.4498	115.4498	110 ²	7002	B
5	MSD on S-1	3	1 =	↓	107.4123	107.4202	107.4202		1079	B/B
6	Sample-2	2	1 =	50	106.3139	106.3220	106.3220		162	8
7	Sample-3	3	1 =	100	105.3551	105.4271	105.4271		720	W ₁
8	Sample-4	4	1 =	↓	99.0748	99.1331	99.1330		582	Y ₂
9	Sample-5	5	1 =	↓	113.8605	113.8800	113.8802		197	6
10	Sample-6	6	1 =	↓	103.9428	103.9780	103.9780		352	14
11	Sample-7	2809-1	1 =	↓	115.1303	115.1521	115.1524		221	10
12	Sample-8	2	1 =	↓	108.5434	108.5430	108.5433		2	IP
13	Sample-9	3	1 =	↓	121.3173	121.3420	121.4417		244	I
14	Sample-10	4	1 =	↓	115.7063	115.7650	115.7650		587	HK
15	LCS	T1115	1 =	↓	107.1671	107.1599	107.1599		428	D ₂
16	Sample-11	5	1 =	↓	104.3088	104.3313	104.3314		226	3
17	Sample-12	6	1 =	↓	114.1411	114.1603	114.1602		191	C ₃
18	Sample-13	7	1 =	↓	115.9112	115.9244	115.9245		133	Y ₉
19	Sample-14	8	1 =	↓	103.5329	103.5329	103.5328		0	Z ₆
20	Sample-15	W4	1 =	↓	104.2613	104.3621	104.3621		1008	7
21	Sample-16	W4D	1 =	↓	106.6522	106.7493	106.7493		971	X
22	Sample-17	2829-1	1 =	↓	107.7078	107.7221	107.7222		144	0
23	Sample-18	2	1 =	↓	111.2781	111.3022	111.3020		239	5
24	Sample-19	3	1 =	↓	114.3693	114.4169	114.4172		479	A
25	Sample-20	4	1 =	↓	111.7908	111.8209	111.8210		302	R
26	1x Dup.	5	1 =	↓	114.3085	114.3040	114.3043		0	H

STD Lot #	$C_{STD}(\mu\text{g/mL}) \times V_{STD}(\text{mL}) / X(\text{g or mL}) = T$	$\frac{W_2 - W_1}{V}$ Spike %	Ctrl Limit (W/S)	PQL/MDL (in ppm)
W- 7618	x 1 = 400 ppm	%	85-115 %/80-120 %	PQL(w) 10
W- "	x 1 = ppm	%	PQL(s) 50
W- 7619	x 1 = ppm	%	90-110 %/85-115 %	MDL(w) 4
W- "	x 1 = ppm	%	MDL(s) 20

Balance Daily Calibration Worksheet

Weight Set S/N: 12006

Calib. Date	Lab Balance					Digital Balance					Analytical Balance					Calib. by
	Balance #	1 g ±0.05g	10 g ±0.1g	200 g ±0.5g	Note (C)	Balance #	1 g ±0.02g	10 g ±0.05g	200 g ±0.10g	Note (D) (C) (AR)	Balance #	1 g ±0.0002g	10 g ±0.0005g	200 g ±0.0010g	Note (D) (C) (AR)	
4/21/03	A-01	Net in Wt			✓	B-01	1.00	10.00	200.01	✓✓✓	C-01	1.0000	10.0000	200.0001	✓✓✓	
	A-02					B-05	1.00	9.99	200.00	✓✓✓	C-02	1.0000	10.0000	200.0000	✓✓✓	
	A-03	1.00	10.00	200.01	✓	B-06	1.00	10.01	200.00	✓✓✓	C-03	1.0000	10.0000	200.0000	✓✓✓	
	A-04					B-07	1.00	10.00	199.99	✓✓✓	C-04	1.0000	10.0000	200.0000	✓✓✓	
	A-					B-					C-					
4/22/03	A-01	Net in Wt			✓	B-01	1.00	10.00	200.01	✓✓✓	C-01	1.0000	10.0000	200.0001	✓✓✓	
	A-02					B-05	1.00	9.99	200.00	✓✓✓	C-02	1.0000	10.0000	200.0000	✓✓✓	
	A-03	1.00	10.00	200.01	✓	B-06	1.00	10.00	200.00	✓✓✓	C-03	1.0000	10.0000	200.0000	✓✓✓	
	A-04					B-07	1.00	10.00	199.99	✓✓✓	C-04	1.0000	10.0000	200.0000	✓✓✓	
	A-					B-					C-					
4/24/03	A-01	Net in Wt			✓	B-01	1.00	10.00	200.00	✓✓✓	C-01	1.0000	9.9999	200.0000	✓✓✓	
	A-02					B-05	1.00	10.01	200.01	✓✓✓	C-02	1.0000	10.0000	200.0002	✓✓✓	
	A-03	1.00	10.00	200.00		B-06	1.00	10.00	200.00	✓✓✓	C-03	1.0000	10.0000	200.0000	✓✓✓	
	A-04					B-07	1.00	9.99	199.99	✓✓✓	C-04	1.0000	10.0000	200.0000	✓✓✓	
	A-					B-					C-					

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

Line	Sample	Sample Type	Level	Method	Data File	Volume	Dilution
1	##03w2496kw ipc 25ppb w7759	Sample		e314-011.met	c:\data\03w2496kw2496k ipc 25ppb	1	1
2	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q01	1	1
3	ccb	Sample		e314-011.met	c:\data\03w2496kw2496k ccb01	1	1
4	lcs 25ppb w7827d	Sample		e314-011.met	c:\data\03w2496kw2496k i01	1	1
5	LCS 18PPB W7685D	Sample		e314-011.met	c:\data\03w2496kw2496k j01	1	1
6	ICCS 4ppb w7827b	Sample		e314-011.met	c:\data\03w2496kw2496k iccs 4ppb	1	1
7	mb	Sample		e314-011.met	c:\data\03w2496kw2496k k01	1	1
8	2805-19 F=1	Sample		e314-011.met	c:\data\03w2496k\2805-19	1	1
9	2809-01 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-01	1	1
10	2809-02 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-02	1	1
11	2809-03 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-03	1	1
12	2809-04 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-04	1	1
13	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q02	1	1
14	ccb	Sample		e314-011.met	c:\data\03w2496kw2496k k02	1	1
15	2809-05 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-05	1	1
16	2809-06 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-06	1	1
17	2809-07 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-07	1	1
18	2809-08 F=1	Sample		e314-011.met	c:\data\03w2496k\2809-08	1	1
19	2819-01 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-01	1	1
20	2819-02 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-02	1	1
21	2819-03 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-03	1	1
22	2819-04 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-04	1	1
23	2819-05 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-05	1	1
24	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q03	1	1
25	CCB	Sample		e314-011.met	c:\data\03w2496kw2496k k03	1	1
26	2819-06 F=1	Sample		e314-011.met	c:\data\03w2496k\2819-06	1	1
27	2809-04 MS 50PPB F=1	Sample		e314-011.met	c:\data\03w2496kw2496k m01	1	1
28	2809-04 MSD 50PPB F=1	Sample		e314-011.met	c:\data\03w2496kw2496k n01	1	1
29	2819-03 MS 50PPB F=1	Sample		e314-011.met	c:\data\03w2496kw2496k m02	1	1
30	2819-03 MSD 50PPB F=1	Sample		e314-011.met	c:\data\03w2496kw2496k n02	1	1
31	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q04	1	1
32	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q05	1	1
33	CCB	Sample		e314-011.met	c:\data\03w2496kw2496k k04	1	1
34	2820-05 F=1	Sample		e314-011.met	c:\data\03w2496k\2820-05	1	1
35	2820-06 F=1	Sample		e314-011.met	c:\data\03w2496k\2820-06	1	1
36	2820-09 F=1	Sample		e314-011.met	c:\data\03w2496k\2820-09	1	1
37	ccv 50ppb w7827e	Sample		e314-011.met	c:\data\03w2496kw2496k q06	1	1
38		Sample		aastopcl.met		1	1

Analyst Wei Wang
Date 4/22-23/03
Instrument IC-K

Line	Sample	Sample Type	Level	Method	Data File	Volume	Dilution
1	Cal blank	Sample		e314-011.met	c:\data\314-011\mb_001.dxd	1	1
2	cal standard 2ppb W7827a	Sample		e314-011.met	c:\data\314-011\std-2pb_002.dxd	1	1
3	cal standard 4ppb W7827b	Sample		e314-011.met	c:\data\314-011\std-4pb_003.dxd	1	1
4	cal standard 10ppb W7827c	Sample		e314-011.met	c:\data\314-011\std-10pb_004.dxd	1	1
5	cal standard 25ppb W7827d	Sample		e314-011.met	c:\data\314-011\std-25pb_005.dxd	1	1
6	cal standard 50ppb W7827e	Sample		e314-011.met	c:\data\314-011\std-50pb_006.dxd	1	1
7	cal standard 75ppb W7827f	Sample		e314-011.met	c:\data\314-011\std-75pb_007.dxd	1	1
8	cal standard 100ppb W7827g	Sample		e314-011.met	c:\data\314-011\std-100pb_008.dxd	1	1
9	ICV 50 ppb w7828a	Sample		e314-011.met	c:\data\314-011\icv-50pb_009.dxd	1	1
10	icb	Sample		e314-011.met	c:\data\314-011\icb_010.dxd	1	1
11	anion 100pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-100_011.dxd	1	1
12	anion 200pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-200_012.dxd	1	1
13	anion 300pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-300_013.dxd	1	1
14	anion 400pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-400_014.dxd	1	1
15	anion 500pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-500_015.dxd	1	1
16	anion 600pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-600_016.dxd	1	1
17	anion 800pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-800_017.dxd	1	1
18	anion 1000pm each ,25pb CLO4	Sample		e314-011.met	c:\data\314-011\mct-1000_018.dxd	1	1
19	anion 400pm each 2pb	Sample		e314-011.met	c:\data\314-011\ipc-2pb_019.dxd	1	1
20	anion 400pm each 4pb	Sample		e314-011.met	c:\data\314-011\ipc-4pb_020.dxd	1	1
21	anion 400pm each 25pb	Sample		e314-011.met	c:\data\314-011\ipc-25pb_021.dxd	1	1
22	ICV 50 ppb	Sample		e314-011.met	c:\data\314-011\icv-50pb	1	1
23	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-02_023.dxd	1	1
24	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-03_024.dxd	1	1
25	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-04	1	1
26	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-05	1	1
27	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-06	1	1
28	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-07	1	1
29	MDL 4pb	Sample		e314-011.met	c:\data\314-011\mdl-08	1	1
30	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
31	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
32	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
33	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
34	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
35	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
36	IDP and IDA 25pb	Sample		e314-011.met	c:\data\314-011\idap-25pb	1	1
37	MCT anion 800pm each, 25pbCLO4	Sample		e314-011.met	c:\data\314-011\ipc-25pb	1	1
38	MCT anion 800pm each, 25pbCLO4	Sample		e314-011.met	c:\data\314-011\ipc-25pb	1	1
39	MCT anion 800pm each, 4pbCLO4	Sample		e314-011.met	c:\data\314-011\ipc-4pb	1	1
40	MCT anion 800pm each, 4pbCLO4	Sample		e314-011.met	c:\data\314-011\ipc-4pb	1	1
41	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s01	1	5
42	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s02	1	5
43	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s03	1	5
44	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s04	1	5
45	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s05	1	5
46	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s06	1	5
47	MDL 20pb soil	Sample		e314-011.met	c:\data\314-011\mdl-s07	1	5
48	standard 25ppb W7827d	Sample		e314-011.met	c:\data\314-011\std-25pb	1	1
49	anion 100pm each,4pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-100-4pb	1	1
50	anion 200pm each ,4pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-200-4pb	1	1
51	anion 300pm each ,4pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-300-4pb	1	1
52	anion 100pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-100-2pb	1	1
53	anion 200pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-200-2pb	1	1
54	anion 300pm each,2pb CLO4	Sample		e314-011.met	c:\data\314-011\lam-300-2pb	1	1
5	1982-01 B S.C 4450us/cm	Sample		e314-011.met	c:\data\314-011\1982-01	1	1
6	1982-01 B S.C 4450us/cm	Sample		e314-011.met	c:\data\314-011\1982-01	1	1
7	1982-02 f=10	Sample		e314-011.met	c:\data\314-011\1982-02_057.dxd	1	2
8		Sample		aastopcl.met		1	10

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

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

Comment:

Analyst DN
 Date 3/21/03
 Instrument J

DIONEX SCHEDULE - C:\DX\SCHEDULE\03W2479.SCH

Inj#	Sample Name	Method	Data File	Vol.	Dil.	Int.Std.
1	##03W2479, W CCVW77	..\E300-063	..\W2479Q01.D01	1	1	1
2	MB RW1408	..\E300-063	..\W2479K01.D02	1	1	1
3	LCS W7768-100X	..\E300-063	..\W2479L01.D03	1	1	1
4	LCSD W7768-100X	..\E300-063	..\W2479J01.D04	1	1	1
5	2809-2 F=1.25	..\E300-063	..\2809-201.D05	1	1.25	1
6	2809-8 F=1.25	..\E300-063	..\2809-801.D06	1	1.25	1
7	2809-1 F=2.5	..\E300-063	..\2809-101.D07	1	2.5	1
8	2809-3 F=2	..\E300-063	..\2809-301.D08	1	2	1
9	2809-4 F=20	..\E300-063	..\2809-401.D09	1	20	1
10	2809-5 F=4	..\E300-063	..\2809-501.D10	1	4	1
11	2809-6 F=4	..\E300-063	..\2809-601.D11	1	4	1
12	CCV2W7767-100X	..\E300-063	..\W2479Q01.D12	1	1	1
13	MB RW1408	..\E300-063	..\W2479K01.D13	1	1	1
14	\$2809-4 MS F=40	..\E300-063	..\W2479M01.D14	1	40	1
15	\$2809-4 MSD F=40	..\E300-063	..\W2479N01.D15	1	40	1
16	2815-1 F=2	..\E300-063	..\2815-101.D16	1	2	1
17	2809-7 F=4	..\E300-063	..\2809-701.D17	1	4	1
18	CCV3W7767-100X	..\E300-063	..\W2479Q01.D18	1	1	1
19	MB RW1408	..\E300-063	..\W2479K11.D19	1	1	1
20	\$2819-3 MS F=40	..\E300-063	..\W2479M01.D20	1	40	1
21	\$2819-3 MSD F=40	..\E300-063	..\W2479N01.D21	1	40	1
22	2819-3 F=20	..\E300-063	..\2819-301.D22	1	20	1
23	2819-4 F=12.5	..\E300-063	..\2819-401.D23	1	12.5	1
24	2819-5 F=4	..\E300-063	..\2819-501.D24	1	4	1
25	2819-6 F=10	..\E300-063	..\2819-601.D25	1	10	1
26	2819-2 F=1.25	..\E300-063	..\2819-201.D26	1	1.25	1
27	2819-1 F=1.25	..\E300-063	..\2819-101.D27	1	1.25	1
28	CCV4W7767-100X	..\E300-063	..\W2479Q01.D28	1	1	1
29		..\STOP.MET		1	1	1

Comment :

LCS/LCSD LOT # W7768

MS/MSD LOT # W7767

ELUENT LOT # W7868

ANALYTICAL METHOD 9056/E300 MATRIX W

Analyst ZC

Date 4/22/03

Instrument 1

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

Chromium (VI) (7196) Worksheet

Batch # 23W 2476 Matrix: W

[Holding Time: 24 hours!!]

Test Date: 4/21/03 Analyst: BW

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

Calibration	STD Lot #	$C_{std} \times V_{std} / V_f = C_i$	A_i	$RF_i = A_i / C_i$	Calibration results	Note
STD-1	W-	x / = mg/L			Least Square [RF]=	Cal. Code:
STD-2	W-	x / = mg/L			Average RF=	
STD-3	W-	x / = mg/L			C.C. = <u>0.9972</u> > 0.995)	
STD-4	W-	x / = mg/L			RSD = % (< 15%)	
STD-5	W-	x / = mg/L			Ref. page	
STD-6	W-	x / = mg/L			<u>A = 0.000 + 0.8360</u>	

Analysis Type	Sample ID or Lot #	Samp. Amt X ₀ (g or mL)	Dilu./Ext X/X ₀ =f ₁	Treat. Ratio V/X=f ₂	540 nm A	Concentration C'=A/RF	C (Sample) C=f ₁ f ₂ C'	Anomaly Note
CCV	Lot: W- <u>7853</u>	Expected Conc.: x	/	= <u>0.25</u> mg/L	<u>0.211</u>	<u>0.25</u> mg/L	REC. %	90-110 %
Method Blank	Bl. Lot: <u>T1115</u>		/X ₀ =	95.0/ =	<u>0.000</u>	mg/L	<u>0.00</u> ppm	
LCS1	Bl. Lot: '1		/X ₀ =	95.0/ =	<u>0.188</u>	mg/L	<u>0.215</u> ppm	
Sample-1	<u>2809-1</u>		/X ₀ =	95.0/ =	<u>0.000</u>	mg/L	<u>0.000</u> ppm	
MS on S-1	<u>4</u>		/X ₀ =	95.0/ =	<u>0.190</u>	mg/L	<u>0.227</u> ppm	
MSD on S-1	<u>4</u>		/X ₀ =	95.0/ =	<u>0.189</u>	mg/L	<u>0.226</u> ppm	
Sample 2	<u>2</u>		/X ₀ =	95.0/ =	<u>0.001</u>	mg/L	<u>0.001</u> ppm	
Sample 3	<u>3</u>		/X ₀ =	95.0/ =	<u>0.001</u>	mg/L	<u>0.001</u> ppm	
Sample 4	<u>4</u>		/X ₀ =	95.0/ =	<u>0.003</u>	mg/L	<u>0.003</u> ppm	
Sample 5	<u>5</u>		/X ₀ =	95.0/ =	<u>0.002</u>	mg/L	<u>0.002</u> ppm	
Sample 6	<u>6</u>		/X ₀ =	95.0/ =	<u>0.001</u>	mg/L	<u>0.001</u> ppm	
Sample 7	<u>7</u>		/X ₀ =	95.0/ =	<u>0.000</u>	mg/L	<u>0.000</u> ppm	
Sample 8	<u>8</u>		/X ₀ =	95.0/ =	<u>0.000</u>	mg/L	<u>0.000</u> ppm	
Sample 9			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 10			/X ₀ =	95.0/ =		mg/L	ppm	
Blank	Lot:		/X ₀ =	95.0/ =		mg/L	ppm	
LCS2	Bl. Lot: <u>T1115</u>		/X ₀ =	95.0/ =	<u>0.192</u>	mg/L	<u>0.230</u> ppm	
Sample 11			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 12			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 13			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 14			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 15			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 16			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 17			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 18			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 19			/X ₀ =	95.0/ =		mg/L	ppm	
Sample 20			/X ₀ =	95.0/ =		mg/L	ppm	
X Dup.	<u>claring 0.25 mg/L</u>		/X ₀ =	95.0/ =	<u>0.209</u>	<u>0.250</u> mg/L	ppm	

STD Lot #	$C_{STD} (\mu\text{g/mL}) \times V_{STD} (\text{mL}) / X (\text{g or mL}) = T$	Spike Rec.	Cl Limit (W/S)	PQL/MDL (in ppm)
W- <u>7757</u>	x / = <u>0.25</u> ppm	%	80-120 %/80-120 %	PQL(w) 0.01
W- '1	x / = ppm	%	PQL(s) 0.05
W- <u>7853</u>	x / = ppm	%	80-120 %/80-120 %	MDL(w) 0.005
W- '1	x / = ppm	%	MDL(s) 0.025

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

Chromium (VI) (7196) Worksheet

Batch # 02W1295 Matrix: W

[Holding Time: 24 hours!!]

Test Date: 1/29/03 Analyst: Bi

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

Calibration	STD Lot #	$C_{std} \times V_{std} / V_f = C_i$	A_i	$RF_i = A_i / C_i$	Calibration results	Note
STD-1	W-7191	x / = 0.000 mg/L	0.000		Least Square [RF]=	Cal. Code:
STD-2	W-	x / = 0.012 mg/L	0.005		Average RF=	A=0.000 + 0.836C
STD-3	W-	x / = 0.020 mg/L	0.011		C.C.=0.997 (> 0.995)	
STD-4	W-	x / = 0.125 mg/L	0.109		RSD= % (< 15%)	
STD-5	W-	x / = 0.250 mg/L	0.214		Ref. page	
STD-6	W- ✓	x / = 0.500 mg/L	0.415			A=0.003 + 0.836C

Analysis Type	Sample ID or Lot #	Samp. Amt X ₀ (g or mL)	Dilu./Ext X/X ₀ =f ₁	Treat. Ratio V/X=f ₂	540 nm A	Concentration C'=A/RF	C (Sample) C=f ₁ f ₂ C'	Anomaly Note
CCV	Lot: W-7076	Expected Conc.: x	1	= 0.25 mg/L	0.216	0.258 mg/L	REC. %	90-110 %
Method Blank	Bl. Lot: T1115		1/X ₀ = 1	95.0/ =	0.000	mg/L	0.00 ppm	
LCS1	Bl. Lot: 1		1/X ₀ =	95.0/ =	0.204	mg/L	0.244 ppm	
Sample-1	1369-1		1/X ₀ =	95.0/ =	0.000	mg/L	0.00 ppm	
MS on S-1	6		1/X ₀ =	95.0/ =	0.223	mg/L	0.266 ppm	
MSD on S-1	6		1/X ₀ =	95.0/ =	0.230	mg/L	0.275 ppm	
Sample 2	2		1/X ₀ =	95.0/ =	0.004	mg/L	0.005 ppm	
Sample 3	3		1/X ₀ =	95.0/ =	0.002	mg/L	0.002 ppm	
Sample 4	4		1/X ₀ =	95.0/ =	0.001	mg/L	0.001 ppm	
Sample 5	5		1/X ₀ =	95.0/ =	0.002	mg/L	0.002 ppm	
Sample 6	6		1/X ₀ = ✓	95.0/ =	0.004	mg/L	0.005 ppm	
Sample 7	-		1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 8			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 9			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 10			1/X ₀ =	95.0/ =		mg/L	ppm	
Blank	Lot:		1/X ₀ =	95.0/ =		mg/L	ppm	
LCS2	Bl. Lot: T1115		1/X ₀ = 1	95.0/ =	0.210	mg/L	0.25 ppm	
Sample 11			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 12			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 13			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 14			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 15			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 16			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 17			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 18			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 19			1/X ₀ =	95.0/ =		mg/L	ppm	
Sample 20			1/X ₀ =	95.0/ =		mg/L	ppm	
MTX Dup.	closing 0.25 mg/L		1/X ₀ =	95.0/ =	0.204	mg/L	0.204 ppm	

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



Applied Physics & Chemistry Laboratory

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

June 25, 2003

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

Dear Leo Williamson,

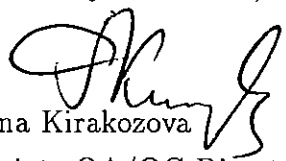
This package contains samples in our Service ID 03-3351 and your project : 04-4428.10 JPL GW Mon-2Q03.

Enclosed please find:

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

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

Respectfully submitted,


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

Applied P & Ch Laboratory

13760 Magnolia Ave. Chino CA 91710

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

Submitted to:

GEOFON, Inc.

Attention: Leo Williamson

22632 Golden Spring Dr Ste 270

Diamond Bar CA 91765

Tel: (909)396-7662 Fax: (909)396-1455

APCL Analytical Report

Service ID #: 801-033351

Received: 05/22/03

Collected by: Leo Williamson

Extracted: 05/22/03

Collected on: 05/22/03

Tested: 05/23/03

Reported: 05/29/03

Sample Description: Water from MW-4

Project Description: 04-4428.10 JPL. GW-Mon. 2Q03.

Analysis of Water Samples

Component Analyzed	Method	Unit	PQL	Analysis Result		
				DUPE-8-2Q03 03-03351-1	EB-14-5/22/03 03-03351-2	MW-4-2 03-03351-3
Dilution Factor				1	1	1
1,4-DIOXANE	8270-SIM	µg/L	1	1	<1	1

Component Analyzed	Method	Unit	PQL	Analysis Result		
				MW-4-3 03-03351-4	MW-4-4 03-03351-5	MW-4-5 03-03351-6
Dilution Factor				1	1	1
1,4-DIOXANE	8270-SIM	µg/L	1	0.4J	<1	<1

PQL: Practical Quantitation Limit. MDL: Method Detection Limit. CRDL: Contract Required Detection Limit

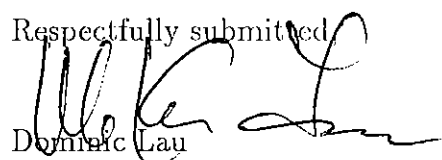
N.D.: Not Detected or less than the practical quantitation limit.

"-": Analysis is not required.

J: Reported between PQL and MDL.

Listed Dilution Factors (DF) are relative to the method default DF. All unlisted DFs are 1.0

Respectfully submitted,


Dominic Lau

Laboratory Director

Applied P & Ch Laboratory

Level C Data Package Deliverables

General Information

Project: JPL

APCL Service ID: 03-3351



Applied P & Ch Laboratory

13760 Magnolia Ave. Chino, CA 91710

Telephone (909)590-1828

Fax (909)590-1498



INCORPORATED
22632 GOLDEN SPRINGS DR., SUITE 270
DIAMOND BAR, CA 91765 • (909) 396-7662 • FAX (909) 396-1455

CHAIN-OF-CUSTODY RECORD

LABORATORY COPY

MW-4 + 0033

LAB COORDINATOR Brad Shojaee PROJECT NAME JPL GW-MON-2403 PROJECT CONTACT Leo W. Williamson PROJECT ADDRESS 4800 Oak Grove Dr PROJECT MANAGER Asrar Fahem	LAB COORDINATOR'S PHONE (909) 396-7662 PROJECT LOCATION MW-4 (SE Facility Reimeter) PROJECT PHONE NUMBER (714) 920-8729 CITY, STATE AND ZIP CODE Pasadena, CA PHONE (909) 396-7662	LAB COORDINATOR'S FAX (909) 396-1455 PROJECT NUMBER 04-4428.10 PROJECT FAX (909) 396-1455 CLIENT US NAVY, 96RDIV PROJECT MANAGER'S FAX (909) 396-1455	LABORATORY SERVICE ID —	LABORATORY CONTACT Kenny Chan LABORATORY PHONE (909) 590-1828 LABORATORY ADDRESS 13760 Magnolia Ave. CITY, STATE AND ZIP CODE Chicago, CA 91710	MAIL REPORT (COMPANY NAME) GEOFON, INC RECIPIENT NAME Leo W. Williamson ADDRESS 22632 Golden Springs Dr. #270 CITY, STATE AND ZIP CODE Diamond Bar, CA. 91765
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Item	Sample Identifier	Matrix			Date	Time	Preserved	# of Cont	QC Level	T.A.T	Analyses	Comments
		H ₂ O	↓	↓								
1	MW-4-5	H ₂ O	5/21/03	915	NONE	2	III	NORMAL	X			
2	MW-4-4			1000					X			
3	MW-4-3			1040					X			
4	MW-4-2			1210					X			
5												
6	EB-14-5/22/03	H ₂ O	5/22/03	1050	NONE	2	III	NORMAL	X			
7	DUPE-8-2403			—		2	IV		X			
8												
9												
10												

3351

SAMPLES COLLECTED BY: Leo W. Williams	COURIER AND AIR BILL NUMBER.	COOLER TEMPERATURE UPON RECEIPT
RELINQUISHED BY: Leo W. Williamson	RECEIVED BY: Asrar Fahem	SAMPLE'S CONDITION UPON RECEIPT
DATE: 5/22/03	DATE: 5/22/03	
TIME: 1350	TIME: 1354	

Case Narrative

Project: JPL GW-Mon. 2Q03./MW-4/04-4428.10

For GEOFON, Inc.

APCL Service No: 03-3351

1. Sample Identification

The sample identifications are listed in the following table:

GEOFON, Inc. Sample ID	APCL Sample ID
MW-4-5	03-03351-6
MW-4-4	03-03351-5
MW-4-3	03-03351-4
MW-4-2	03-03351-3
EB-14-5/22/03	03-03351-2
DUPE-8-2Q03	03-03351-1

2. Analytical Methodology

Samples are analyzed by EPA methods
8270-SIM (1,4-Dioxane),

3. Holding Time

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

4. Preservation

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

5. Tele-log

None

6. Anomaly

None

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

Respectfully submitted,



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

Sample Receiving Checklist

3357

APCL ServiceID: _____ Client Name/Project: Calson Inc

1. Sample Arrival

Date/Time Received 5/22/03 1355 Date/Time Opened 5/22/03 1355 By (name): Paul Kon

Custody Transfer: Client Golden State UPS US Mail FedEx APCL Empl: Adam Wood

2. Chain-of-Custody (CoC)

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

3. Shipping Container/Cooler

Cooler Used? # of 1 Cooled by: Ice Blue Ice Dry Ice None
Temp °C 2.2°C

(Cooler temperature measured from temp blank if present, otherwise measured from the cooler).

Cooler Custody Seal? Absent Intact Tampered?

4. Sample Preservation

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

5. Holding-time Requirements

pH 24hr BACT 6/24hr Cr^{VI} 24hr NO₃⁻ 48hr BOD 48hr
 Cl₂ ASAP Turbidity 48hr DO ASAP Fe(II) ASAP
 HT Expired? Client notified?

6. Sample Container Condition

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

7. Turn Around Time

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

8. Sample Matrix

Drinking H₂O Other Liq Soil Wipe Polymer Air Other: _____
 Ground H₂O Sludge Filter Oil/Petro Paint W. Water Extract Unknown

9. Pre-Login Check List Completed & OK?

ALL OK? (if not, attach docs) Client Contact? (Name: _____) Date/Time: _____

Received/Checked by: Paul Kon Date: 22 May 2003 Time: 7:40 a.m.

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

Applied P & Ch Laboratory

13760 Magnolia Ave. Chino CA 91710

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

Sample Login: Check List

03-03351 (0470_ 148) (2202777_ 148)

05/22/03

Part 1: General Information

- | | | | |
|--------------------------|--------------------------|----------------------|--|
| <input type="checkbox"/> | Company Information | Name: | <i>GEOFON, Inc.</i> |
| | | Address: | <i>22632 Golden Spring Dr Ste 270 ,Diamond Bar ,CA 91765</i> |
| <input type="checkbox"/> | Project Information | Project Description: | <i>JPL</i> |
| | | Project #: | <i>04-4428.10</i> |
| <input type="checkbox"/> | Billing Information | P.O. #: | |
| | | Bill Address: | <i>22632 Golden Spring Dr Ste 270 ,Diamond Bar ,CA 91765</i> |
| | | Lab Project ID: | |
| | | Client Database #: | <i>3</i> |
| <input type="checkbox"/> | Receiving Information | Who Received Sample? | <i>Paul Kou</i> |
| | | Receiving Date/Time: | <i>05/22/03 1355</i> |
| | | COC No. | |
| <input type="checkbox"/> | Shipping Information | Shipping Company | <i>APCL pick up</i> |
| | | Packing Information: | <i>Cooler/Ice Chester</i> |
| | | Cooler Temperature: | <i>2.2 °C</i> |
| <input type="checkbox"/> | Container Information | Container Provider: | <i>Client</i> |
| <input type="checkbox"/> | Sampling Information | Sampling Person: | |
| | | Sampling Company: | <i>Client</i> |
| <input type="checkbox"/> | Turn-Around-Time Option: | | <i>Rush 5 working day(s)</i> |
| <input type="checkbox"/> | QC Option: | | <i>NEESA C</i> |
| <input type="checkbox"/> | Disposal Option: | | <i>Not specify</i> |
-

Part 2: Sample Information

Seq. #	Sample ID (on COC)	Sample Sub-ID	APCL Sample ID	Matrix	Cont- tainer	Preser- vative	Vol, ml Am. g	# of Replica	Condition G, L, B	Collected mmddyy	Hold ?	Composite Group	TAT Days	
1	MW-4-5	Dioxane	03-03351-6	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>
2	MW-4-4	Dioxane	03-03351-5	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>
3	MW-4-3	Dioxane	03-03351-4	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>
4	MW-4-2	Dioxane	03-03351-3	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>
5	EB-14-5/22/03	Dioxane	03-03351-2	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>
6	DUPE-8-2Q03	Dioxane	03-03351-1	W	G		1000	2	G	052203	N	0	7	<input type="checkbox"/>

Part 3: Analysis Information

Test Items:	<input type="checkbox"/> 524.2	Volatile Organic Compounds
	<input type="checkbox"/> 7196A	Chromium (VI)
	<input type="checkbox"/> 314.0/300.0	Perchlorate, low level
	<input type="checkbox"/> 300.0	Chloride Cl ⁻ by IC
	<input type="checkbox"/> 300.0	Sulfate (SO ₄ ⁻), by IC
	<input type="checkbox"/> 300.0/SM4500NO3	Nitrate (NO ₃ ⁻) as N by IC
	<input type="checkbox"/> SM2320B	Carbonate
	<input type="checkbox"/> SM2320B	Bicarbonate
	<input type="checkbox"/> 9040B/150.1	pH
	<input type="checkbox"/> 160.1	Solids, Total Dissolved (TDS)
	<input type="checkbox"/> 200.7/6010B	Sodium, Na, by ICP
	<input type="checkbox"/> 200.7/6010B	Calcium, Ca, by ICP
	<input type="checkbox"/> 200.7/6010B	Potassium, K, by ICP
	<input type="checkbox"/> 200.7/6010B	Magnesium, Mg, by ICP
	<input type="checkbox"/> 200.7/6010B	Iron, Fe, by ICP
	<input type="checkbox"/> 206.2/7060A	Arsenic, As, by GFAA
	<input checked="" type="checkbox"/> 8270-SIM	1,4-Dioxane

Seq. #	Client's Sample ID (as given on COC)	Sample Sub-ID	APCL Sample ID	Matrix	524.2	CHROMIUM	PERCH	CL	SO4	NO3	CARBON	BICARB	
1	MW-4-5	Dioxane	03-03351-6	W									<input type="checkbox"/>
2	MW-4-4	Dioxane	03-03351-5	W									<input type="checkbox"/>
3	MW-4-3	Dioxane	03-03351-4	W									<input type="checkbox"/>
4	MW-4-2	Dioxane	03-03351-3	W									<input type="checkbox"/>
5	EB-14-5/22/03	Dioxane	03-03351-2	W									<input type="checkbox"/>
6	DUPE-8-2Q03	Dioxane	03-03351-1	W									<input type="checkbox"/>

Level C Data Package Deliverables

8270-SIM



Applied P & Ch Laboratory
13760 Magnolia Ave. Chino, CA 91710
Telephone (909)590-1828
Fax (909)590-1498

Applied P & Ch Laboratory
Organic Analysis Results for Method 8270-SIM

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/22/2003
Project ID: JPL GW-Mon. 2Q03.	Service ID: 33351	Collected by:
Sample ID: 03G2623-MB-01	Lab Sample ID: 03G2623-MB-01	Received Date: 05/22/2003
Sample Type: Method Blank	Sample Matrix: Water	Moisture %: -
Anal. Method: 8270-SIM	Prep. Method: 3520	Instrument ID: GC/MS: M
Batch No: 03G2623	Prep. Date: 05/22/03	Anal. Date: 05/23/03
Data File Name: G2623K01	Prep. No: 1 of 1	Anal. Time: 13:10
Extract Vol. 1.0 mL	Sample Amount: 1000 mL	Dilution Factor: 1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	<1	U
Internal Standard				Control Limit, %	IS Rec.%	
1	1,4-DIOXANE-D8	17647-74-4		50-200	92	
# of out-of-control					0	

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

Qualifier: U - Not Detected or less than MDL	E - Exceed calibration range
J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)	B - A positive value was found in the method blank
	D - Diluted

Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623K01.D Sample : F=.001
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 13:10 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 23 14:57 2003 Multiplr: 0.001000
 Print Time : Fri May 23 14:57 2003
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
1	1.4-Dioxane-d8	4.83	4.80	0.006	96	66	296.242	20.00		Dev (Min)	
										0.03	

System Monitoring Compounds (Surrogate) %Recovery

Target Compounds Qvalue

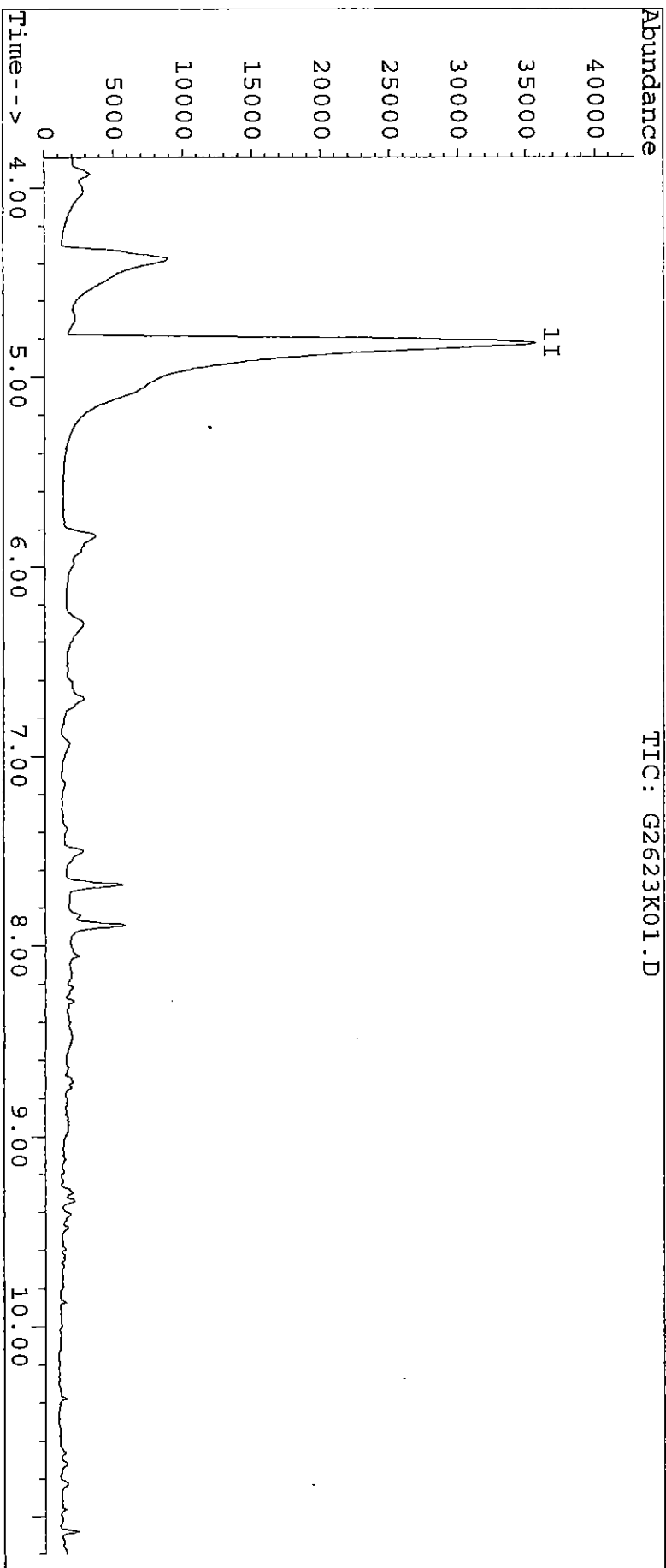
 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623K01.D Vial: 1
Acq On : 23 May 03 1:10 pm Operator: Andy Huang
Sample : F=.001 Inst : GC/MS - M
Misc : Multiplr: 1.00
Quant Time: May 23 14:57 2003 Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Fri May 23 12:56:26 2003
Response via : Multiple Level Calibration



Applied P & Ch Laboratory
Organic Analysis Results for Method 8270-SIM

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/22/2003
Project ID: JPL GW-Mon. 2Q03.	Service ID: 33351	Collected by:
Sample ID: DUPE-8-2Q03	Lab Sample ID: 03-3351-1	Received Date: 05/22/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 8270-SIM	Prep. Method: 3520	Instrument ID: GC/MS: M
Batch No: 03G2623	Prep. Date: 05/22/03	Anal. Date: 05/23/03
Data File Name: 3351-01	Prep. No: 1 of 1	Anal. Time: 21:45
Extract Vol. 1.0 mL	Sample Amount: 1000 mL	Dilution Factor: 1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	1	
Internal Standard				Control Limit, %	IS Rec. %	
1	1,4-DIOXANE-D8	17647-74-4		50-200	95	
# of out-of-control					0	

Qualifier: U - Not Detected or less than MDL
 J - Less than RL (PQL, EQL or CRDL), but greater than MDL, or an estimated result (e.g. for TIC)
 E - Exceed calibration range
 B - A positive value was found in the method blank
 D - Diluted

Data Filename: C:\HPCHEM\1\DATA\03G2623\3351-01.D Sample : F=.001
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 21:45 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 27 13:50 2003 Multiplr: 0.001000
 Print Time : Tue May 27 13:50 2003
 Miscleaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
Internal Standards											
1	1,4-Dioxane-d8	4.77	4.80	-0.004	96	66	306.192	20.00			Dev(Min) -0.02

System Monitoring Compounds (Surrogate) %Recovery

Target Compounds	<<< I1 : ISTD ID = 1 >>>	4.87	4.88	-0.002	88	58	19.664	1.06	1.1	78	Qvalue
2	1,4-Dioxane										✓

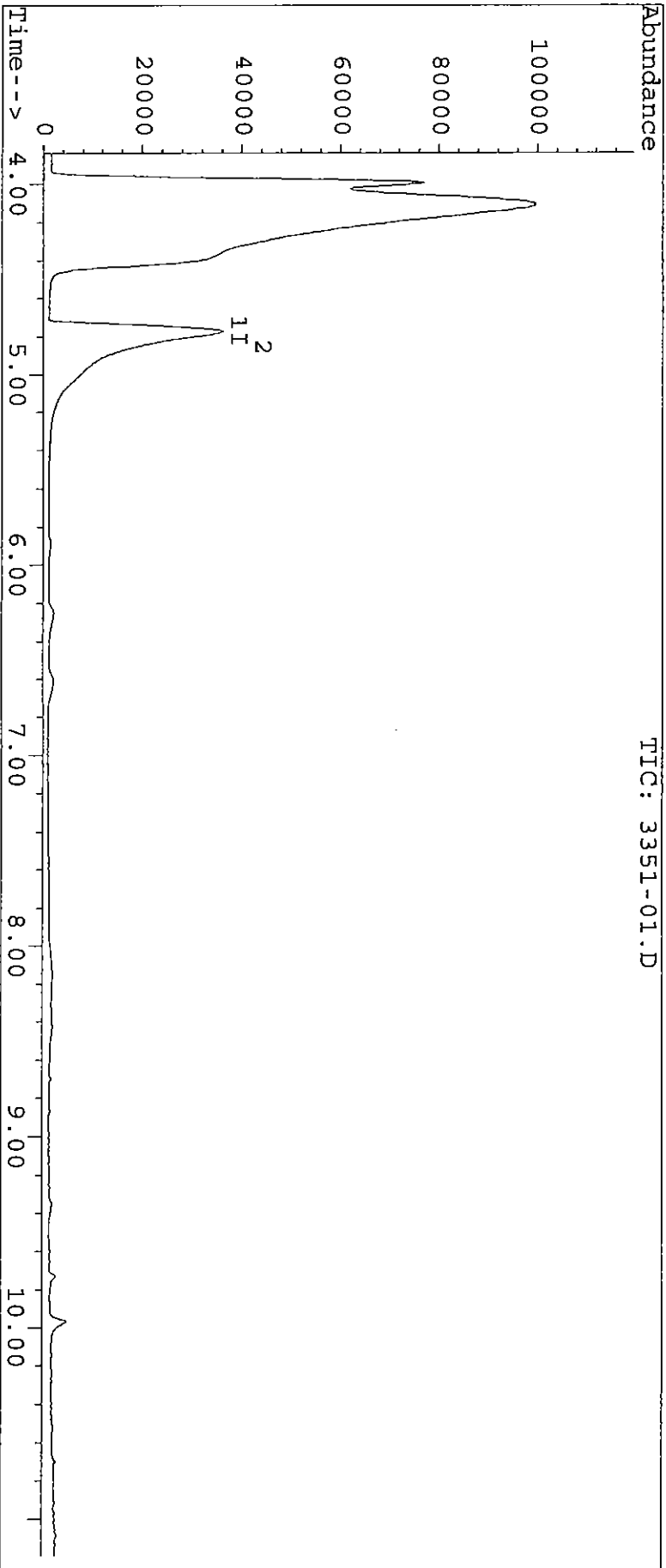
Jim S/20/03

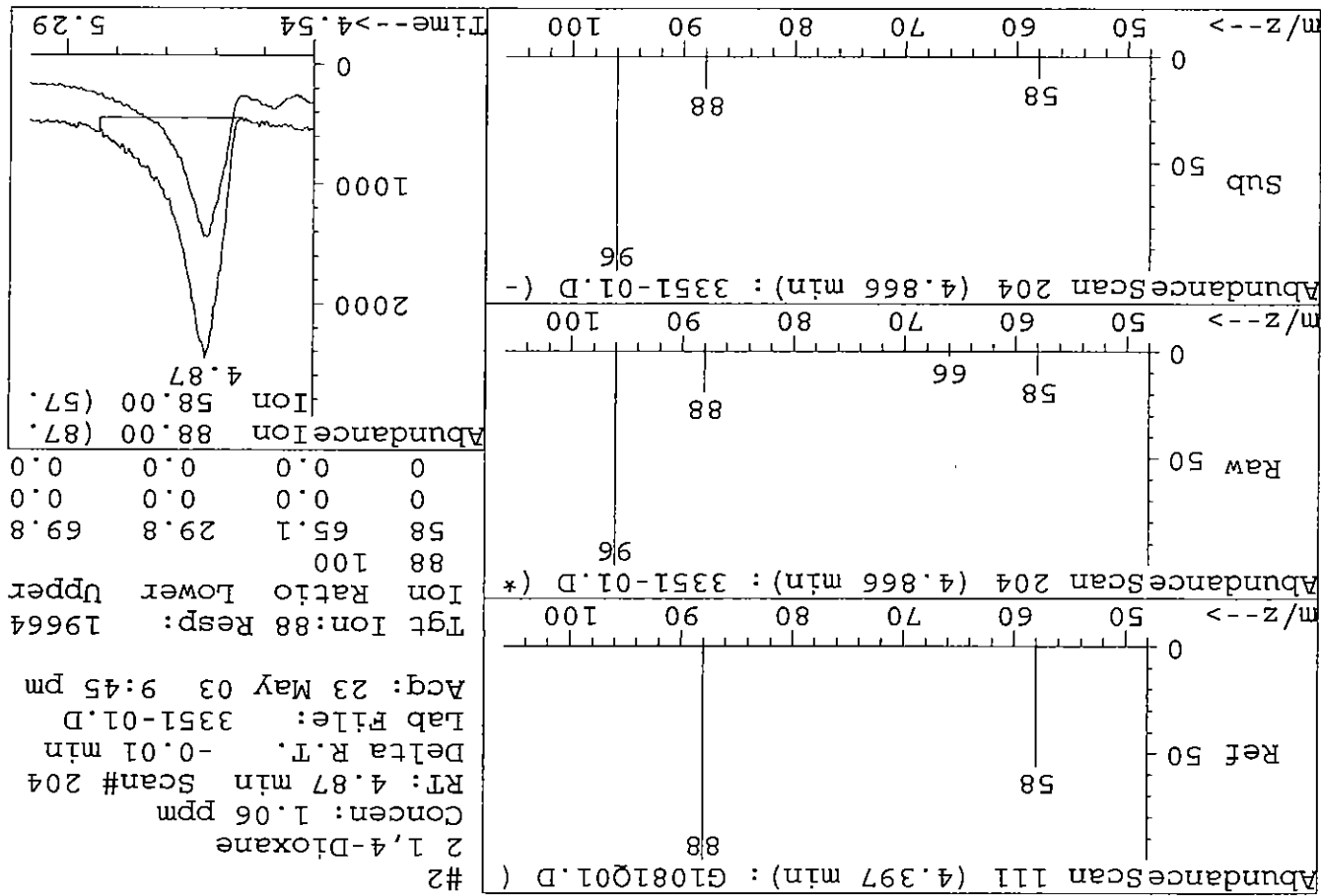
 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\3351-01.D
Acq On : 23 May 03 9:45 pm
Sample : F=.001
Misc :
Quant Time: May 27 13:50 2003
Vial: 2
Operator: Andy Huang
Inst : GC/MS - M
Multiplr: 1.00
Quant Results File: quant.res
Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Fri May 23 12:56:26 2003
Response via : Multiple Level Calibration





Applied P & Ch Laboratory
Organic Analysis Results for Method 8270-SIM

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/22/2003
Project ID: JPL GW-Mon. 2Q03.	Service ID: 33351	Collected by:
Sample ID: EB-14-5/22/03	Lab Sample ID: 03-3351-2	Received Date: 05/22/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 8270-SIM	Prep. Method: 3520	Instrument ID: GC/MS: M
Batch No: 03G2623	Prep. Date: 05/22/03	Anal. Date: 05/23/03
Data File Name: 3351-02	Prep. No: 1 of 1	Anal. Time: 22:05
Extract Vol. 1.0 mL	Sample Amount: 1000 mL	Dilution Factor: 1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	<1	U
Internal Standard				Control Limit, %	IS Rec. %	
1	1,4-DIOXANE-D8	17647-74-4		50-200	93	
	# 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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Organic Analysis Results for Method 8270-SIM

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Collection Date:	05/22/2003
Project ID:	JPL GW-Mon. 2Q03.	Service ID:	33351	Collected by:	
Sample ID:	MW-4-2	Lab Sample ID:	03-3351-3	Received Date:	05/22/2003
Sample Type:	Field Sample	Sample Matrix:	Water	Moisture %:	-
Anal. Method:	8270-SIM	Prep. Method:	3520	Instrument ID:	GC/MS: M
Batch No:	03G2623	Prep. Date:	05/22/03	Anal. Date:	05/23/03
Data File Name:	3351-03	Prep. No:	1 of 1	Anal. Time:	22:26
Extract Vol.	1.0 mL	Sample Amount:	1000 mL	Dilution Factor:	1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	1	
Internal Standard				Control Limit, %	IS Rec. %	
1	1,4-DIOXANE-D8	17647-74-4		50-200	93	
# of out-of-control					0	

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

Organic Analysis Results for Method 8270-SIM

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Collection Date:	05/22/2003
Project ID:	JPL GW-Mon. 2Q03.	Service ID:	33351	Collected by:	
Sample ID:	MW-4-3	Lab Sample ID:	03-3351-4	Received Date:	05/22/2003
Sample Type:	Field Sample	Sample Matrix:	Water	Moisture %:	--
Anal. Method:	8270-SIM	Prep. Method:	3520	Instrument ID:	GC/MS: M
Batch No:	03G2623	Prep. Date:	05/22/03	Anal. Date:	05/23/03
Data File Name:	3351-04	Prep. No:	1 of 1	Anal. Time:	22:47
Extract Vol.	1.0 mL	Sample Amount:	1000 mL	Dilution Factor:	1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	0.4	J
Internal Standard				Control Limit, %	IS Rec.%	
1	1,4-DIOXANE-D8	17647-74-4		50-200	85	
	# of out-of-control				0	

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

Organic Analysis Results for Method 8270-SIM

Client Name:	GEOFON, Inc.	Project No:	04-4428.10	Collection Date:	05/22/2003
Project ID:	JPL GW-Mon. 2Q03.	Service ID:	33351	Collected by:	
Sample ID:	MW-4-4	Lab Sample ID:	03-3351-5	Received Date:	05/22/2003
Sample Type:	Field Sample	Sample Matrix:	Water	Moisture %:	-
Anal. Method:	8270-SIM	Prep. Method:	3520	Instrument ID:	GC/MS: M
Batch No:	03G2623	Prep. Date:	05/22/03	Anal. Date:	05/23/03
Data File Name:	3351-05	Prep. No:	1 of 1	Anal. Time:	23:08
Extract Vol.	1.0 mL	Sample Amount:	1000 mL	Dilution Factor:	1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	<1	U
Internal Standard				Control Limit, %	IS Rec. %	
1	1,4-DIOXANE-D8	17647-74-4		50-200	96	
	# 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

Applied P & Ch Laboratory
Organic Analysis Results for Method 8270-SIM

Client Name: GEOFON, Inc.	Project No: 04-4428.10	Collection Date: 05/22/2003
Project ID: JPL GW-Mon. 2Q03.	Service ID: 33351	Collected by:
Sample ID: MW-4-5	Lab Sample ID: 03-3351-6	Received Date: 05/22/2003
Sample Type: Field Sample	Sample Matrix: Water	Moisture %: -
Anal. Method: 8270-SIM	Prep. Method: 3520	Instrument ID: GC/MS: M
Batch No: 03G2623	Prep. Date: 05/22/03	Anal. Date: 05/23/03
Data File Name: 3351-06	Prep. No: 1 of 1	Anal. Time: 23:29
Extract Vol. 1.0 mL	Sample Amount: 1000 mL	Dilution Factor: 1

#	Component Name	CAS No	Unit	RL	Result	Qualifier
1	1,4-DIOXANE	123-91-1	µg/L	1	<1	U
Internal Standard				Control Limit, %	IS Rec.%	
1	1,4-DIOXANE-D8	17647-74-4		50-200	99	
# of out-of-control					0	

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

Qualifier: U - Not Detected or less than MDL 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
--	---

FORM-3C

Applied P & Ch Laboratory

Lab Control Spike/Lab Control Spike Duplicate Recovery for Method 8270-SIM

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33351
Project ID: JPL GW-Mon. 2Q03.	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03G2623	
LCS Filename: G2623L01	Date Analyzed: 052303	Time Analyzed: 13:31
LCSD Filename: G2623J01	Date Analyzed: 052303	Time Analyzed: 13:54

Spiked Components	Unit	Spike Added	Concentration		LCS Rec% #	QC Limit, % REC
			Unspiked	LCS		
1,4-DIOXANE	µg/L	20	0	18.7	94	40-140
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	LCSD Concentration	LCSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
1,4-DIOXANE	µg/L	20	18.7	94	0	30	40-140
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623I01.D Sample : F=.001
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 13:31 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 23 14:28 2003 Multiplr: 0.001000
 Print Time : Fri May 23 14:28 2003
 Miscellaneous :

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

Internal Standards

1	1,4-Dioxane-d8	4.81	4.80	0.003	96	66	246.833	20.00		Dev(Min)	
---	----------------	------	------	-------	----	----	---------	-------	--	----------	--

System Monitoring Compounds (Surrogate)

Target Compounds	<<< I1 : ISTD ID = 1 >>>	4.89	4.88	0.003	88	58	280.604	18.71	18.7	80	Qvalue
2	2,1,4-Dioxane										

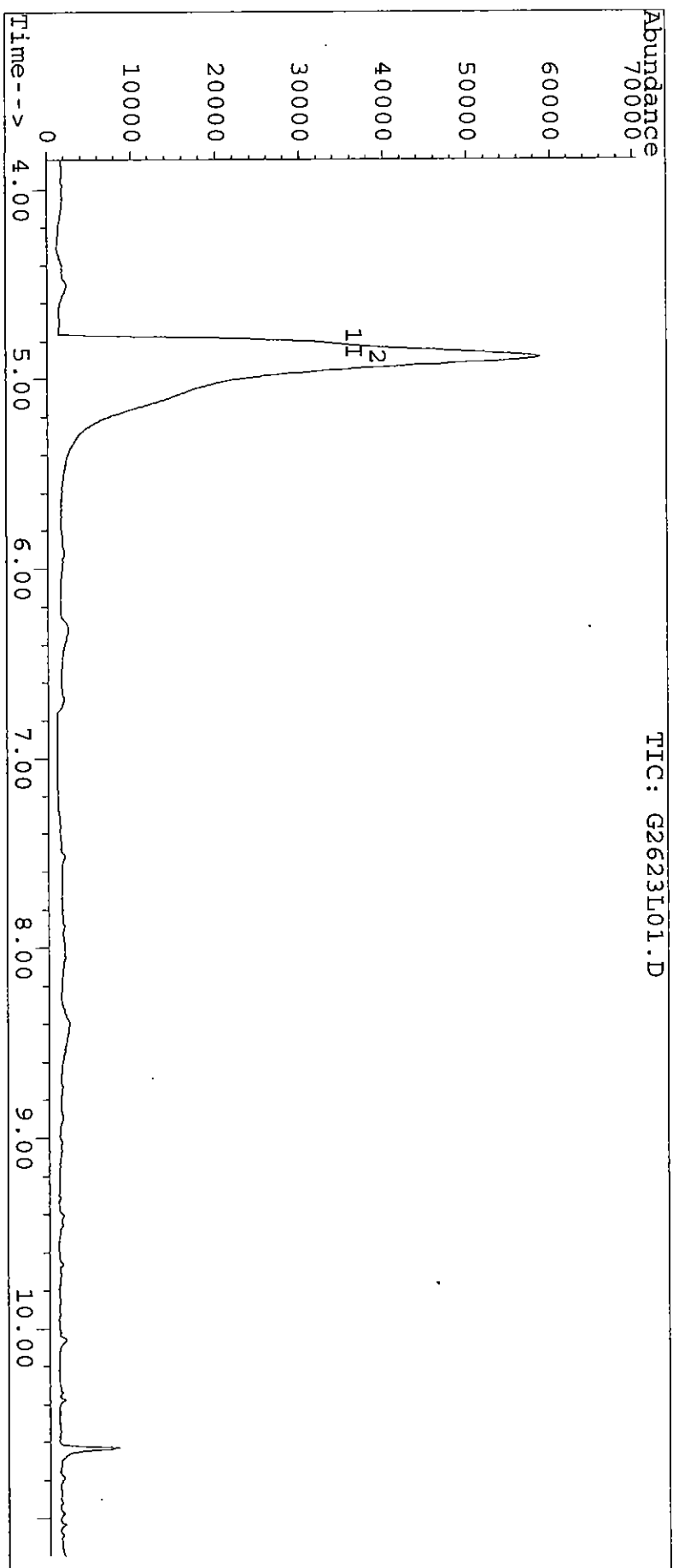
 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623L01.D Vial: 2
Acq On : 23 May 03 1:31 pm Operator: Andy Huang
Sample : F=.001 Inst : GC/MS - M
Misc : Multiplr: 1.00
Quant Time: May 23 14:28 2003 Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Fri May 23 12:56:26 2003
Response via : Multiple Level Calibration



Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623J01.D Sample : F=.001
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 13:54 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 23 14:28 2003 Multiplr: 0.001000
 Print Time : Fri May 23 14:28 2003
 Miscelaneous :

ID	Component Name	R.T.	RT0	DRRT	QIon	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
----	----------------	------	-----	------	------	----	---------	--------	-------	---------	------

Internal Standards

1 1.4-Dioxane-d8 4.83 4.80 0.006 96 66 231.813 20.00 Dev(Min) 0.03

System Monitoring Compounds (Surrogate)

Target Compounds

<<< I1 : ISTD ID = 1 >>>

2 2 1,4-Dioxane 4.91 4.88 0.006 88 58 263.999 18.74 18.7 80 %Recovery

Qvalue

 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623J01.D

Vial: 3

Acq On : 23 May 03 1:54 pm

Operator: Andy Huang

Sample : F=.001

Inst : GC/MS - M

Misc :

Multiplr: 1.00

Quant Time: May 23 14:28 2003

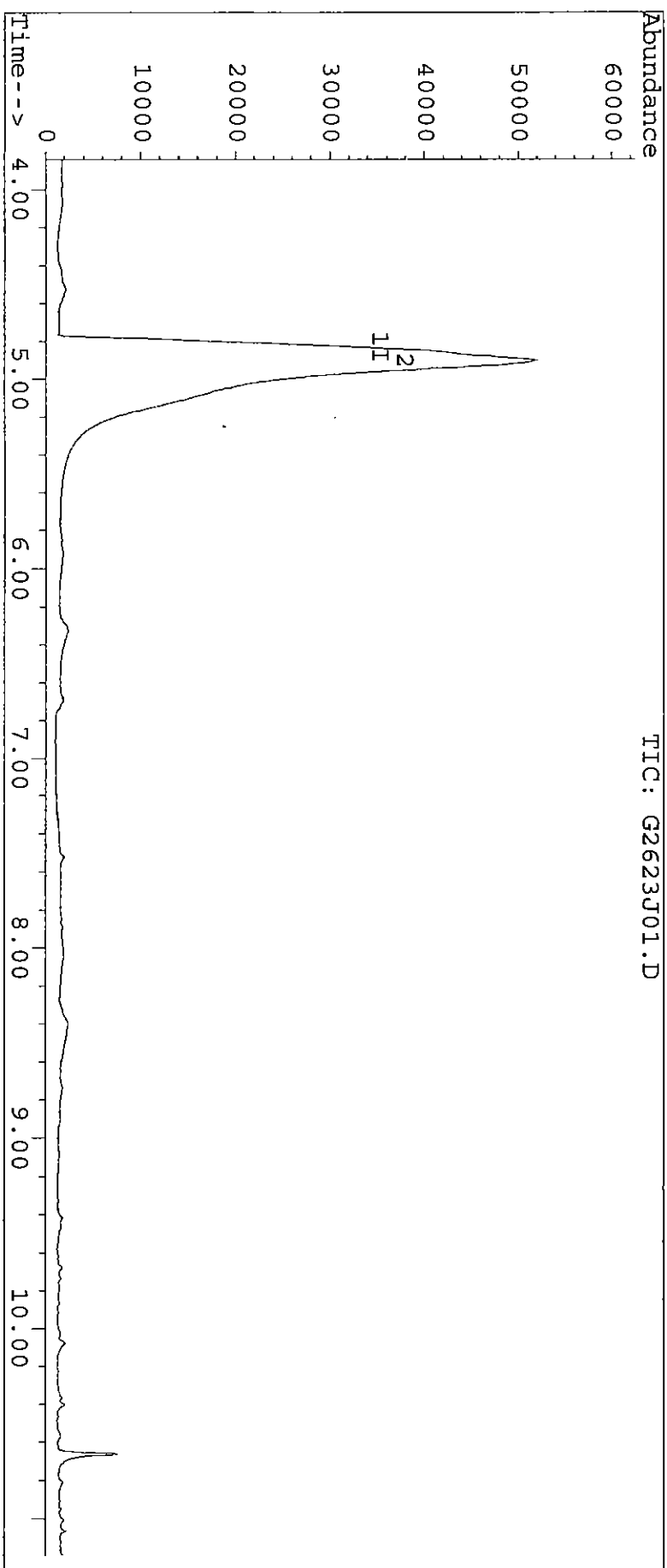
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Fri May 23 12:56:26 2003

Response via : Multiple Level Calibration



FORM-3C

Applied P & Ch Laboratory

Matrix Spike/Matrix Spike Duplicate Recovery for Method 8270-SIM

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33351
Project ID: JPL GW-Mon. 2Q03.	Project No: 04-4428.10	Sample Matrix: Water
	Batch No: 03G2623	
MS Filename: G2623M02	Date Analyzed: 052303	Time Analyzed: 15:01
MSD Filename: G2623N02	Date Analyzed: 052303	Time Analyzed: 15:23
MS Sample No: 04MW-01-1-GW	Sample Lab ID: 03-3318-4	

Spiked Components	Unit	Spike Added	Concentration		MS Rec% #	QC Limit, % REC
			Unspiked	MS		
1,4-DIOXANE	µg/L	20	0	19.4	97	40-140
# of Out-of-control					0	

Spiked Components	Unit	Spike Added	MSD Concentration	MSD Rec% #	RPD% #	QC Limit, %	
						RPD	REC
1,4-DIOXANE	µg/L	20	19.4	97	0	30	40-140
# of Out-of-control				0	0		

Column to be used to flag recovery and RPD values:

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

Comments: _____

Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623M02.D Sample : f=.001 \$3318-04
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 15:01 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 23 15:46 2003 Multiplr: 0.001000
 Print Time : Fri May 23 15:46 2003
 Miscellaneous :

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

Internal Standards

Dev (Min)
0.03

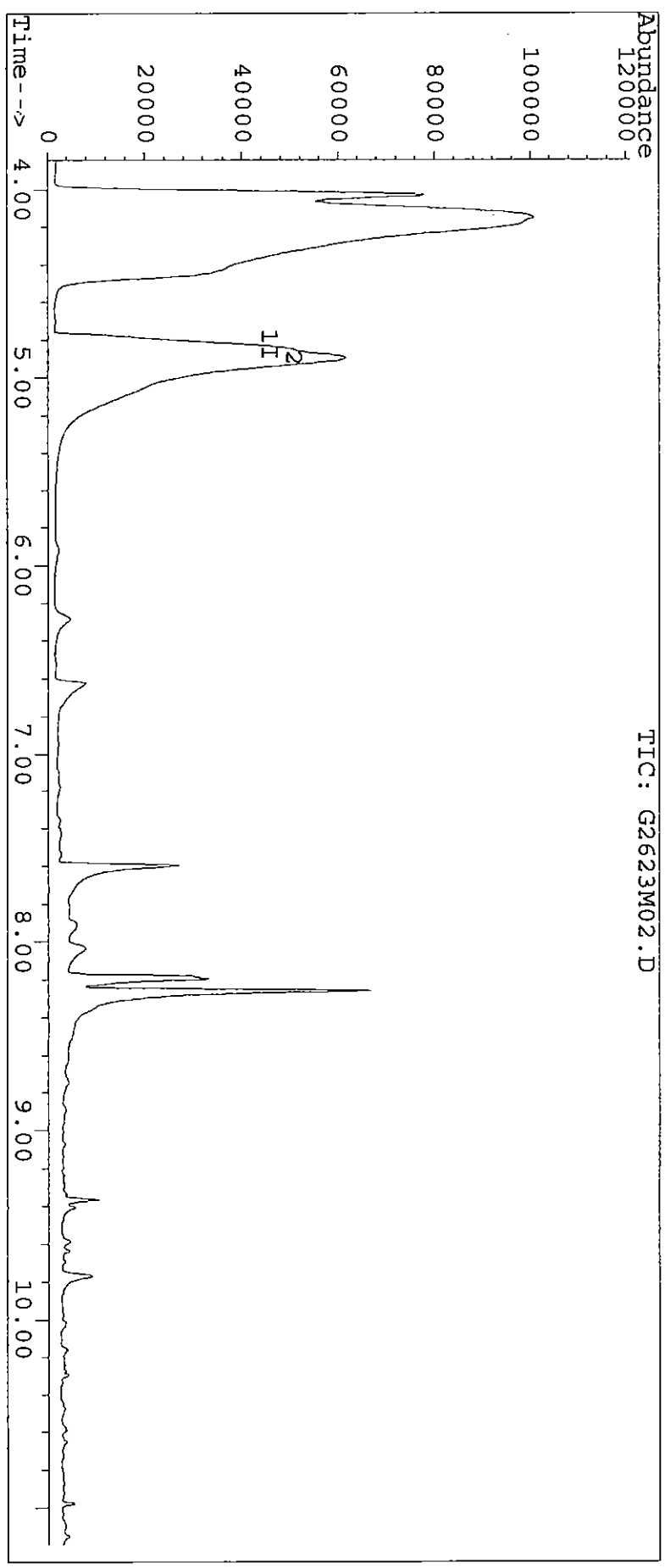
%Recovery

1	1,4-Dioxane-d8	4.82	4.80	0.005	96	66	266.456	20.00				
System Monitoring Compounds (Surrogate)												
Target Compounds												
<<<	I1 : ISTD ID = 1	>>>										Qvalue
2	2,1,4-Dioxane	4.90	4.88	0.005	88	58	314.103	19.40	19.4	84		

= qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623M02.D Vial: 6
Acq On : 23 May 03 3:01 pm Operator: Andy Huang
Sample : F=.001 \$3318-04 Inst : GC/MS - M
Misc : Quant Results File: quant.res
Quant Time: May 23 15:46 2003
Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Fri May 23 12:56:26 2003
Response via : Multiple Level Calibration



Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623N02.D Sample : F=.001 \$3318-04
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 15:23 2003
 Method Update: Fri May 23 12:56 2003 RF via : Multiple Level Calibration
 Quant. Time : May 23 15:46 2003 Operator: Andy Huang
 Print Time : Fri May 23 15:46 2003 Multiplr: 0.001000
 Miscellaneous :

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

Internal Standards

Dev (Min)
0.02

System Monitoring Compounds (Surrogate)

%Recovery

Target Compounds	<<< I1 : ISTD ID = 1 >>>	4.90	4.88	0.005	88	58	319.174	19.36	19.4	85	Qvalue
------------------	--------------------------	------	------	-------	----	----	---------	-------	------	----	--------

= qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623N02.D

Vial: 7

Acq On : 23 May 03 3:23 pm

Operator: Andy Huang

Sample : F=.001 \$3318-04

Inst : GC/MS - M

Misc :

Multiplr: 1.00

Quant Time: May 23 15:46 2003

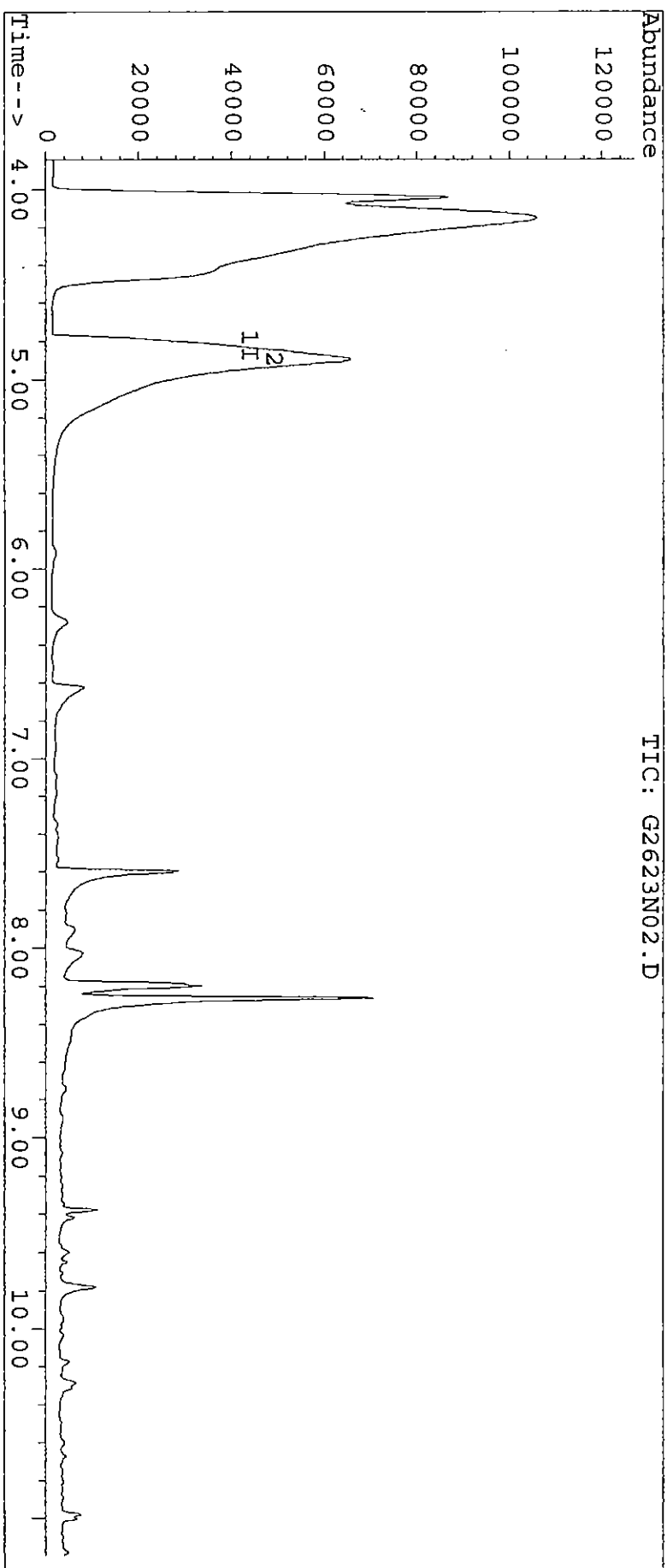
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Fri May 23 12:56:26 2003

Response via : Multiple Level Calibration



TIC: G2623N02.D

FORM-4B

Applied P & Ch Laboratory

Method Blank Summary for Method 8270-SIM

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCL
Case No:	SAS No:	Service ID: 33351
Project ID: JPL GW-Mon. 2Q03.	Project No: 04-4428.10	Analysis Date: 05/23/03
Sample ID: 03G2623-MB-01	Sample Matrix: Water	Analysis Time: 13:10
Lab Sample ID: 03G2623-MB-01	Batch No: 03G2623	Instrument ID: GC/MS: M
	Data File Name: G2623K01	GC Column: DB-5.625
		Column ID: 0.25 mm

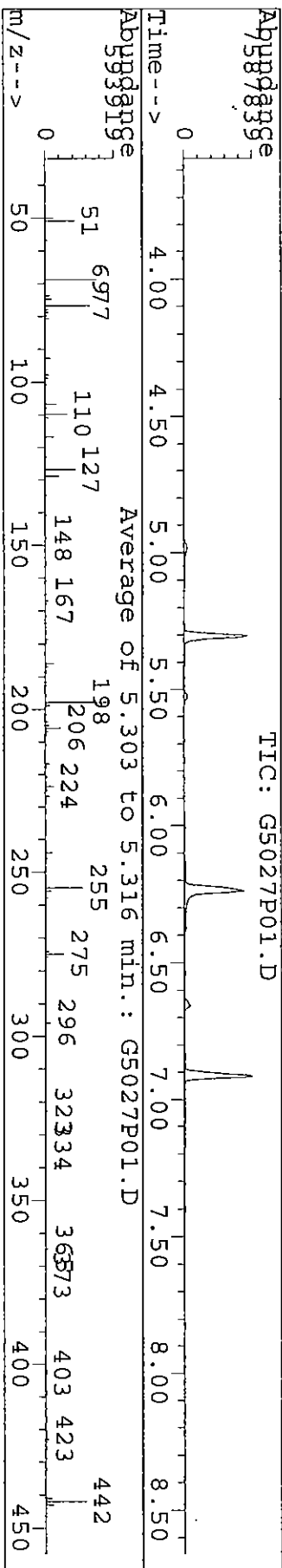
This Method Blank applies to the following samples and QC samples:

#	Client Sample No	Lab Sample ID	Sample Type	Data Filename	Analysis Date	Analysis Time
1	03G2623-LCS-01	03G2623-LCS-01	Lab Control Spike	G2623L01	05/23/03	13:31
2	03G2623-LSD-01	03G2623-LSD-01	Lab Control Spike Duplicate	G2623J01	05/23/03	13:54
3	04MW-01-1-GWMS	03-3318-4MS	Matrix Spike	G2623M02	05/23/03	15:01
4	04MW-01-1-GWMSD	03-3318-4MSD	Matrix Spike Duplicate	G2623N02	05/23/03	15:23
5	DUPE-8-2Q03	03-3351-1	Field Sample	3351-01	05/23/03	21:45
6	EB-14-5/22/03	03-3351-2	Field Sample	3351-02	05/23/03	22:05
7	MW-4-2	03-3351-3	Field Sample	3351-03	05/23/03	22:26
8	MW-4-3	03-3351-4	Field Sample	3351-04	05/23/03	22:47
9	MW-4-4	03-3351-5	Field Sample	3351-05	05/23/03	23:08
10	MW-4-5	03-3351-6	Field Sample	3351-06	05/23/03	23:29
11						
12						
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						
23						
24						
25						

Data File : C:\HPCHEM\1\DATA\02G5027\G5027P01.D
 Acq On : 17 Dec 02 2:19 pm
 Sample : ##02G5027,w dftpp gc14349
 Misc :

Vial: 99
 Operator: Andy Huang
 Inst : GC/MS - M
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
 Title : * Applied P & Ch Lab * GC/MS 8270



Peak Apex is scan: AVERAGE

Target Mass	Rel. Mass	to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result
51	198	69	30	80	44.9	254307	PASS
68	69	198	0	2	0.0	0	PASS
69	198	69	1	100	78.3	443243	PASS
70	69	198	0	2	0.3	1277	PASS
127	198	69	25	75	46.2	261139	PASS
197	198	198	0	1	0.0	0	PASS
198	198	198	100	100	100.0	565803	PASS
199	198	198	5	9	6.7	37981	PASS
275	198	198	10	30	27.4	155040	PASS
365	198	198	1	100	3.7	20870	PASS
441	443	198	1	99	78.6	52501	PASS
442	198	443	40	110	61.8	349717	PASS
443	442	198	15	24	19.1	66784	PASS

5B
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name : APPLIED P & CH LAB Contract: _____
 Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: _____
 Lab File ID: G5027P01 DFTPP Injection Date: 12/17/02
 Instrument ID: GCMS-M DFTPP Injection Time: 1419

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30.0 - 60.0% of mass 198	44.9
68	Less than 2.0% of mass 69	0.0 (0.0)1
69	Mass 69 relative abundance	78.3
70	Less than 2.0% of mass 69	0.2 (0.3)1
127	40.0 - 60.0% of mass 198	46.2
197	Less than 1.0% of mass 198	0.0
198	Base Peak, 100 % relative abundance	100.0
199	5.0 - 9.0% of mass 198	6.7
275	10.0 - 30.0% of mass 198	27.4
365	Greater than 0.75% of mass 198	3.7
441	Present, but less than mass 443	9.3 (78.6)3
442	40.0 - 100.0% of mass 198	61.8
443	17.0 - 23.0% of mass 442	11.8 (19.1)2

1-Value is % mass 69 2-Value is % mass 442 3-Value is % mass 443

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	SSTD040	DS006-40	DS006-040.D	12/17/02	1438
02	SSTD030	DS006-30	DS006-030.D	12/17/02	1457
03	SSTD020	DS006-20	DS006-020.D	12/17/02	1516
04	SSTD010	DS006-10	DS006-10.D	12/17/02	1536
05	SSTD001	DS006-01	DS006-01.D	12/17/02	1555
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
 Title : * Applied P & Ch Lab * GC/MS 8270
 Last Update : Tue Dec 17 16:08:23 2002
 Response via : Initial Calibration

Calibration Files
 30 =M06-030.D 20 =M06-020.D 10 =M06-010.D
 1 =M06-001.D 40 =M06-040.D

Compound	30	20	10	1	40	Avg	%RSD
1) 1 1.4-Dioxane-d8							
2) 2 1,4-Dioxane	1.151	1.165	1.216	1.379	1.165	1.215	7.81

-----ISTD-----

(#) = Out of Range
 DIOSIM06.M Tue Dec 17 16:08:40 2002 5972

Data Filename: C:\HPCHEM\1\DATA\02G5027\M06-040.D Sample : 40ppm gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 14:38 2002 RF via : Multiple Level Calibration
 Method Update: Fri Oct 18 22:02 2002 Operator: Andy Huang
 Quant. Time : Dec 17 15:07 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 15:07 2002
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	QIon	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
Internal Standards											
1	1,4-Dioxane-d8	4.36	4.37	-0.001	96	66	477.763	20.00		Dev (Min)	
System Monitoring Compounds (Surrogate)											
Target Compounds											
<<<	I1 : ISTD ID = 1	>>>								Qvalue	
2	2,1,4-Dioxane	4.43	4.44	-0.003	88	58	1113.531	38.33	38334.5	76	#

 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

Data File : C:\HPCHEM\1\DATA\02G5027\M06-040.D

Vial: 1

Acq On : 17 Dec 02 2:38 pm

Operator: Andy Huang

Sample : 40ppm gcl14554

Inst : GC/MS - M

Misc :

Multiplr: 1.00

Quant Time: Dec 17 15:07 2002

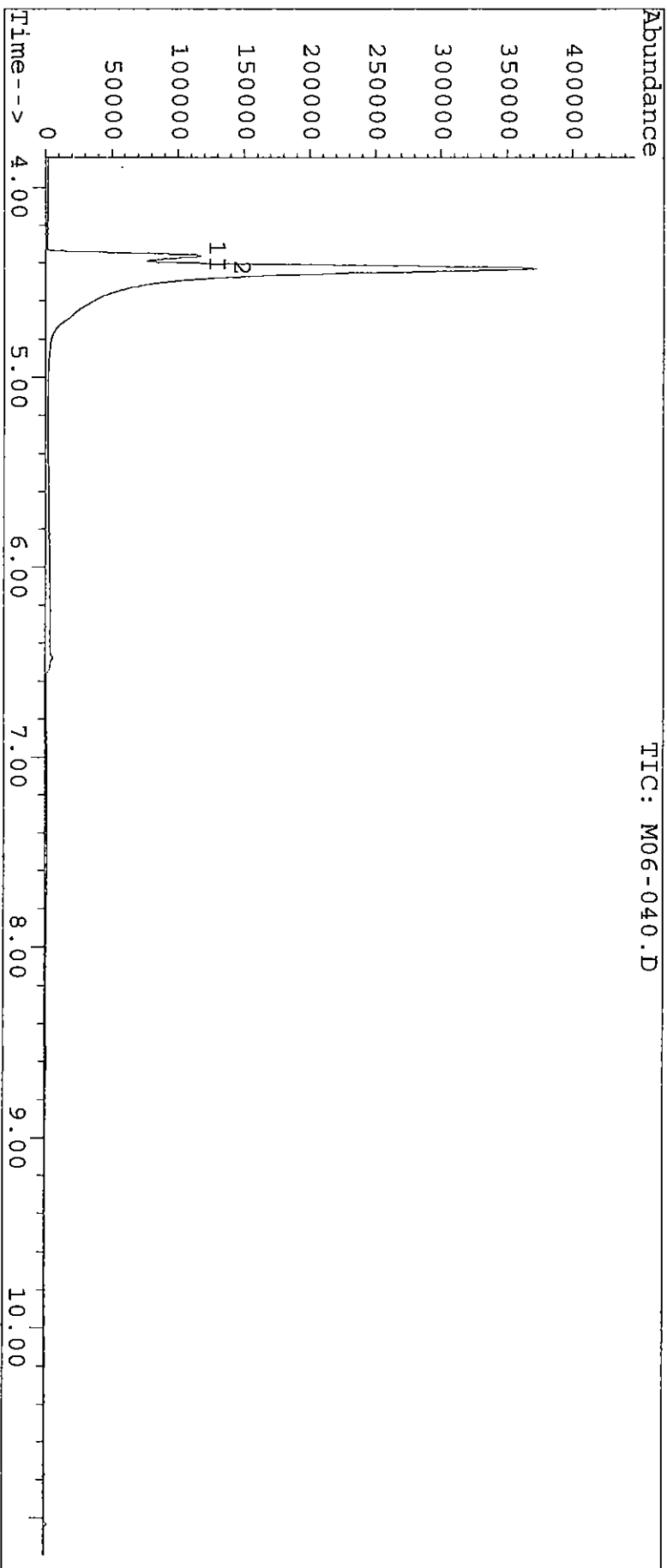
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Fri Oct 18 22:02:00 2002

Response via : Multiple Level Calibration



Data Filename: C:\HPCHEM\1\DATA\02G5027\M06-030.D Sample : 30ppm gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 14:57 2002 RF via : Multiple Level Calibration
 Method Update: Tue Dec 17 15:08 2002 Operator: Andy Huang
 Quant. Time : Dec 17 15:17 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 15:17 2002
 Miscleaneous :

ID	Component Name	R.T.	RT0	DRRT	QIon	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
1	Internal Standards										
1	1,4-Dioxane-d8	4.37	4.36	0.000	96	66	513.693	20.00		Dev (Min)	0.00

System Monitoring Compounds (Surrogate) %Recovery

Target Compounds	Qvalue
<<< I1 : ISTD ID = 1 >>>	
2 2 1,4-Dioxane	4.44 4.43 0.002 88 58 886.552 28.47 28474.2 98

 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\02G5027\M06-030.D

Vial: 2

Acq On : 17 Dec 02 2:57 pm

Operator: Andy Huang

Sample : 30ppm gcl4554

Inst : GC/MS - M

Misc :

Multiplr: 1.00

Quant Time: Dec 17 15:17 2002

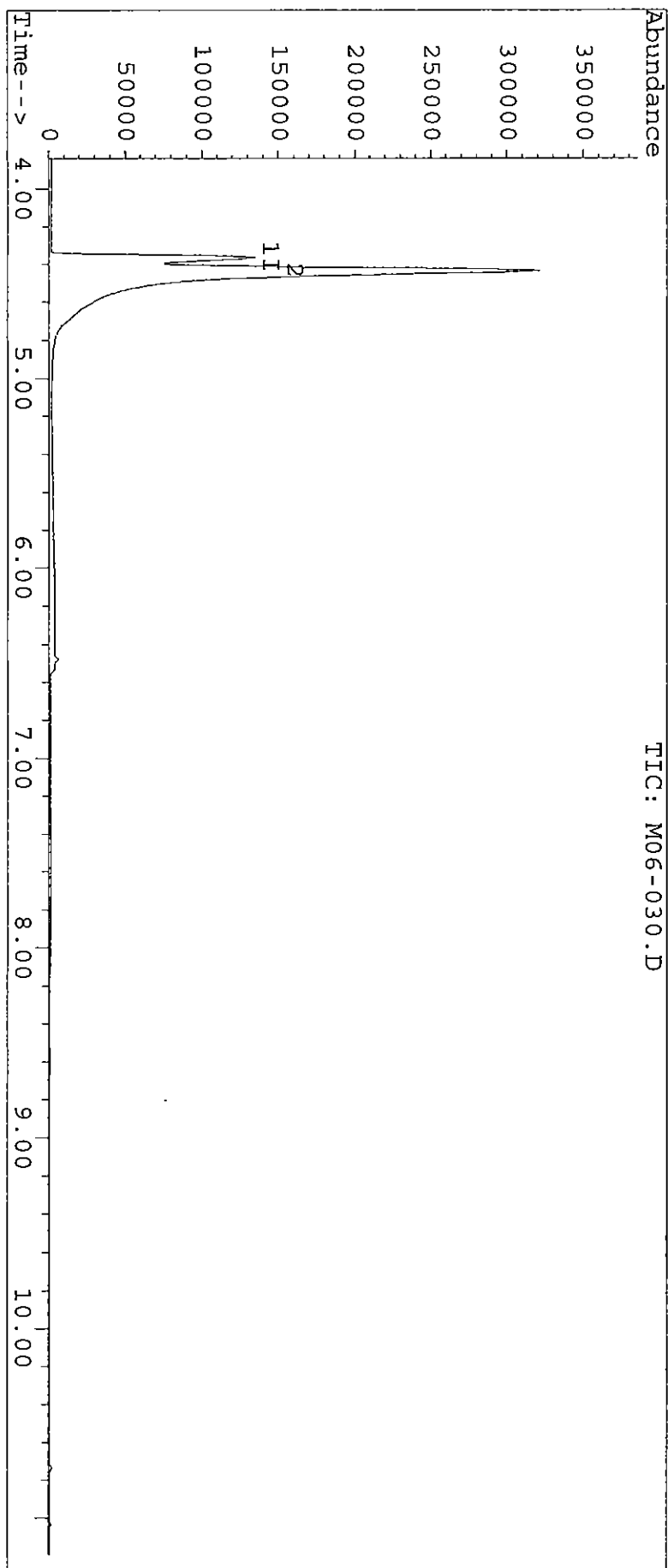
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Tue Dec 17 15:08:08 2002

Response via : Multiple Level Calibration



Data Filename: C:\HPCHEM\1\DATA\02G5027\M06-020.D Sample : 20ppm gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 15:16 2002 RF via : Multiple Level Calibration
 Method Update: Tue Dec 17 15:17 2002 Operator: Andy Huang
 Quant. Time : Dec 17 15:35 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 15:35 2002
 Miscleaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
Internal Standards											
1	1,4-Dioxane-d8	4.36	4.36	0.000	96	66	465.038	20.00		Dev (Min)	
										0.00	

System Monitoring Compounds (Surrogate)

%Recovery

Target Compounds	Qvalue
<<< I1 : ISTD ID = 1 >>>	
2 2 1,4-Dioxane	4.43 4.43 0.000 88 58 541.825 19.40 19401.4 97

= qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

Data File : C:\HPCHEM\1\DATA\02G5027\M06-020.D

Vial: 3

Acq On : 17 Dec 02 3:16 pm

Operator: Andy Huang

Sample : 20ppm gcl4554

Inst : GC/MS - M

Misc :

Multiplr: 1.00

Quant Time: Dec 17 15:35 2002

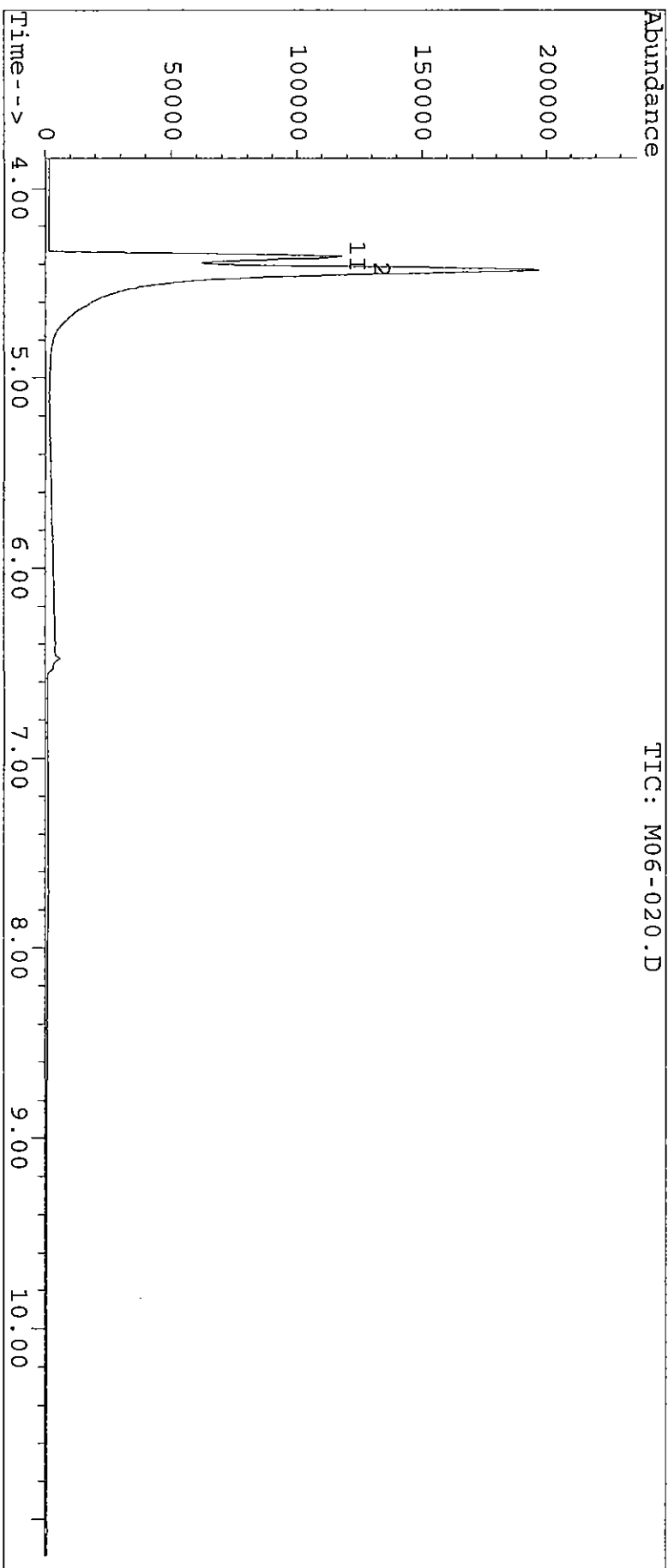
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Tue Dec 17 15:17:15 2002

Response via : Multiple Level Calibration



Data Filename: C:\HPCHEM\1\DATA\02G5027\M06-010.D Sample : 10ppm gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 15:36 2002 RF via : Multiple Level Calibration
 Method Update: Tue Dec 17 15:36 2002 Operator: Andy Huang
 Quant. Time : Dec 17 15:51 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 15:51 2002
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	QIon	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
----	----------------	------	-----	------	------	----	---------	--------	-------	---------	------

Internal Standards
 1 1.4-Dioxane-d8 4.36 4.36 0.000 96 66 464.928 20.00 Dev(Min) 0.00

System Monitoring Compounds (Surrogate)
 %Recovery

Target Compounds
 <<< I1 : ISTD ID = 1 >>> Qvalue

2 2 1,4-Dioxane 4.43 4.43 0.000 88 58 282.615 10.16 10158.2 100

= qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\02G5027\M06-010.D

Acq On : 17 Dec 02 3:36 pm

Sample : 10ppm gc14554

Misc :

Quant Time: Dec 17 15:51 2002

Vial: 4

Operator: Andy Huang

Inst : GC/MS - M

Multiplr: 1.00

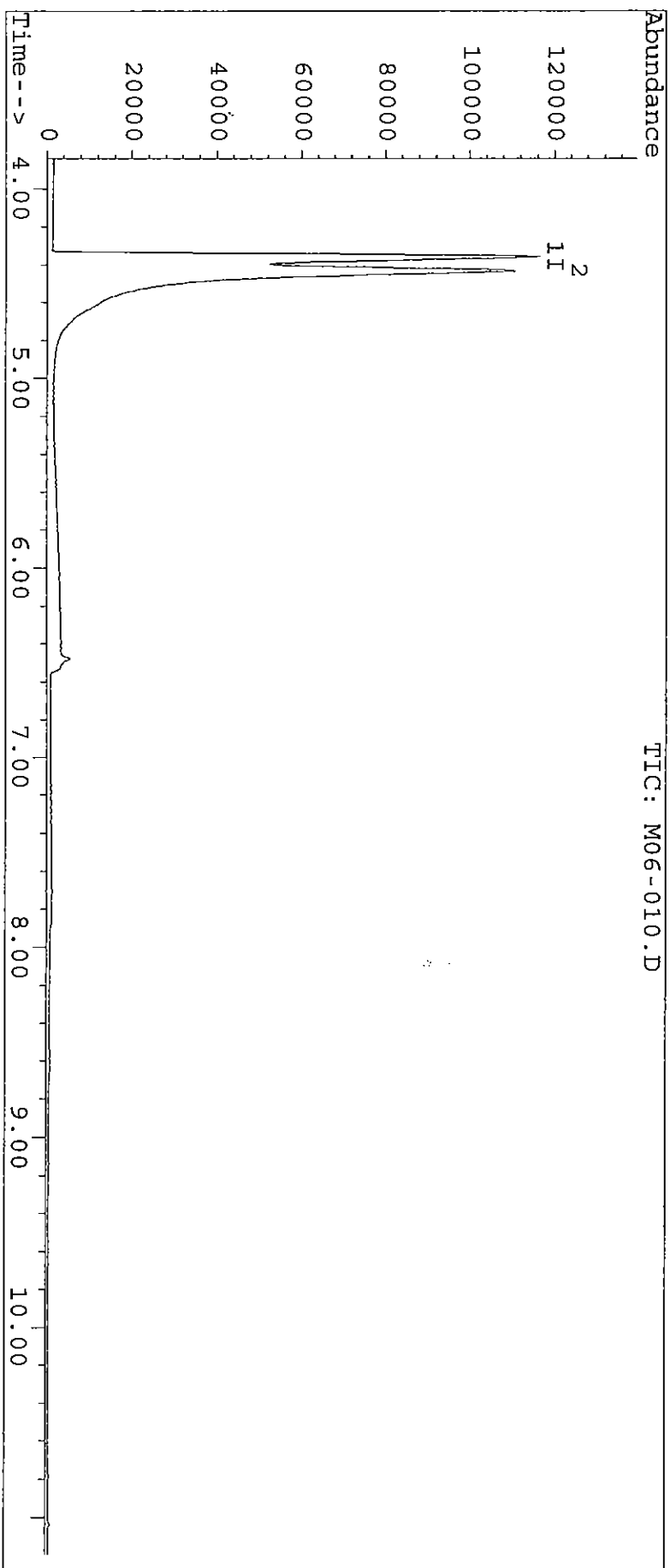
Quant Results File: quant.res

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Tue Dec 17 15:36:28 2002

Response via : Multiple Level Calibration



M06-010.D DIOSIM06.M

Tue Dec 17 15:51:30 2002

5972

Data Filename: C:\HPCHEM\1\DATA\02G5027\M06-001.D Sample : 1ppm gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 15:55 2002 RF via : Multiple Level Calibration
 Method Update: Tue Dec 17 16:00 2002 Operator: Andy Huang
 Quant. Time : Dec 17 16:08 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 16:08 2002
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
1	1,4-Dioxane-d8	4.38	4.36	0.004	96	66	515.201	20.00		Dev (Min) 0.02	

System Monitoring Compounds (Surrogate) %Recovery

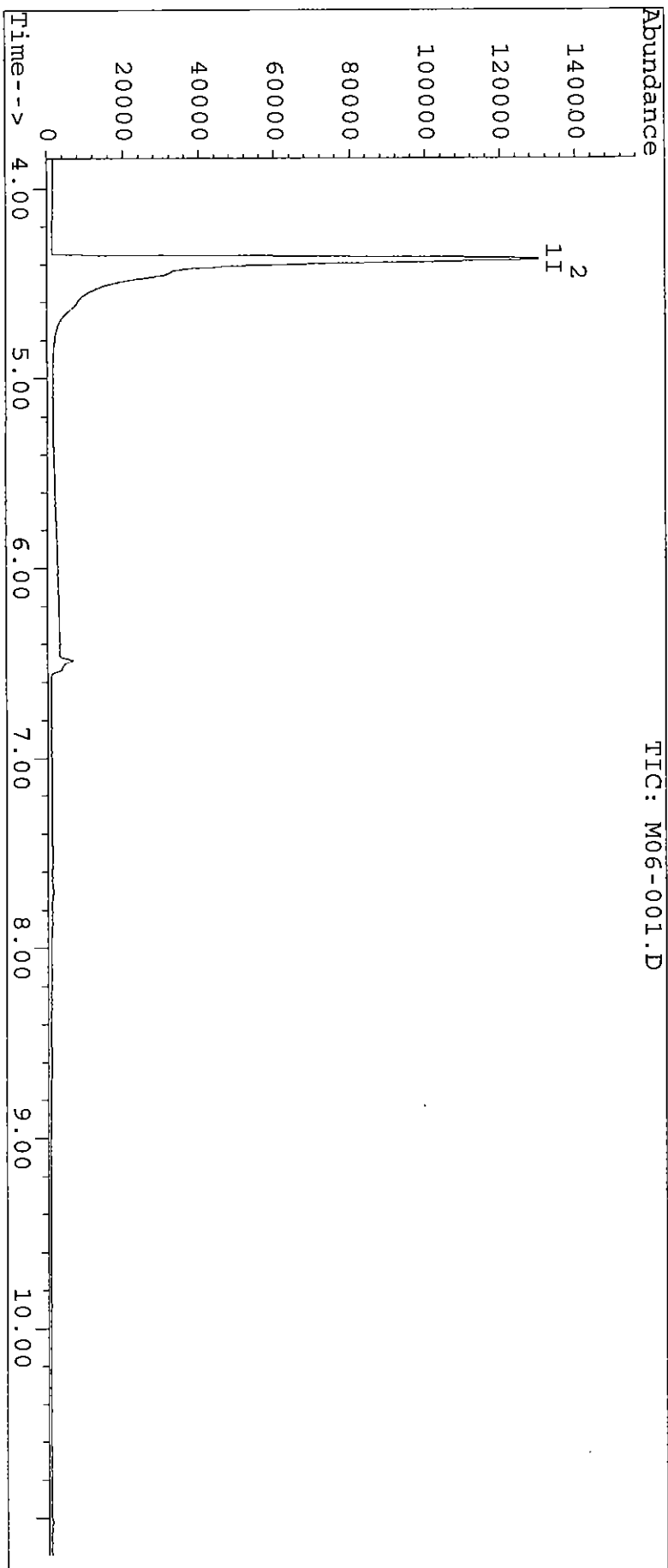
Target Compounds	<<< I1 : ISTD ID = 1 >>>	4.46	4.43	0.007	88	58	35.522	1.14	1137.6	98	Qvalue
2	2,1,4-Dioxane										

 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

 Page 1

Quantitation Report

Data File : C:\HPCHEM\1\DATA\02G5027\M06-001.D Vial: 5
Acq On : 17 Dec 02 3:55 pm Operator: Andy Huang
Sample : 1ppm gc14554 Inst : GC/MS - M
Misc : Quant Results File: quant.res
Quant Time: Dec 17 16:08 2002
Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Tue Dec 17 16:00:39 2002
Response via : Multiple Level Calibration



Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\02G5027\G5027Q01.D
Acq On : 17 Dec 02 4:14 pm
Sample : gcl4554
Misc :

Vial: 100
Operator: Andy Huang
Inst : GC/MS - M
Multiplier: 1.00

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Tue Dec 17 16:08:23 2002
Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 200%

Compound	AvgRRF	CCRF	%Dev Area	Dev (min)
1 I 1 1.4-Dioxane-d8	1.000	1.000	0.0	86 0.00
2 1,4-Dioxane	1.215	1.370	-12.7	102 0.00

(#) = Out of Range
G5027Q01.D DIOSIM06.M SPC's out = 0 CCC's out = 0
Tue Dec 17 16:30:02 2002 5972

Data Filename: C:\HPCHEM\1\DATA\02G5027\G5027Q01.D Sample : gc14554
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : Dec 17 16:14 2002 RF via : Multiple Level Calibration
 Method Update: Tue Dec 17 16:08 2002 Operator: Andy Huang
 Quant. Time : Dec 17 16:29 2002 Multiplr: 1.000000
 Print Time : Tue Dec 17 16:29 2002
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	Qion	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
Internal Standards											
1	1,4-Dioxane-d8	4.37	4.36	0.002	96	66	402.190	20.00		Dev(Min)	0.00

System Monitoring Compounds (Surrogate)

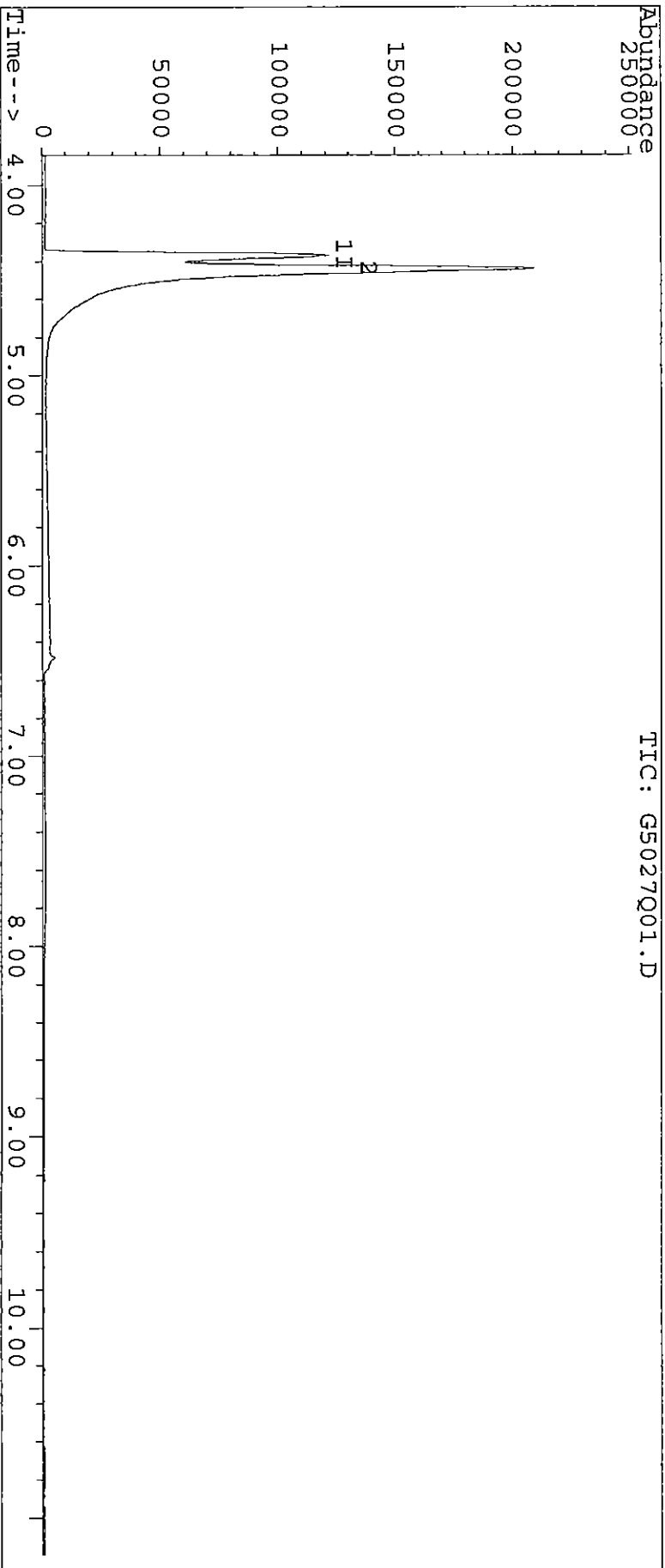
Target Compounds

<<<	I1 : ISTD ID = 1	>>>								Qvalue	
2	2 1,4-Dioxane	4.44	4.43	0.002	88	58	551.016	22.55	22549.4	98	

 # = qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

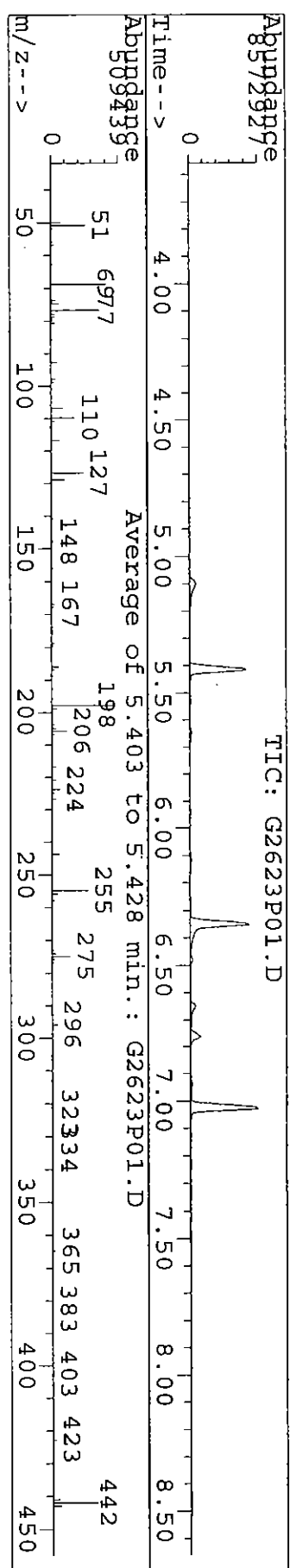
Data File : C:\HPCHEM\1\DATA\02G5027\G5027Q01.D
Acq On : 17 Dec 02 4:14 pm
Sample : gc14554
Misc :
Quant Time: Dec 17 16:29 2002
Vial: 100
Operator: Andy Huang
Inst : GC/MS - M
Multiplr: 1.00
Quant Results File: quant.res
Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Tue Dec 17 16:08:23 2002
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\03G2623\G2623P01.D
 Acq On : 23 May 03 12:09 pm
 Sample : ##03g2623,w dftpp gcl4349
 Misc :

Vial: 99
 Operator: Andy Huang
 Inst : GC/MS - M
 Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
 Title : * Applied P & Ch Lab * GC/MS 8270



Peak Apex is scan: AVERAGE

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result
51	198	30	80	53.8	261003	PASS
68	69	0	2	0.0	0	PASS
69	198	1	100	85.2	413160	PASS
70	69	0	2	0.1	536	PASS
127	198	25	75	50.2	243334	PASS
197	198	0	1	0.0	0	PASS
198	198	100	100	100.0	485182	PASS
199	198	5	9	6.8	32853	PASS
275	198	10	30	27.6	134080	PASS
365	198	1	100	3.9	18694	PASS
441	443	1	99	79.8	51403	PASS
442	198	40	110	69.7	338086	PASS
443	442	15	24	19.0	64403	PASS

FORM-5B

Applied P & Ch Laboratory

**Semivolatile Organic Instrument Performance Check for Method 8270-SIM
Decafluorotriphenylphosphine (DFTPP), Part II**

Client Name: GEOFON, Inc.	Contract No:	Lab Code: APCI
Case No:	SAS No:	Service ID: 033351
Project ID: JPL GW-Mon. 2Q03.	DFTPP Inj. Date: <u>05/23/03</u>	Batch No: 03G2623
	DFTPP Inj. Time: <u>12:09</u>	Sequence No: 03G2623
Project No: 04-4428.10	Instrument ID: M	GC Column: DB-5.625
Data File Name: G2623P01		Column ID: 0.25 mm

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS AND STANDARDS:

#	Client Sample No	Lab Sample ID	Data File Name	Date Analyzed	Time Analyzed
1	03G2623-CCV-01	03G2623-CCV-01	G2623Q01	05/23/03	12:48
2	03G2623-MB-01	03G2623-MB-01	G2623K01	05/23/03	13:10
3	03G2623-LCS-01	03G2623-LCS-01	G2623L01	05/23/03	13:31
4	03G2623-LSD-01	03G2623-LSD-01	G2623J01	05/23/03	13:54
5	04MW-01-1-GWMS	03-3318-4MS	G2623M02	05/23/03	15:01
6	04MW-01-1-GWMSD	03-3318-4MSD	G2623N02	05/23/03	15:23
7	DUPE-8-2Q03	03-3351-1	3351-01	05/23/03	21:45
8	EB-14-5/22/03	03-3351-2	3351-02	05/23/03	22:05
9	MW-4-2	03-3351-3	3351-03	05/23/03	22:26
10	MW-4-3	03-3351-4	3351-04	05/23/03	22:47
11	MW-4-4	03-3351-5	3351-05	05/23/03	23:08
12	MW-4-5	03-3351-6	3351-06	05/23/03	23:29
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					
25					

Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\03G2623\G2623Q01.D
Acq On : 23 May 03 12:48 pm
Sample : gcl5155
Misc :
Vial: 100
Operator: Andy Huang
Inst : GC/MS - M
Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M
Title : * Applied P & Ch Lab * GC/MS 8270
Last Update : Fri May 23 12:56:26 2003
Response via : Multiple Level Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 200%

Compound	AvgRRF	CCRF	%Dev	Area%	Dev(min)
1 1.4-Dioxane-d8	1.000	1.000	0.0	69	0.00
2 1,4-Dioxane	1.215	1.157	4.8	69	0.00

(#) = Out of Range
G2623Q01.D DIOSIM06.M
SPCC's out = 0
CCC's out = 0
Fri May 23 14:56:39 2003
5972

Data Filename: C:\HPCHEM\1\DATA\03G2623\G2623Q01.D Sample : gc15155
 Method : C:\HPCHEM\1\METHODS\DIOSIM06.M Inst. : GC/MS - M
 Acq. Time : May 23 12:48 2003 RF via : Multiple Level Calibration
 Method Update: Fri May 23 12:56 2003 Operator: Andy Huang
 Quant. Time : May 23 14:56 2003 Multiplr: 1.000000
 Print Time : Fri May 23 14:56 2003
 Miscellaneous :

ID	Component Name	R.T.	RT0	DRRT	QIon	Q1	RF/1000	C0,ppm	C,ppb	Quality	Note
----	----------------	------	-----	------	------	----	---------	--------	-------	---------	------

Internal Standards

Dev (Min)
0.00

%Recovery

1	1,4-Dioxane-d8	4.80	4.80	0.000	96	66	322.849	20.00			
2	1,4-Dioxane	4.88	4.88	0.000	88	58	373.629	19.05	19047.7	75	

Qvalue

= qualifier out of range, m = manual integration, ? = RT coelution, * = DRRT > 0.06
 5972

Quantitation Report

Vial: 100

Operator: Andy Huang

Inst : GC/MS - M

Multiplier: 1.00

Quant Results File: quant.res

Data File : C:\HPCHEM\1\DATA\03G2623\G2623Q01.D

Acq On : 23 May 03 12:48 pm

Sample : gc15155

Misc :

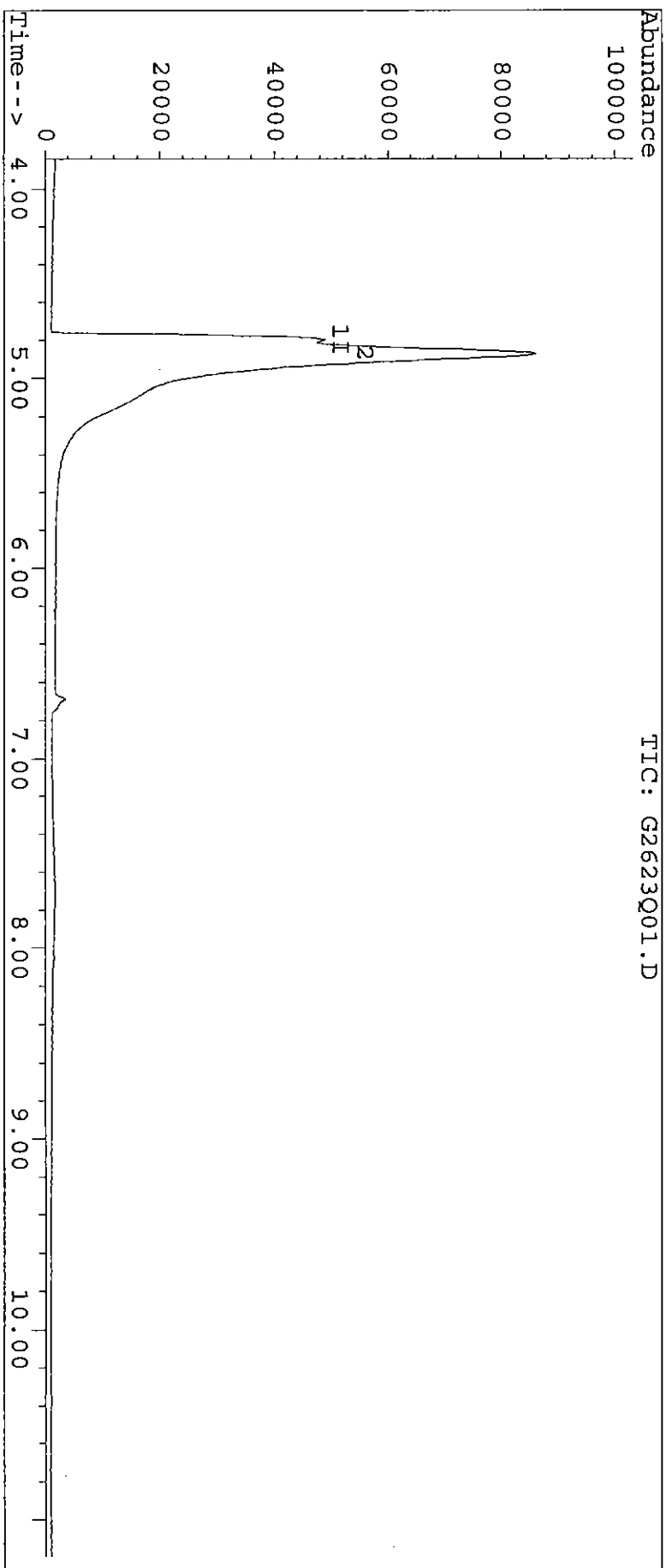
Quant Time: May 23 14:56 2003

Method : C:\HPCHEM\1\METHODS\DIOSIM06.M

Title : * Applied P & Ch Lab * GC/MS 8270

Last Update : Fri May 23 12:56:26 2003

Response via : Multiple Level Calibration



FORM-8B

Applied P & Ch Laboratory

Internal Standard Area and RT Summary for Method 8270-SIM

Client Name: GEOFON, Inc.

Contract No:

Lab Code: APCL

Case No:

SAS No:

Service ID: 033351

Project ID: JPL GW-Mon. 2Q03.

Project No: 04-4428.10

Sample Matrix: Water

CCV Data File: G2623Q01

Instrument ID: M

Batch No: 03G2623

#	Client Sample No	Lab Sample ID	Analysis Date & Time	IS-1	
				Area #	RT #
12 Hour CCV STD			05/23/03 12:48	322849	4.80
CCV Upper Limit				645698	5.30
CCV Lower Limit				161424	4.30
1	03G2623-MB-01	03G2623-MB-01	05/23/03 13:10	296242	4.83
2	03G2623-LCS-01	03G2623-LCS-01	05/23/03 13:31	246833	4.81
3	03G2623-LSD-01	03G2623-LSD-01	05/23/03 13:54	231813	4.83
4	04MW-01-1-GWMS	03-3318-4MS	05/23/03 15:01	266456	4.82
5	04MW-01-1-GWMSD	03-3318-4MSD	05/23/03 15:23	271289	4.82
6	DUPE-8-2Q03	03-3351-1	05/23/03 21:45	306192	4.77
7	EB-14-5/22/03	03-3351-2	05/23/03 22:05	299480	4.78
8	MW-4-2	03-3351-3	05/23/03 22:26	301802	4.78
9	MW-4-3	03-3351-4	05/23/03 22:47	275682	4.81
10	MW-4-4	03-3351-5	05/23/03 23:08	310393	4.79
11	MW-4-5	03-3351-6	05/23/03 23:29	320050	4.79
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					

IS-1 = 1,4-DIOXANE-D8

Area Upper Limit = +100% of CCV internal standard area

Area Lower Limit = - 50% of CCV internal standard area

RT Upper Limit = +0.50 minutes of CCV internal standard RT

RT Lower Limit = - 0.50 minutes of CCV internal standard RT

Column used to flag values outside QC limits with an asterisk.

* Values outside of QC limits

Applied P & Ch Laboratory

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Organic Sample Preparation Logbook

Batch # D3G 2623 Matrix: W Target method: 1,4 Dioxane EXT Method: 3520C

Date: 5-22-03

Solvent: CH₂Cl₂ Lot # 027471; Exchange Solvent: N/A Lot # N/A

Analyst: HL

Lot #: Surr.: GC-NA Conc.: NA ppm Na₂SO₄: Rc0078

Op. #	Sample Type	Sample ID #	Smpl Amnt (mL or g)	Surr. Vol 1/2 mL	Extract sub-ID	Final Vol. mL	Dilution F=V1/X	Note & Anomaly
3431	MB	03G2623-K01	1000.0	NA	1,4 Dioxane	1.00	0.001	
3432	LCS	-201	1000.0					
3433	LCSD	-201	1000.0					
3434	Sample-1	3318-1	1000.0					
3435	MS	-1	1000.0					
3436	MSD	-1	1000.0					
3437	Sample-2	3318-2	1000.0					
3438	Sample-3	-3	1000.0					
3439	Sample-4	-4	1000.0					
3440	Sample-5	-5	1000.0					
3441	Sample-6	-6	1000.0					
3442	Sample-7	-7	1000.0					
3443	Sample-8	3301-2	1000.0					
3444	Sample-9	3304-1	1000.0					
3445	Sample-10	3334-1	1000.0					
3446	Sample-11	-2	1000.0					
3447	Sample-12	-3	1000.0					
3448	Sample-13	-4	1000.0					
3449	Sample-14	-5	1000.0					
3450	Sample-15	3351-1	1000.0					
3451	Sample-16	-2	1000.0					
3452	Sample-17	-3	1000.0					
3453	Sample-18	-4	1000.0					
3454	Sample-19	-5	1000.0					
3455	Sample-20	-6	1000.0					
	MTX Dup.							
	LCS'							
	LCSD'							
	MS'	3318-4	1000.0	NA	1,4 Dioxane	1.00	0.001	
	MSD'	-4	1000.0	1/2				

5/22/03 HR

Op #	STD Lot #	C _{std} (ppm) x V _{std} (mL) / X(g or mL) = T	Op #	STD Lot #	C _{std} (ppm) x V _{std} (mL) / X(g or mL) = T
3432	GC-15154	1000 x 20.0 x 1.00 / X = ppb	3433	GC-15154	1000 x 20.0 x 1.00 / X = ppb
3435	GC	1000 x x / X = ppb	3436	GC	1000 x x / X = ppb
	GC	1000 x x / X = ppb		GC	1000 x x / X = ppb
3459	GC	1000 x x / X = ppb	3460	GC	1000 x x / X = 551 ppb

HL

Applied P & Ch Laboratory

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Organic Sample Preparation

Extraction method: 1 3510-separate funnel; 2 3520-L-L continue extraction; 3 3550-Ultrasonic; 4 LUFT-Shaker; 5 3540-Soxhlet;

6 Solid phase Extraction; 7 Microextraction; 8 SFE; 9 Dilution; 10 Concentration; 11 Other, specify.

Service ID: 3351 Batch # 262623 Matrix W GC, GC/MS Method 1,4-Dioxane Ext Method 3520C

Sample Extraction	OP. by <u>HL</u>	Date <u>5/20/03</u>	Surro Lot <u>GC-NA</u>	Conc. <u>NA</u> pp
	LCS Lot <u>GC-14768</u>	Conc. <u>20.0</u> pp	MS Lot <u>GC-14768</u>	Conc. <u>20.0</u> pp
	Ext. solv. <u>CH₂Cl₂</u>	Lot # <u>027471</u>	Exchange solv. <u>NA</u>	Lot # <u>NA</u>
Sample ID	<u>-1</u> <u>-2</u> <u>-3</u> <u>-4</u> <u>-5</u> <u>-6</u>			
Sample Matrix	<u>W</u>			
Sample Amount, X (g/mL)	<u>1000.0</u> <u>1000.0</u> <u>1000.0</u> <u>1000.0</u> <u>1000.0</u> <u>1000.0</u>			
Any TCLP, EP, WET, SPLP ?				
Surr. STD used, mL	<u>NA</u>			
Spike STD used, mL				
[Extracted at pH]	<u>1</u> <u>1</u> <u>1</u> <u>1</u> <u>1</u> <u>1</u>			
Solvent amount, mL	<u>200</u> <u>X</u> <u>X</u> <u>X</u> <u>X</u> <u>X</u>			<u>X</u> <u>X</u>
Final volume, V _i mL	<u>1000</u>			
Extraction DF f _e = V/X	<u>0.001</u>			
Sub-ID of extracts	<u>1,4-Dioxane</u>			
Anomaly Footnote [†] :				
Sample Cleanup	Op. by _____	Date <u>/ /</u>	(For details, see Cleanup worksheet)	
Cleanup method:	_____			
Cleanup DF, f _c	_____			
Preparation DF, f ₁ = f _e f _c	_____			

Pre-injection*:	Op. by _____	Date <u>/ /</u>	IS Lot : <u>GC-</u>	Conc. _____ pp
2nd dilution, f ₂ (1)	_____			
V _i , μL (40)	_____			
Single Extract GC, GC/MS Analysis:				
V ₂ = V _i /f ₂ , μL (40/f ₂)	_____			
V _{solvent} = V _i - V ₂ , μL	_____			
Double Extract GC, GC/MS Analysis:				
V _{>} = 1/2 V _i /f ₂ , μL (20/f ₂)	_____			
V _{<} = 1/2 V _i f _{<} /f _{>} , μL (20 f _{<} /f _{>})	_____			
V _{solvent} = V _i - V _{>} - V _{<}	_____			
IS volume, μL (10)	_____			
Final Conc. of IS (40 ppm)	_____			
Total F = 2 f ₁ f _{>} or f ₁ f ₂	_____			

[†] Extraction Anomaly Footnote:

1. Difficult to concentrate to 1.0 mL.	5. Too much precipitate during extraction.	9. Sample has strong stink
2. Extract is dark color	6. There is an oil layer above the sample.	10. Some solvent suspected
3. Extract is sticky	7. Mixture of soil & water.	11. special color, specify
4. Heterogenous matrix.	8. BN-extract is interfered, A-extract only.	

* For double extract analysis (ABN), f_> = max[f_A, f_B] and f_< = min[f_A, f_B], where f_A, f_B are f₁ for A and BN extracts, respectively. The values in () is the typical case for 625/8270 ABN analysis.

Organic Sample Preparation

Extraction method: 1 3510-separate funnel; 2 3520-L-L continue extraction; 3 3550-Ultrasonic; 4 LUFT-Shaker; 5 3540-Soxhlet;
 6 Solid phase Extraction; 7 Microextraction; 8 SFE; 9 Dilution; 10 Concentration; 11 Other, specify.

Service ID: 3318 Batch # 0362623 Matrix W GC, GC/MS Method 14-Dioxane Ext Method 3520C

Sample Extraction	OP. by <u>ide</u>	Date <u>5/22/03</u>	Surro Lot <u>GC-NA</u>	Conc. <u>NA</u> pp
	LCS Lot <u>GC-14768</u>	Conc. <u>20.0</u> pp	MS Lot <u>GC-14768</u>	Conc. <u>20.0</u> pp
	Ext. solv. <u>CH₂Cl₂</u>	Lot # <u>027471</u>	Exchange solv. <u>NA</u>	Lot # <u>NA</u>
Sample ID	<u>-K01 -L01</u>	<u>-J01 -I MS</u>	<u>-1 MSD -4MS</u>	<u>-4 MSD -1</u>
Sample Matrix	<u>W</u>			
Sample Amount, X (g/mL)	<u>1000.0</u>	<u>1000.0</u>	<u>1000.0</u>	<u>1000.0</u>
Any TCLP, EP, WET, SPLP ?	_____			
Surr. STD used, mL	<u>NA</u>			
Spike STD used, mL	<u>100</u>			
[Extracted at pH]	<u>1</u>	<u>1</u>	<u>1</u>	<u>1</u>
Solvent amount, mL	<u>200</u>	<u>X</u>	<u>X</u>	<u>X</u>
Final volume, V _i mL	<u>1000</u>			
Extraction DF f _e = V/X	<u>0.001</u>			
Sub-ID of extracts	<u>1,4-Dioxane</u>			
Anomaly Footnote [†] :	_____			
Sample Cleanup	Op. by _____	Date <u>/ /</u>	(For details, see Cleanup worksheet)	
Cleanup method:	_____			
Cleanup DF, f _c	_____			
Preparation DF, f ₁ = f _e f _c	_____			

Pre-injection*:	Op. by _____	Date <u>/ /</u>	IS Lot : <u>GC-</u>	Conc. _____ pp
2nd dilution, f ₂ (1)	_____	_____	_____	_____
V _i , μL (40)	_____	_____	_____	_____
Single Extract GC, GC/MS Analysis:				
V ₂ = V _i /f ₂ , μL (40/f ₂)	_____	_____	_____	_____
V _{solvent} = V _i - V ₂ , μL	_____	_____	_____	_____
Double Extract GC, GC/MS Analysis:				
V _{>} = 1/2 V _i /f ₂ , μL (20/f ₂)	_____	_____	_____	_____
V _{<} = 1/2 V _i f _{<} /f _{>} , μL (20 f _{<} /f _{>})	_____	_____	_____	_____
V _{solvent} = V _i - V _{>} - V _{<}	_____	_____	_____	_____
IS volume, μL (10)	_____	_____	_____	_____
Final Conc. of IS (40 ppm)	_____	_____	_____	_____
Total F = 2 f ₁ f _{>} or f ₁ f ₂	_____	_____	_____	_____

[†] Extraction Anomaly Footnote:

1. Difficult to concentrate to 1.0 mL.	5. Too much precipitate during extraction.	9. Sample has strong stink
2. Extract is dark color	6. There is an oil layer above the sample.	10. Some solvent suspected
3. Extract is sticky	7. Mixture of soil & water.	11. special color, specify
4. Heterogenous matrix.	8. BN-extract is interfered, A-extract only.	

* For double extract analysis (ABN), f_> = max[f_A, f_B] and f_< = min[f_A, f_B], where f_A, f_B are f₁ for A and BN extracts, respectively. The values in () is the typical case for 625/8270 ABN analysis.

M

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Semi-VOC Analysis Logbook

Sequence # 13672623

Starting Date: 5/23/03 Time 12:09 pm Analyst: JW

Seq. type: Cal. Ini. Batch Middle Final Continue Study

Datafile Path: C:\APCH\2003\1\DATA\0367

Routine Maintenance: Replace Septum Replace Liner Replace Seal Cut Guid Column Cut Column Others

Op. #	Batch-No	MTX	S Type*	Sample ID	Method**	f ₂	F	A-#	Datafile	OK ?†	Note††
1121			SP	672623P01	DFTPP6VS			99	672623P01	✓	5/23/03 12:09 pm
1122			CCV	Q01	DIOSIMob			100	Q01	✓	GC14349
1123	03672623	S	MB	K01			0.001	1	K01	✓	
1124			LCS	L01				2	L01	✓	
1125			LCSB	J01				3	J01	✓	
1126			MS	M01				4	M01	✓	#3318-01
1127			MSD	N01				5	N01	✓	1
1128			MS	M02				6	M02	✓	#3318-04
1129			MSB	N02				7	N02	✓	1
1130				3318-01				8	3318-01	✓	
1131				02				9	02	✓	
1132				03				10	03	✓	
1133				04				11	04	✓	
1134				05				12	05	✓	
1135				06				13	06	✓	
1136				07				14	07	✓	
1137				3304-01				15	3304-01	✓	
1138				3301-02				16	3301-02	✓	
1139				3334-01				17	3334-01	✓	
1140				02				18	02	✓	
1141				03				19	03	✓	
1142				04				20	04	✓	
1143				05				1	05	✓	
1144				3351-01				2	3351-01	✓	
1145				02				3	02	✓	
1146				03				4	03	✓	
1147				04				5	04	✓	
1148				05				6	05	✓	
1149				06				7	06	✓	
1150											

JW 5-27-03

Footnote/Anomaly:

* Sample type such as, SP(BFB, or DFTPP), CCV, CVA, CVB, LCS, LCSB, MS, MSD, MD, CLS etc. For field samples leave as blank.
 ** Method name: [MMMM][S][nnn]. [MMMM] - method code, per SOP C-88. Special codes: DSL2: Diesel #2; FP30: Fuel Finger Print (C8-C30); FP40: Fuel Finger Print (C8-C40); EGLY: Ethylene Glycol; FORM: Formaldehyde; MEOH: Methyl Alcohol. [S] - APCL system code; [nnn] - initial calibration number.
 † Specify the result of the injection is accepted (✓), or denied (X), or need Re-run (R) or use as a reference data (ref).
 †† Lot # for CCV/Closing, Time for SP (e.g. DFTPP) must be recorded.

Supervisor Initial [Signature]

Semi-VOC Analysis Logbook

Sample # 026502

Starting Date: 12-17-02 Time 2:19 pm Analyst: J.Y.

Sample type: Cal. Ini. Batch Middle Final Continue Study
 Maintenance: Replace Septum Replace Liner Replace Seal Cut Guid Column Cut Column Others

Datafile Path: _____

Batch-No	MTX	S Type*	Sample ID	Method**	f ₂	F	A-#	Datafile	OK ?†	Note††
		SP	G5027P01	DFTPP625			99	G5027P01	✓	12-17-02 2:19 pm
		Cali-	M06-040	DIOSIM06			1	M06-040	✓	Koppin GC14554
			-030				2	-030	✓	30
			-020				3	-020	✓	20
			-010				4	-010	✓	10
			-001				5	-001	✓	1
		CCV	G5027Q01				100	G5027Q01	✓	GC14554
026502	W	MR	K01			0.001	6	K01	-	
		LCS	L01				7	L01	-	
		LCSD	J01				8	J01	-	
		MS	M01				9	M01	-	\$6556-03
		MSD	N01				10	N01	-	1
		MS'	M02			0.000962	11	M02	-	\$6587-01
		MSD'	N02				12	N02	-	1
			6587-01				13	6587-01	-	
			-02				14	-02	-	
			-03				15	-03	-	
			6556-01			0.002	16	6556-01	-	
			-02			0.001	17	-02	-	
			-03				18	-03	-	
			6588-01			0.000962	19	6588-01	-	
			6603-01				20	6603-01	-	
			-02				21	-02	-	
			6558-01				22	6558-01	-	
			-03			0.001	23	-03	-	
			6586-02				24	6586-02	-	
			-04				25	-04	-	
			-05				26	-05	-	
			-06				27	-06	-	Reverse side
			-08				28	-08	↓	✓

Footnote/Anomaly:

* Sample type such as, SP(BFB, or DFTPP), CCV, CVA, CVB, LCS, LCSD, MS, MSD, MD, CLS etc. For field samples leave as blank.
 ** Method name: [MMMM][S][nnn]. [MMMM] - method code, per SOP C-88. Special codes: DSL2: Diesel #2; FP30: Fuel Finger Print (C8-C30); FP40: Fuel Finger Print (C8-C40); EGLY: Ethylene Glycol; FORM: Formaldehyde; MEOH: Methyl Alcohol. [S] - APCL system code; [nnn] - initial calibration number.
 † Specify the result of the injection is accepted (✓), or denied (X), or need Re-run (R) or use as a reference data (ref).
 †† Lot # for CCV/Closing, Time for SP (e.g. DFTPP) must be recorded.

Supervisor Initial J.Y.