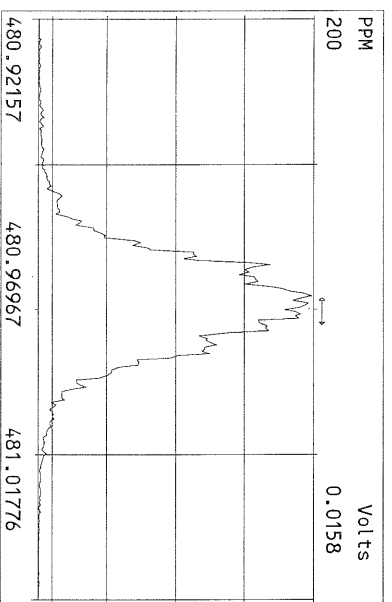
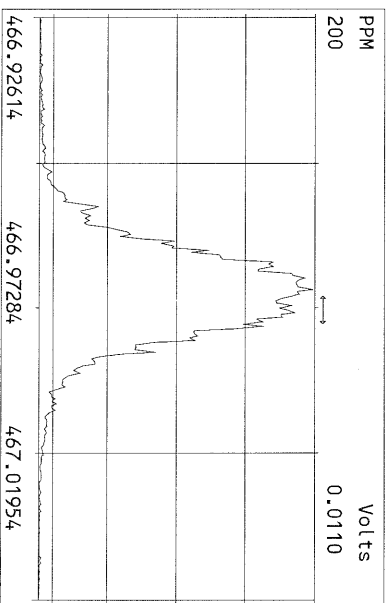
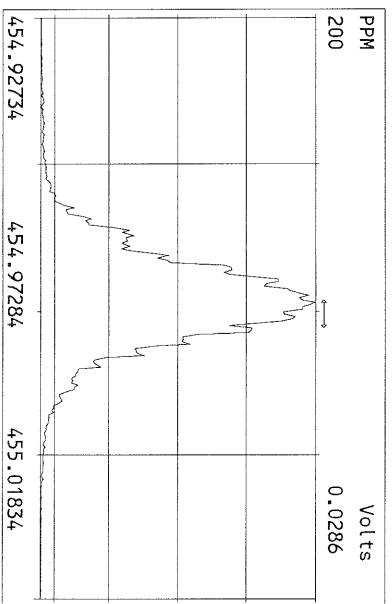
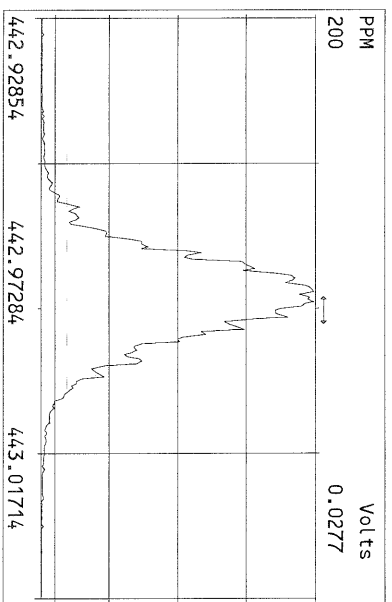
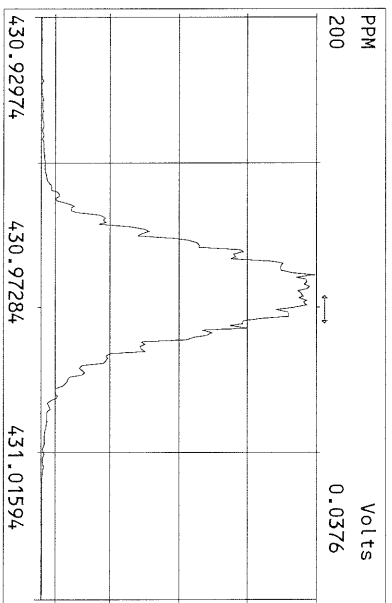
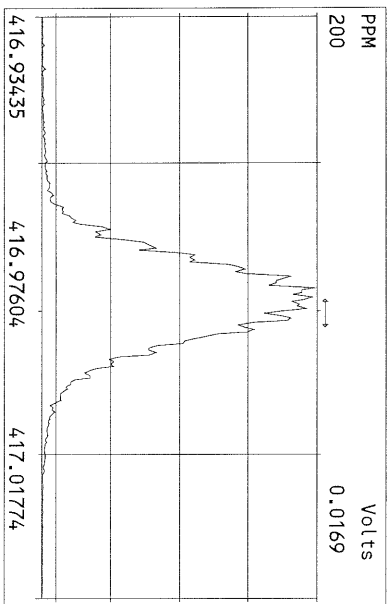
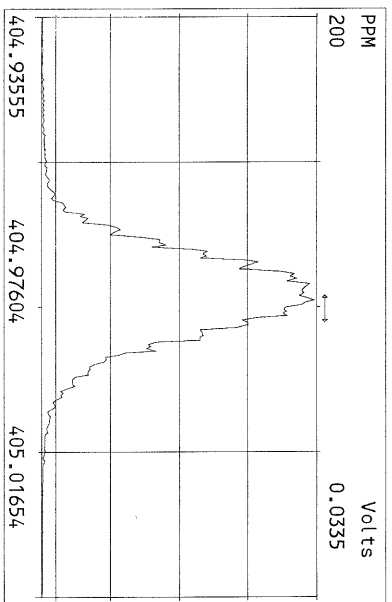
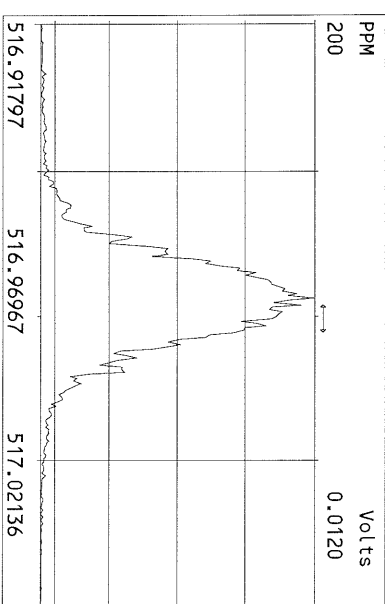
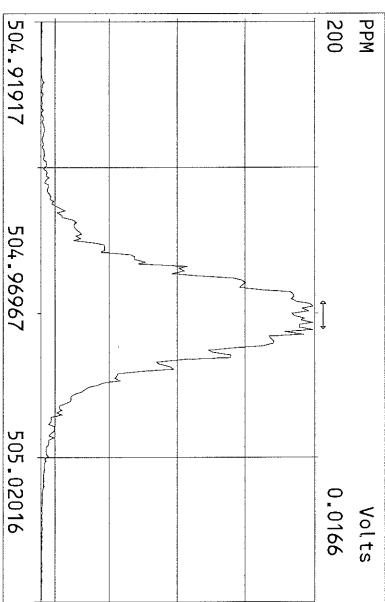
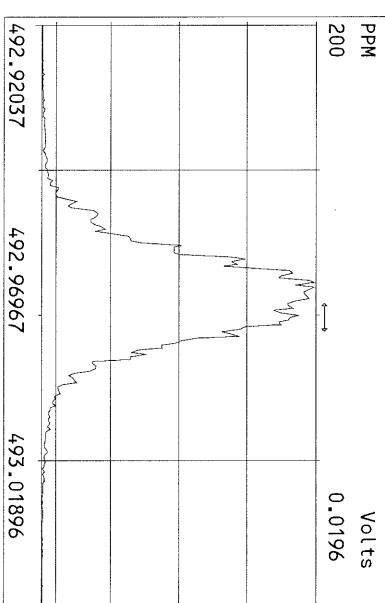
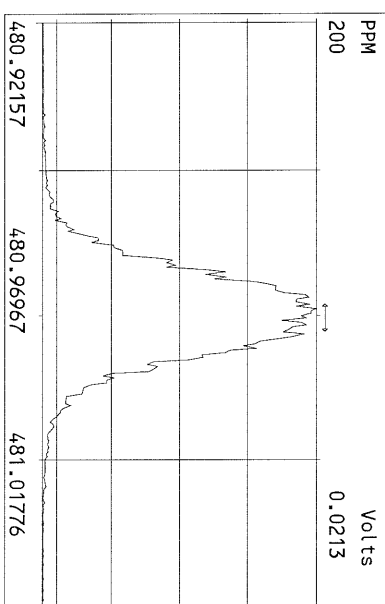
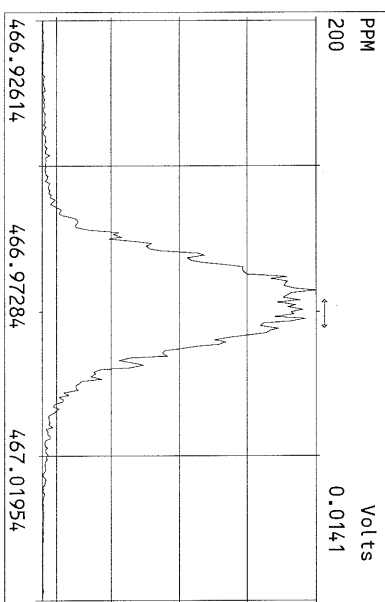
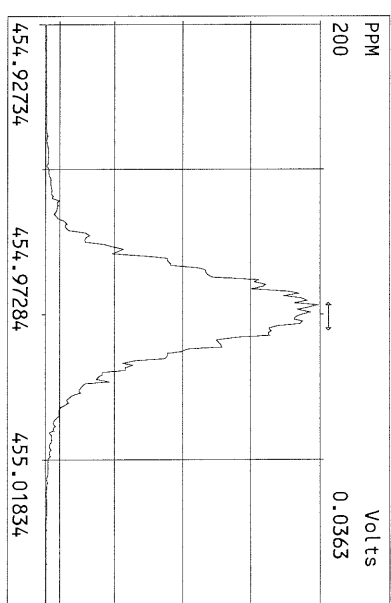
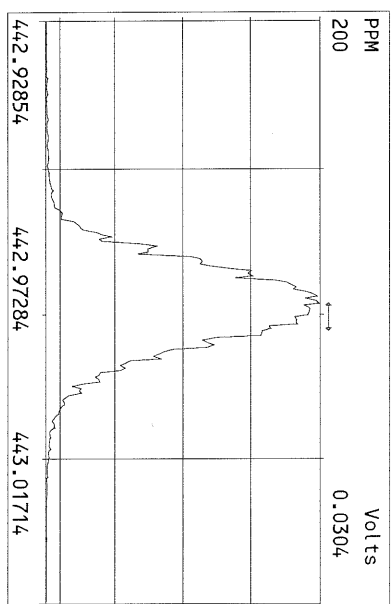
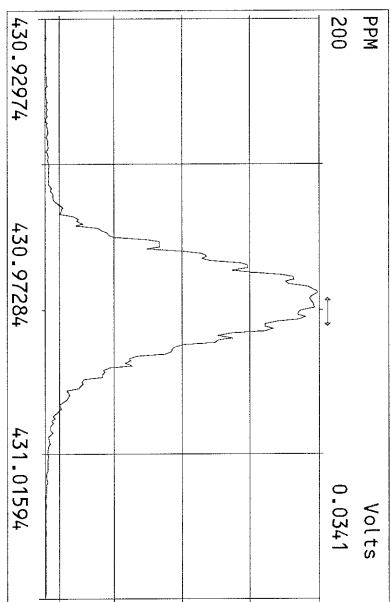


Peak Locate Examination: 9-APR-2011:08:32 File:08APR11M_RES_CHECK
Experiment:OCD Function:4 Reference:PFK





USEPA - ITD

FORM 4A

PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3 GC Column ID: DB5

VER Data Filename: 08APR11M Sam:16 Analysis Date: 9-APR-11 04:06:14

NATIVE ANALYTES	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
2,3,7,8-TCDD	M/M+2	0.79	0.65-0.89	y	10.8	7.80 - 12.9 ✓
1,2,3,7,8-PeCDD	M+2/M+4	1.56	1.32-1.78	y	52.3	39.0 - 65.0 ✓
1,2,3,4,7,8-HxCDD	M+2/M+4	1.29	1.05-1.43	y	48.7	39.0 - 64.0 ✓
1,2,3,6,7,8-HxCDD	M+2/M+4	1.29	1.05-1.43	y	51.0	39.0 - 64.0 ✓
1,2,3,7,8,9-HxCDD	M+2/M+4	1.30	1.05-1.43	y	52.6	41.0 - 61.0 ✓
1,2,3,4,6,7,8-HpCDD	M+2/M+4	0.92	0.88-1.20	y	49.4	43.0 - 58.0 ✓
OCDD	M+2/M+4	0.87	0.76-1.02	y	100	79.0 - 126 ✓
2,3,7,8-TCDF	M/M+2	0.67	0.65-0.89	y	11.4	8.40 - 12.0 ✓
1,2,3,7,8-PeCDF	M+2/M+4	1.60	1.32-1.78	y	52.8	41.0 - 60.0 ✓
2,3,4,7,8-PeCDF	M+2/M+4	1.61	1.32-1.78	y	51.4	41.0 - 60.0 ✓
1,2,3,4,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	49.8	45.0 - 56.0 ✓
1,2,3,6,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	50.5	44.0 - 57.0 ✓
2,3,4,6,7,8-HxCDF	M+2/M+4	1.28	1.05-1.43	y	49.4	44.0 - 57.0 ✓
1,2,3,7,8,9-HxCDF	M+2/M+4	1.24	1.05-1.43	y	50.7	45.0 - 56.0 ✓
1,2,3,4,6,7,8-HpCDF	M+2/M+4	1.06	0.88-1.20	y	49.2	45.0 - 55.0 ✓
1,2,3,4,7,8,9-HpCDF	M+2/M+4	1.05	0.88-1.20	y	51.2	43.0 - 58.0 ✓
OCDF	M+2/M+4	0.88	0.76-1.02	y	96.2	63.0 - 159 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

Analyst: Date: 4/11/11

USEPA - ITD

FORM 4B
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 08APR11M Sam:16

Analysis Date: 9-APR-11 04:06:14

LABELLED COMPOUNDS	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
13C-2,3,7,8-TCDD	M/M+2	0.81	0.65-0.89	y	93.1	82.0 - 121 ✓
13C-1,2,3,7,8-PeCDD	M+2/M+4	1.77	1.32-1.78	y	89.8	62.0 - 160 ✓
13C-1,2,3,4,7,8-HxCDD	M+2/M+4	1.26	1.05-1.43	y	87.8	85.0 - 117 ✓
13C-1,2,3,6,7,8-HxCDD	M+2/M+4	1.26	1.05-1.43	y	102	85.0 - 118 ✓
13C-1,2,3,4,6,7,8-HpCDD	M+2/M+4	1.05	0.88-1.20	y	116	72.0 - 138 ✓
13C-OCDD	M+2/M+4	0.90	0.76-1.02	y	203	96.0 - 415 ✓
13C-2,3,7,8-TCDF	M/M+2	0.85	0.65-0.89	y	101	71.0 - 140 ✓
13C-1,2,3,7,8-PeCDF	M+2/M+4	1.67	1.32-1.78	y	107	76.0 - 130 ✓
13C-2,3,4,7,8-PeCDF	M+2/M+4	1.71	1.32-1.78	y	109	77.0 - 130 ✓
13C-1,2,3,4,7,8-HxCDF	M/M+2	0.49	0.43-0.59	y	91.5	76.0 - 131 ✓
13C-1,2,3,6,7,8-HxCDF	M/M+2	0.47	0.43-0.59	y	99.5	70.0 - 143 ✓
13C-2,3,4,6,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	98.7	73.0 - 137 ✓
13C-1,2,3,7,8,9-HxCDF	M/M+2	0.47	0.43-0.59	y	99.1	74.0 - 135 ✓
13C-1,2,3,4,6,7,8-HpCDF	M/M+2	0.46	0.37-0.51	y	112	78.0 - 129 ✓
13C-1,2,3,4,7,8,9-HpCDF	M/M+2	0.48	0.37-0.51	y	114	77.0 - 129 ✓
13C-OCDF	M+2/M+4	0.93	0.76-1.02	y	209	96.0 - 415 ✓
CLEANUP STANDARD (4)						
37Cl-2,3,7,8-TCDD					8.36	7.80 - 12.8 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

(4) No ion abundance ratio; report concentration found.

Analyst: _____

Date: _____

FORM 5
PCDD/PCDF RT WINDOW AND ISOMER SPECIFICITY STANDARDS

Lab Name: Frontier Analytical Laboratory Episode No.:
 Contract No.: SAS No.:
 Instrument ID: FAL3 Initial Calibration Date: 3/7/11
 RT Window Data Filename: 08APR11M Sam:16 Analysis Date: 9-APR-11 Time: 04:06:14
 DB-5 IS Data Filename: 08APR11M Sam:16 Analysis Date: 9-APR-11 Time: 04:06:14
 DB-225 IS Data Filename: Analysis Date: Time:

DB-5 RT WINDOW DEFINING STANDARDS RESULTS

ISOMERS	ABSOLUTE RT	ISOMERS	ABSOLUTE RT
1,3,6,8-TCDD (F)	24:35 ✓	1,3,6,8-TCDF (F)	23:13 ✓
1,2,8,9-TCDD (L)	28:33 ✓	1,2,8,9-TCDF (L)	28:46 ✓
1,2,4,7,9-PeCDD (F)	30:28 ✓	1,3,4,6,8-PeCDF (F)	28:38 ✓
1,2,3,8,9-PeCDD (L)	34:02 ✓	1,2,3,8,9-PeCDF (L)	34:29 ✓
1,2,4,6,7,9-HxCDD (F)	36:23 ✓	1,2,3,4,6,8-HxCDF (F)	35:30 ✓
1,2,3,7,8,9-HxCDD (L)	39:27 ✓	1,2,3,7,8,9-HxCDF (L)	40:02 ✓
1,2,3,4,6,7,9-HpCDD (F)	43:04 ✓	1,2,3,4,6,7,8-HpCDF (F)	42:33 ✓
1,2,3,4,6,7,8-HpCDD (L)	44:28 ✓	1,2,3,4,7,8,9-HpCDF (L)	45:23 ✓

(F) = First eluting isomer (DB-5); (L) = Last eluting isomer (DB-5)

ISOMER SPECIFICITY (IS) TEST STANDARD RESULTS

% VALLEY HEIGHT
BETWEEN
COMPARED PEAKS (1)

<25%

(1) To meet contract requirement, %Valley Height Between Compared Peaks shall not exceed 25% (section 15.4.2.2, Method 1613).

Analyst: _____

Date: 4/11/11

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 9-APR-11 04:06:14

CS3 or VER Data Filename: 08APR11M

Sam:16

NATIVE ANALYTES	RETENTION TIME REFERENCE	RRT	RRT QC LIMITS (1)
2,3,7,8-TCDD	13C-2,3,7,8-TCDD	1.001	0.999-1.002 ✓
2,3,7,8-TCDF	13C-2,3,7,8-TCDF	1.001	0.999-1.003 ✓
1,2,3,7,8-PeCDD	13C-1,2,3,7,8-PeCDD	1.000	0.999-1.002 ✓
1,2,3,7,8-PeCDF	13C-1,2,3,7,8-PeCDF	1.001	0.999-1.002 ✓
2,3,4,7,8-PeCDF	13C-2,3,4,7,8-PeCDF	1.000	0.999-1.002 ✓
LABELED COMPOUNDS			
37Cl-2,3,7,8-TCDD	13C-1,2,3,4-TCDD	1.023	0.989-1.052 ✓
13C-2,3,7,8-TCDD		1.022	0.976-1.043 ✓
13C-2,3,7,8-TCDF		0.994	0.923-1.103 ✓
13C-1,2,3,7,8-PeCDD		1.239	1.000-1.567 ✓
13C-1,2,3,7,8-PeCDF		1.175	0.923-1.203 ✓
13C-2,3,4,7,8-PeCDF		1.224	0.923-1.303 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified in Table 2, Method 1613.

Analyst: Date:

USEPA - ITD

FORM 6B

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 9-APR-11 04:06:14

CS3 or VER Data Filename: 08APR11M

Sam:16

NATIVE ANALYTES	RETENTION TIME REFERENCE	RRT	RRT QC LIMITS (1)
1,2,3,4,7,8-HxCDD	13C-1,2,3,4,7,8-HxCDD	1.000	0.999-1.001 ✓
1,2,3,6,7,8-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.001	0.998-1.004 ✓
1,2,3,7,8,9-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.012	1.000-1.019 ✓
1,2,3,4,7,8-HxCDF	13C-1,2,3,4,7,8-HxCDF	1.001	0.999-1.001 ✓
1,2,3,6,7,8-HxCDF	13C-1,2,3,6,7,8-HxCDF	1.000	0.997-1.005 ✓
2,3,4,6,7,8-HxCDF	13C-2,3,4,6,7,8-HxCDF	1.001	0.999-1.001 ✓
1,2,3,7,8,9-HxCDF	13C-1,2,3,7,8,9-HxCDF	1.000	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDD	13C-1,2,3,4,6,7,8-HpCDD	1.001	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDF	13C-1,2,3,4,6,7,8-HpCDF	1.000	0.999-1.001 ✓
1,2,3,4,7,8,9-HpCDF	13C-1,2,3,4,7,8,9-HpCDF	1.001	0.999-1.001 ✓
OCDD	13C-OCDD	1.001	0.999-1.001 ✓
OCDF	13C-OCDF	1.000	0.999-1.001 ✓
LABELED COMPOUNDS			
13C-1,2,3,4,7,8-HxCDD	13C-1,2,3,7,8,9-HxCDD	0.984	0.977-1.000 ✓
13C-1,2,3,6,7,8-HxCDD		0.989	0.981-1.003 ✓
13C-1,2,3,4,7,8-HxCDF		0.949	0.944-0.970 ✓
13C-1,2,3,6,7,8-HxCDF		0.954	0.949-0.975 ✓
13C-2,3,4,6,7,8-HxCDF		0.978	0.959-1.021 ✓
13C-1,2,3,7,8,9-HxCDF		1.015	0.977-1.047 ✓
13C-1,2,3,4,6,7,8-HpCDD		1.127	1.086-1.130 ✓
13C-1,2,3,4,6,7,8-HpCDF		1.079	1.043-1.085 ✓
13C-1,2,3,4,7,8,9-HpCDF		1.150	1.057-1.154 ✓
13C-OCDD		1.270	1.032-1.311 ✓
13C-OCDF		1.279	1.000-1.311 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified in Table 2, Method 1613.

Analyst: _____

Date: _____

Frontier Analytical Laboratory - Acquisition Log

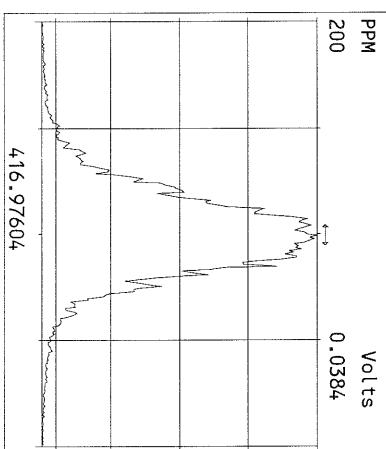
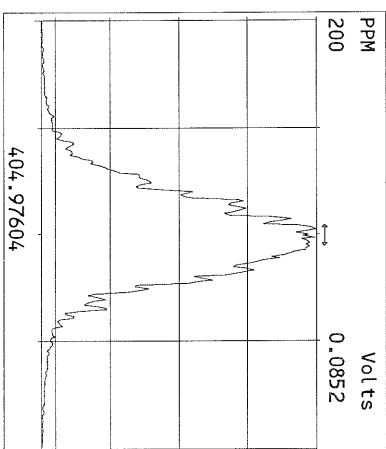
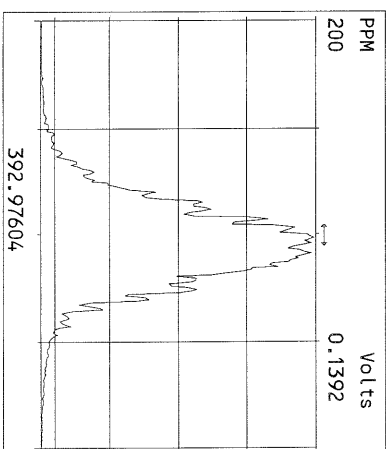
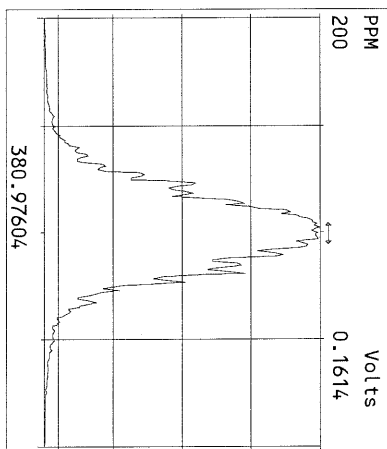
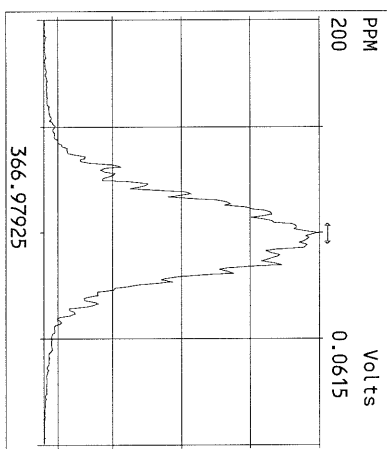
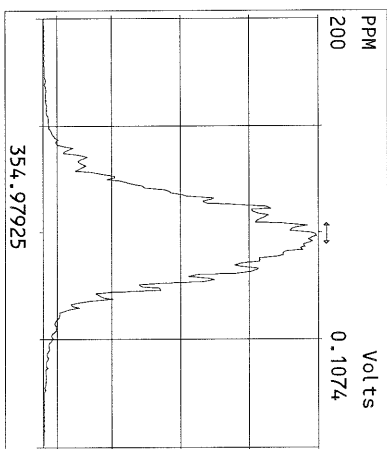
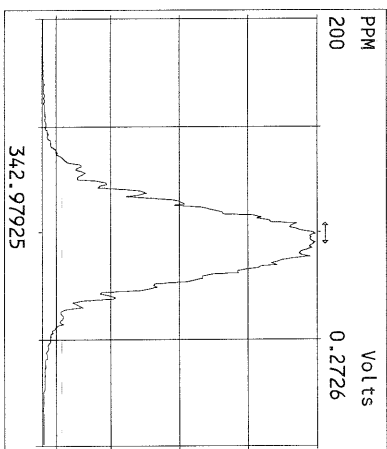
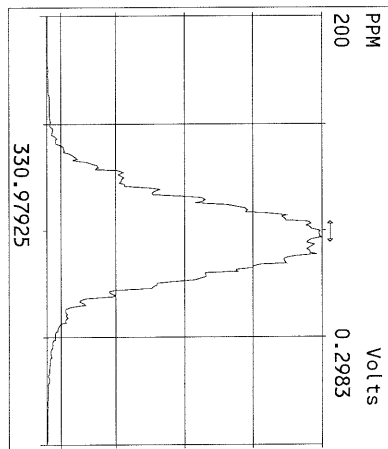
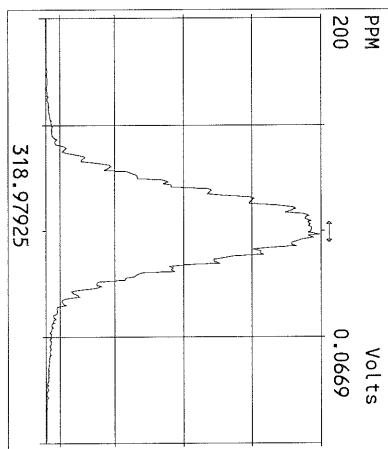
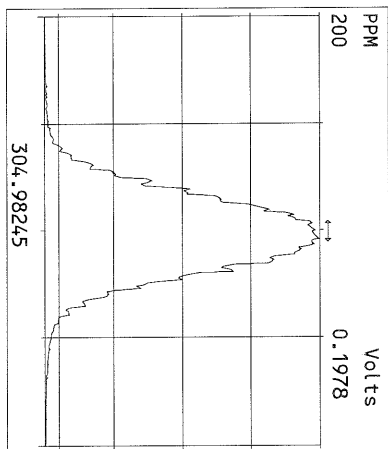
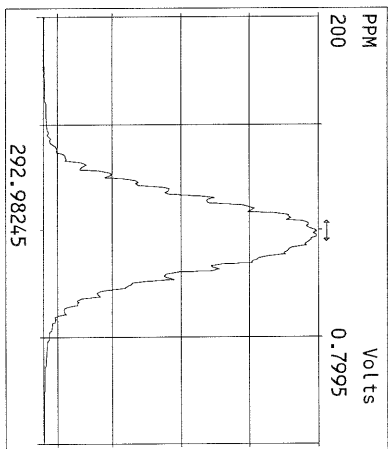
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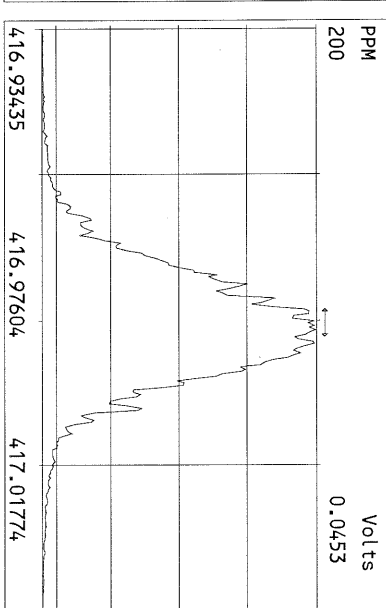
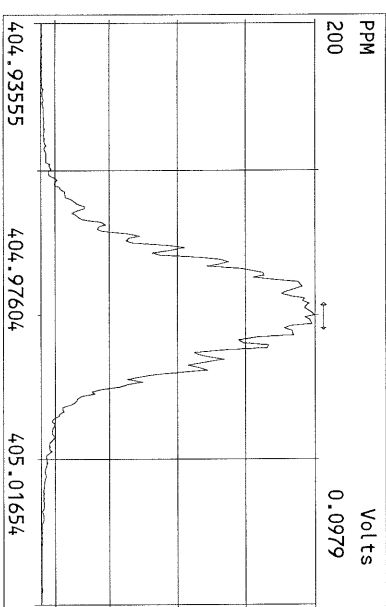
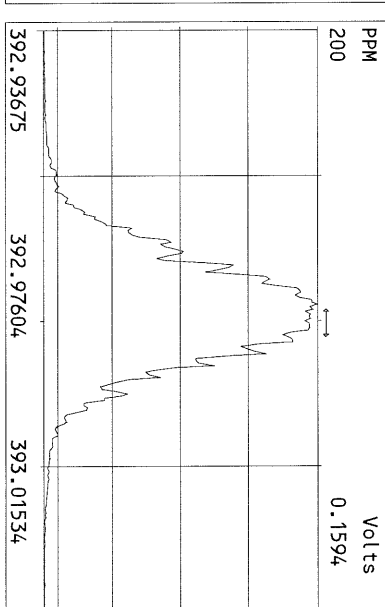
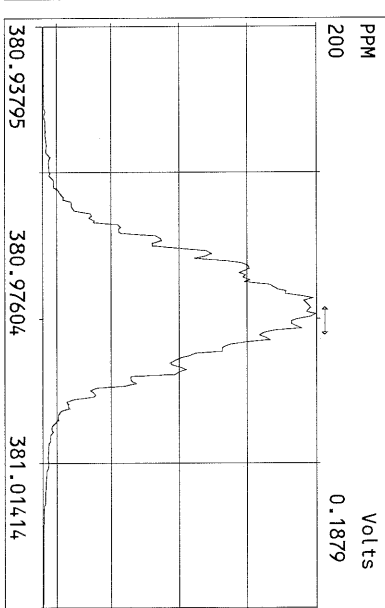
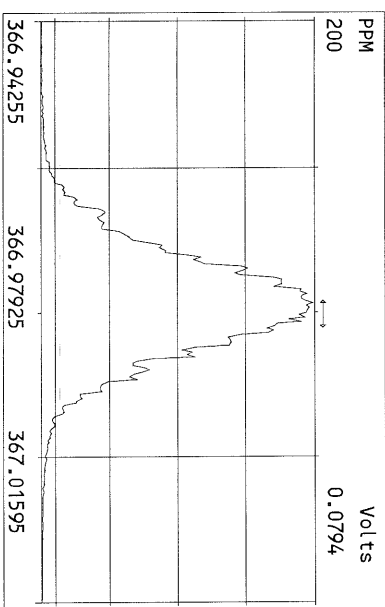
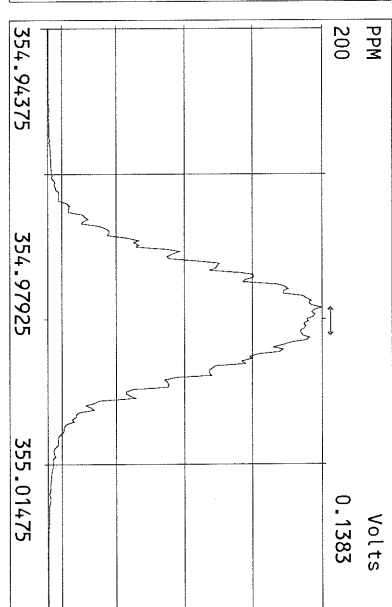
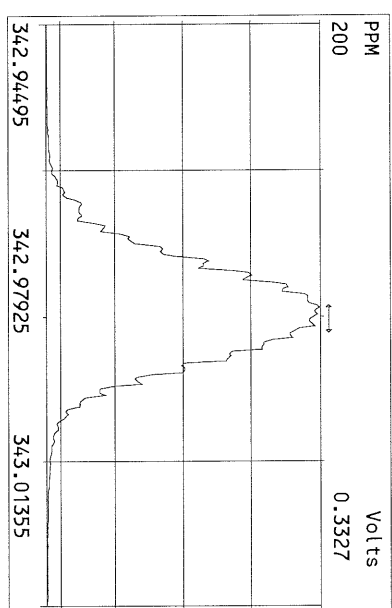
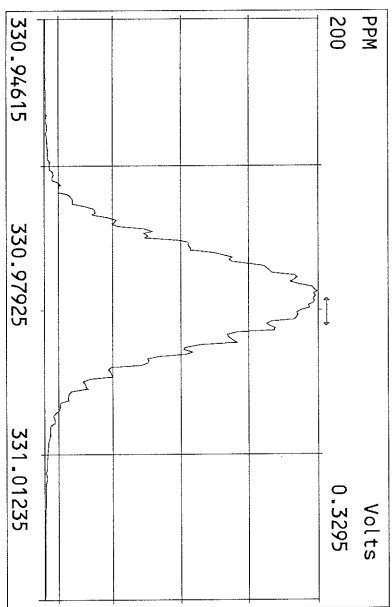
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08APR11M 12	6697-002-0001-SA	1031033-02	9-APR-11 00:24:44	ST040811M1	ST040811M2	TC
08APR11M 13	6701-010-0001-SA	LL-SED1-0-15-032911-ER	9-APR-11 01:20:06	ST040811M1	ST040811M2	TC
08APR11M 14	6684-002-0001-SA	OM-7	9-APR-11 02:15:29	ST040811M1	ST040811M2	TC
08APR11M 15	SB040811M1	Solvent Blank	9-APR-11 03:10:48	ST040811M1	ST040811M2	TC
08APR11M 16	ST040811M2	1613 CS3 100511J	9-APR-11 04:06:14	ST040811M1	ST040811M2	TC

84/11/11

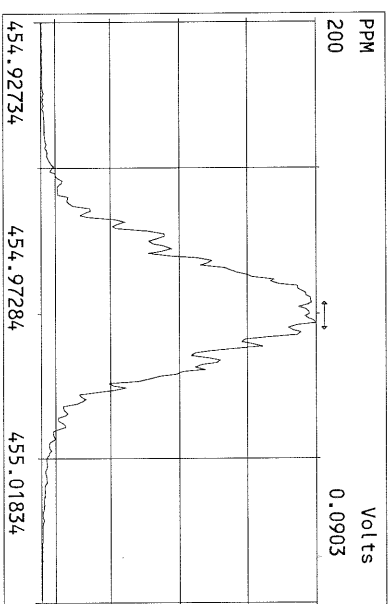
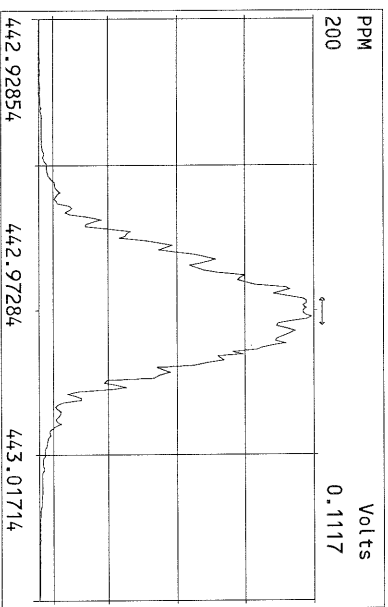
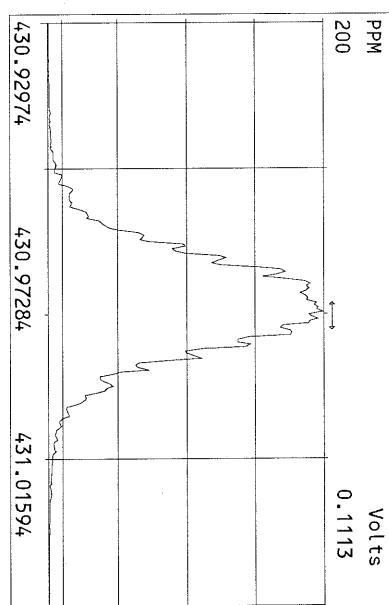
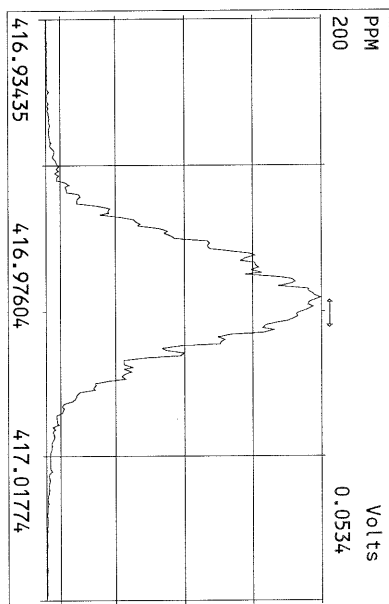
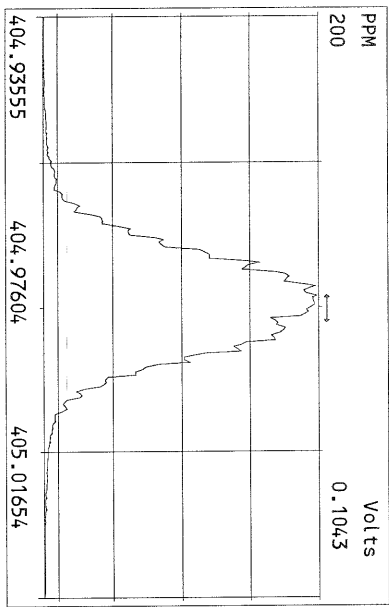
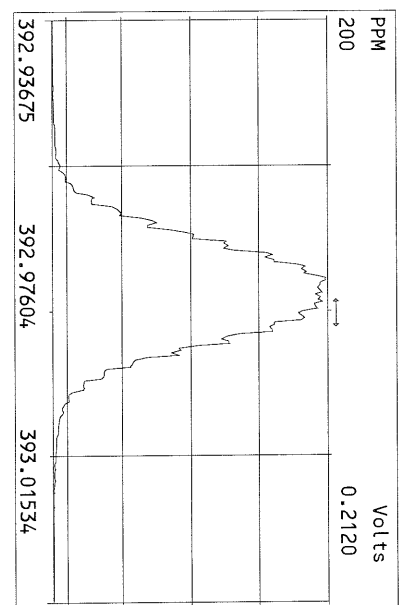
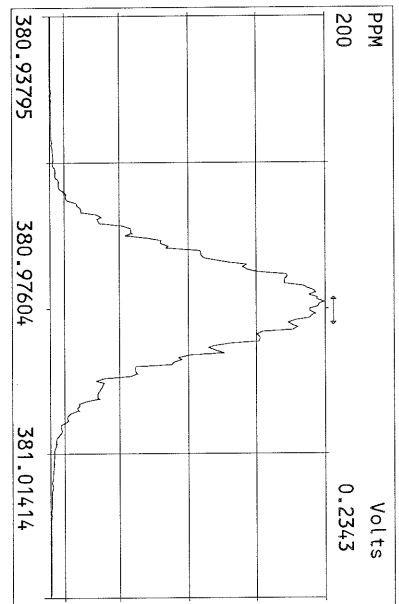
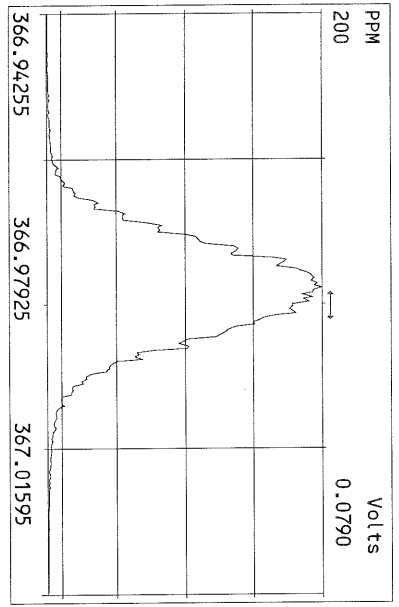
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Date: _____

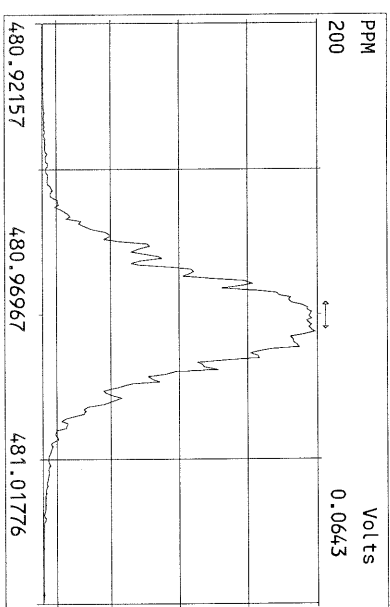
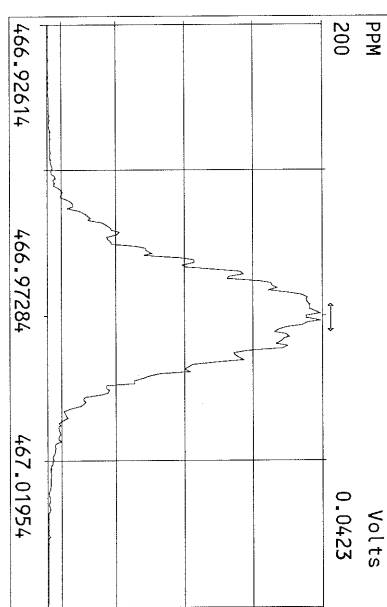
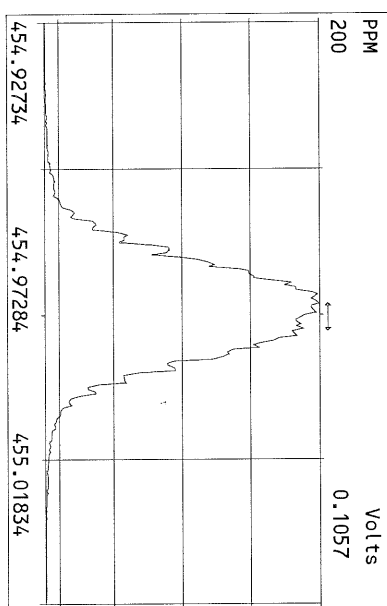
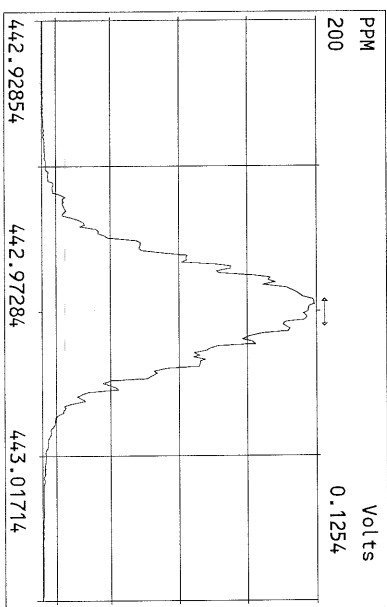
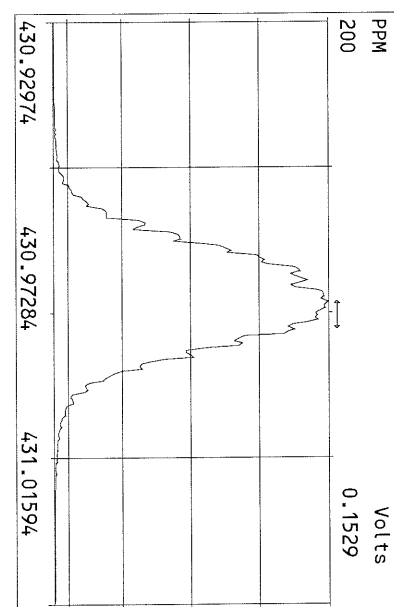
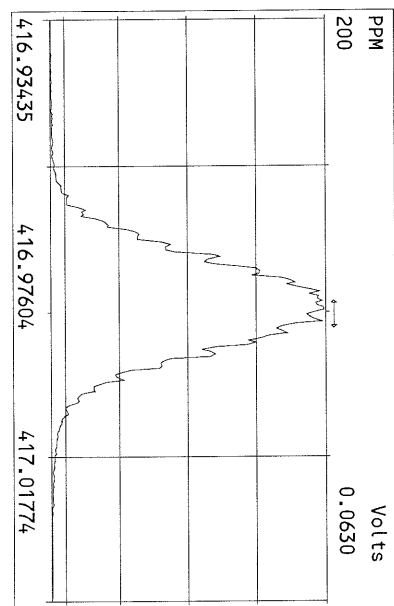
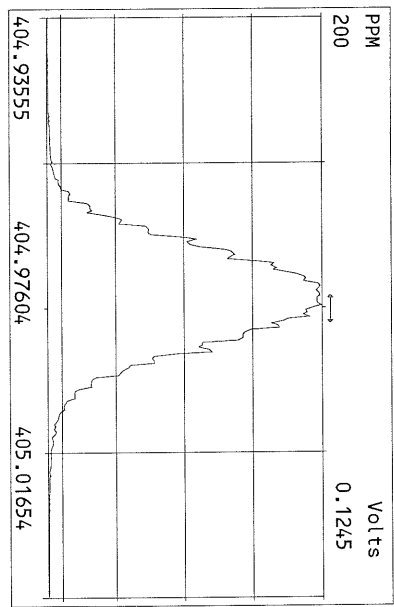


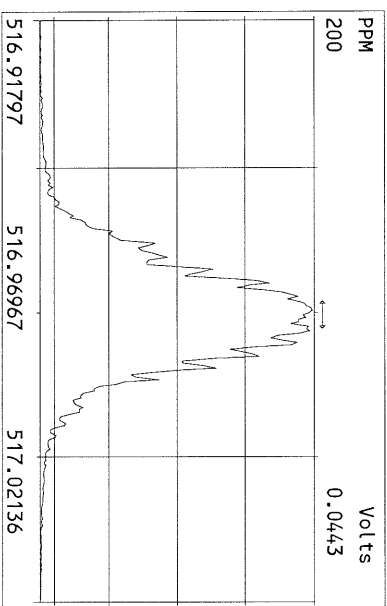
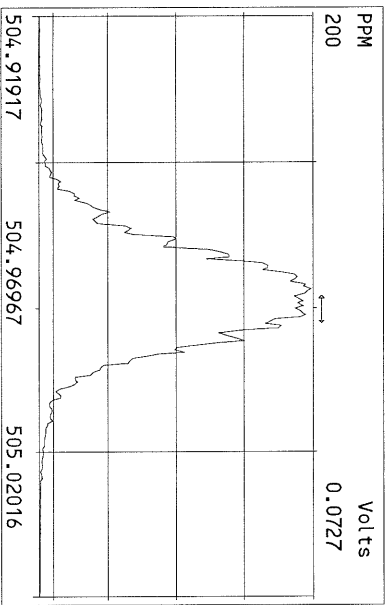
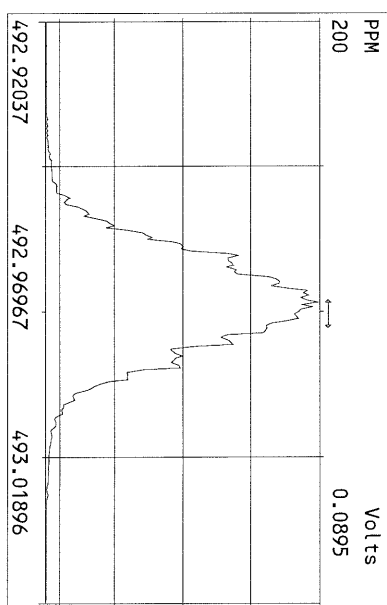
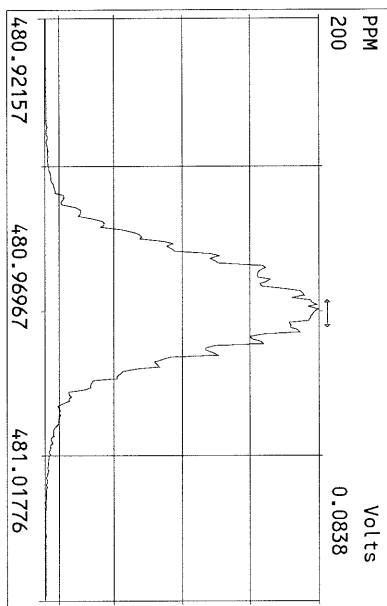
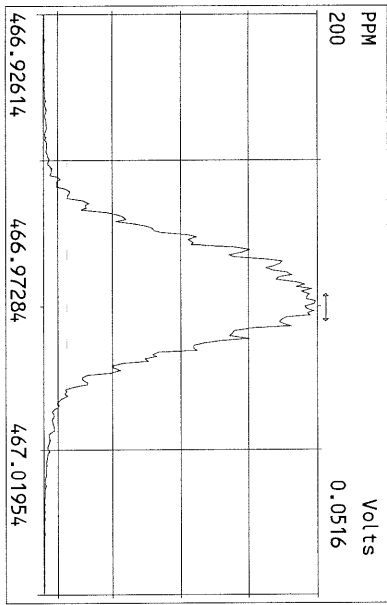
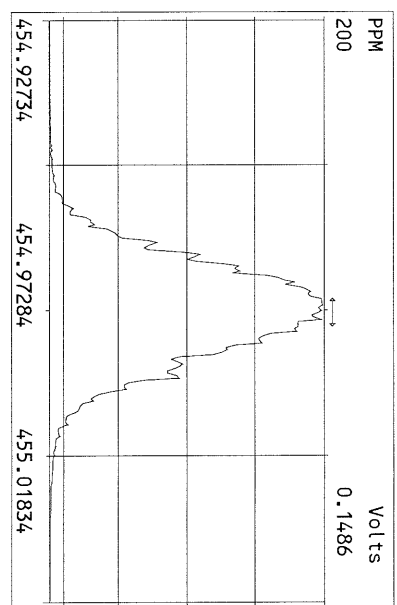
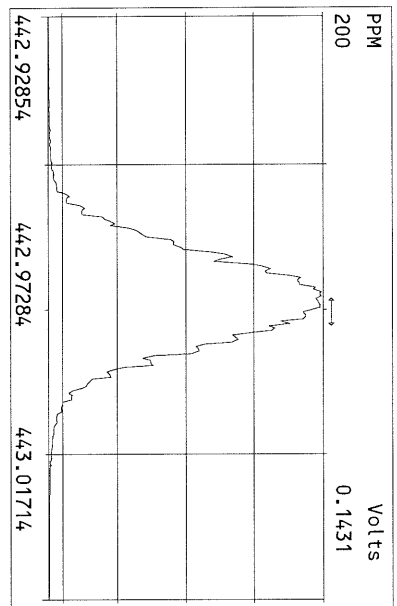
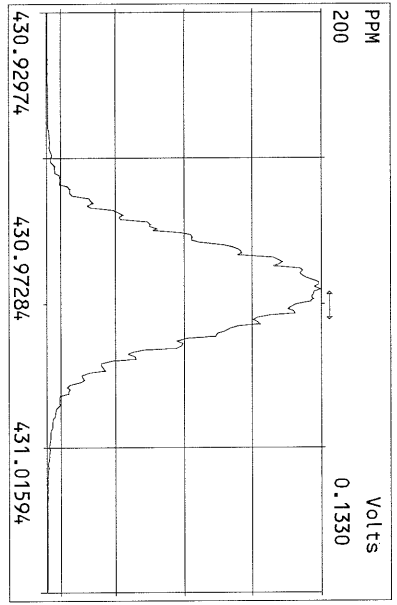


Peak Locate Examination: 8-APR-2011:14:14 File:08APR11M
Experiment:00DD Function:3 Reference:PFK

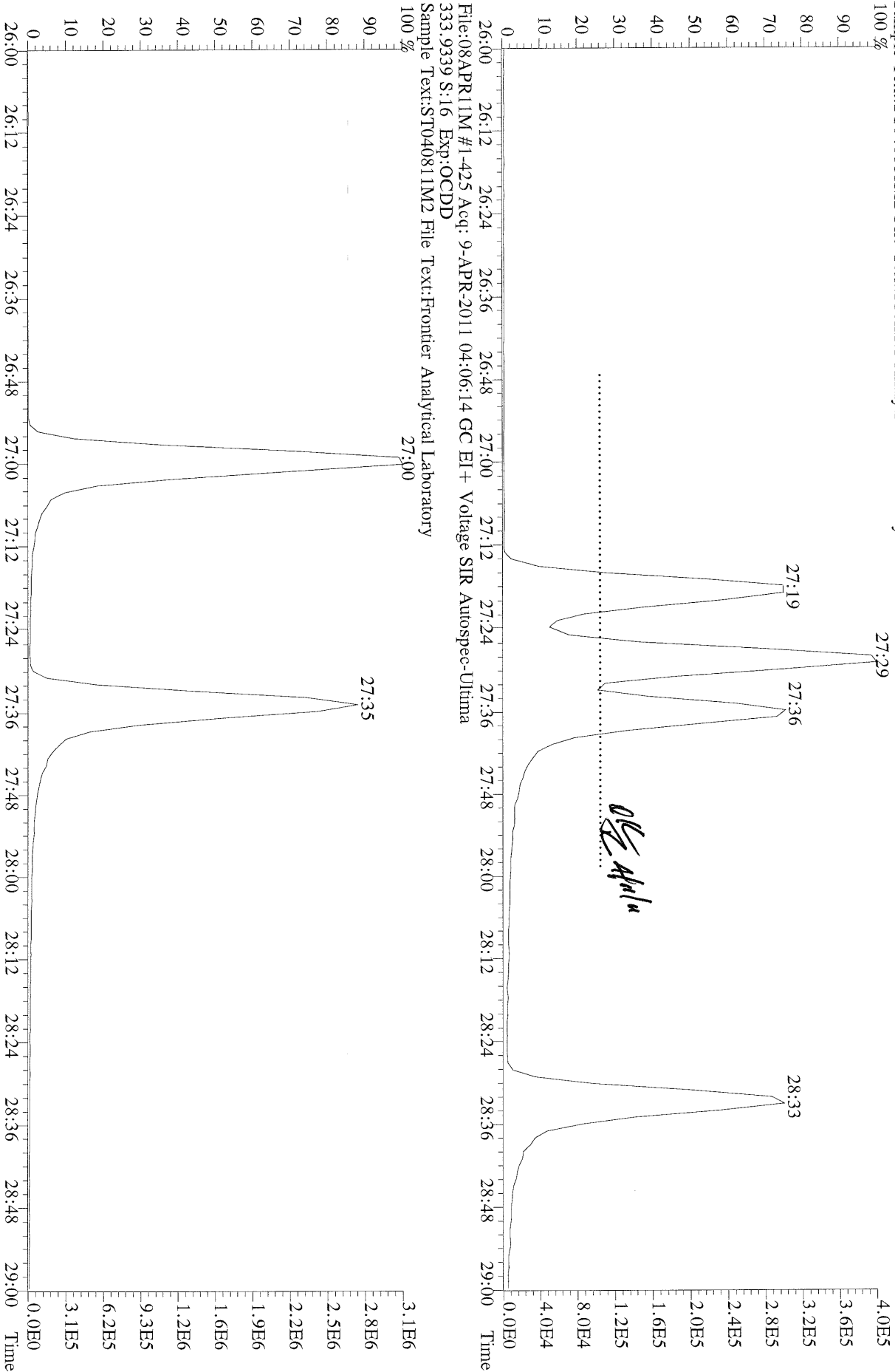


Peak Locate Examination: 8-APR-2011:14:14 File:08APR11M
Experiment:0CDD Function:4 Reference:PFK

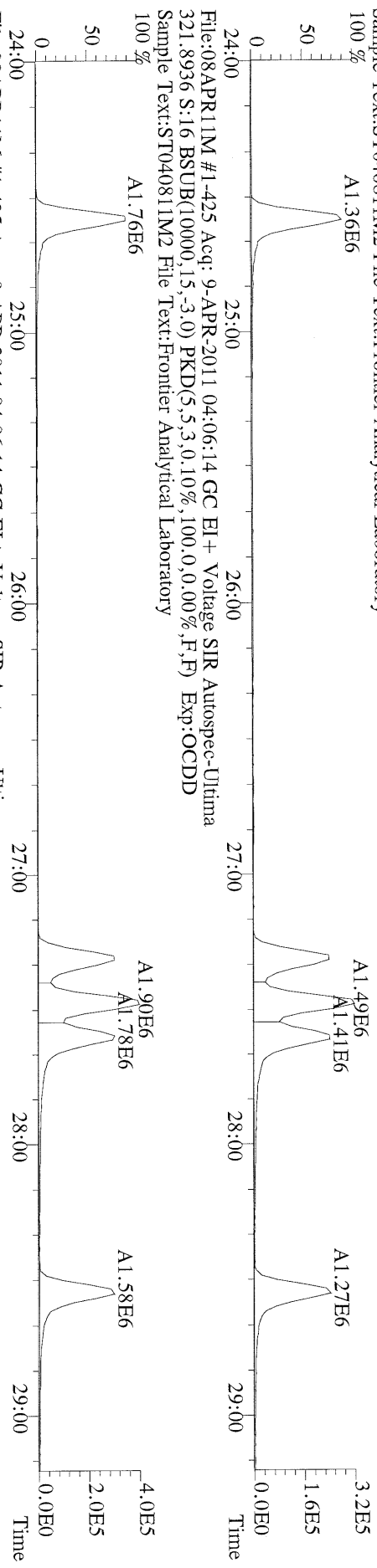




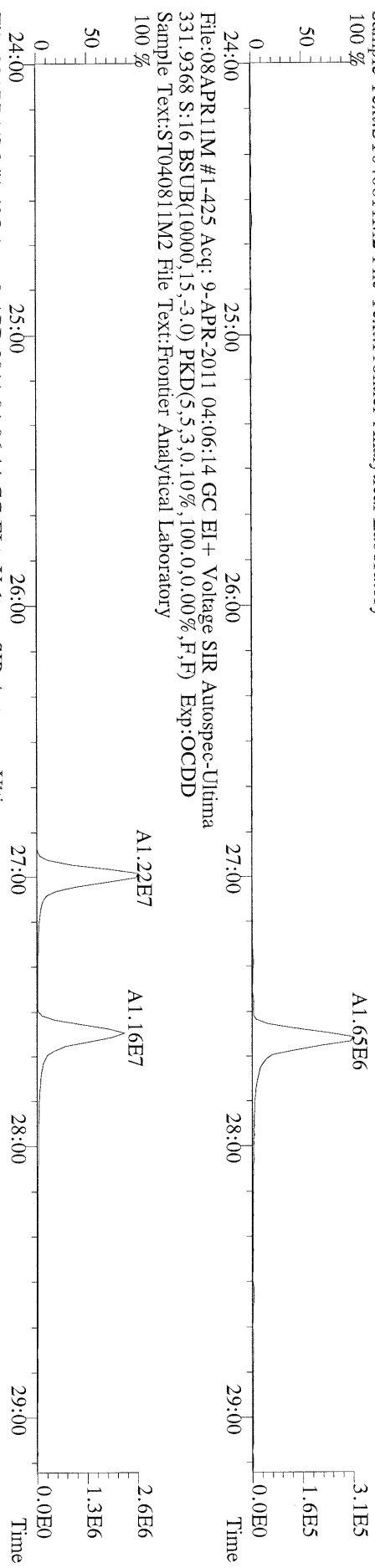
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321.8936 S:16 Exp:OCDD
Sample Text:ST04081M2 File Text:Frontier Analytical Laboratory
100%



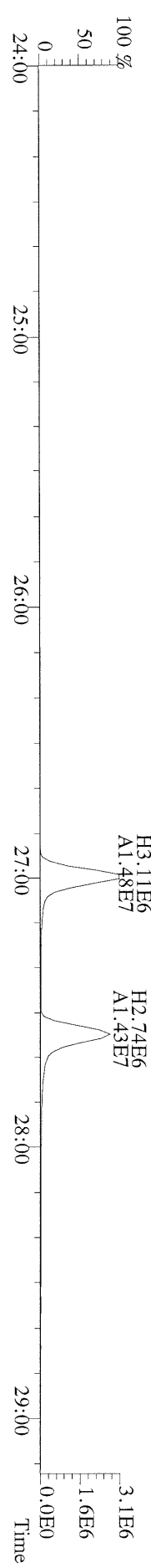
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319.8965 S:16 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



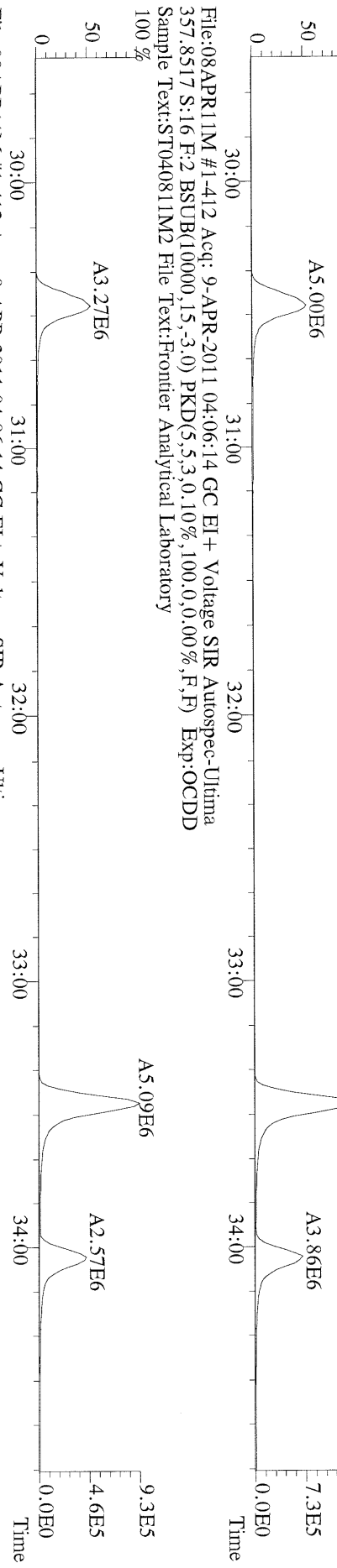
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327.8847 S:16 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



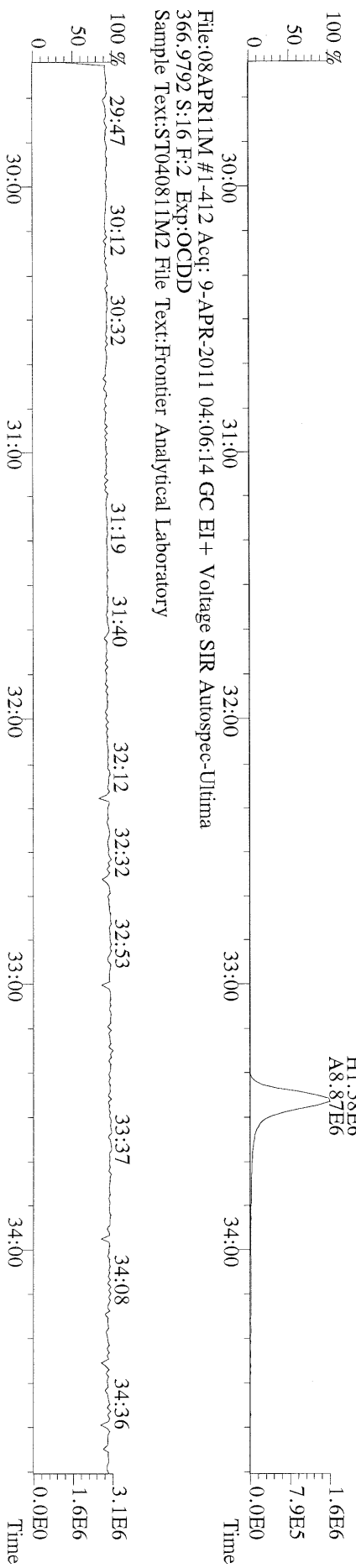
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333.9339 S:16 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
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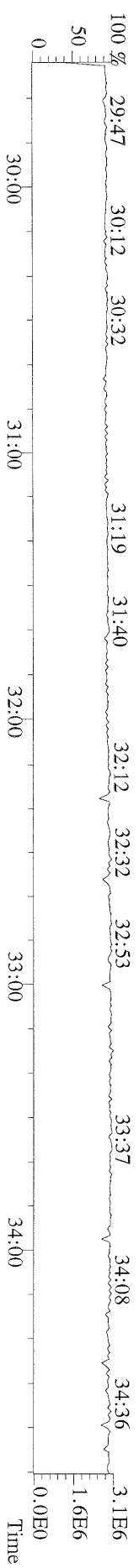
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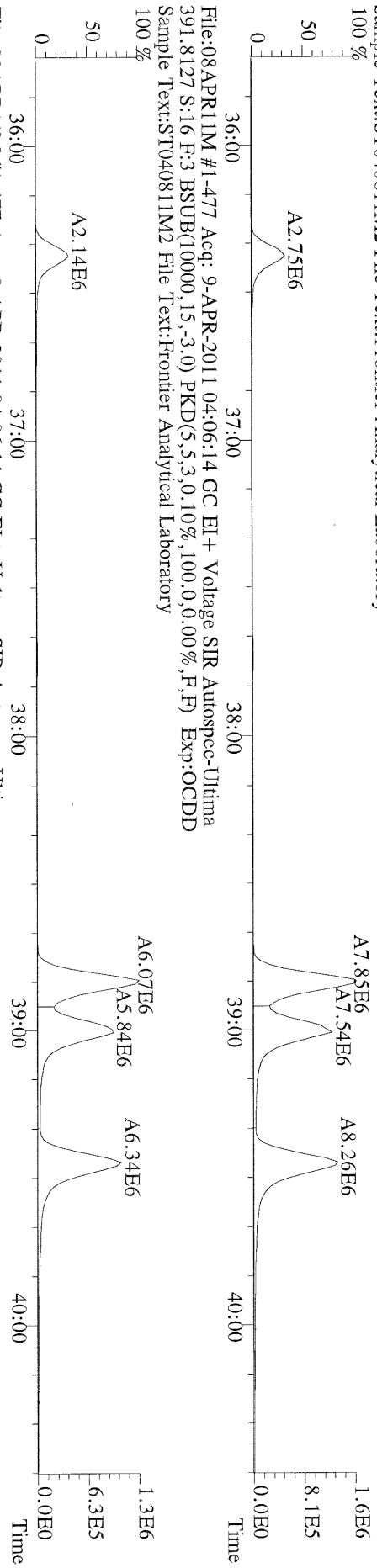
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 367.8949 S:16 F:2 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



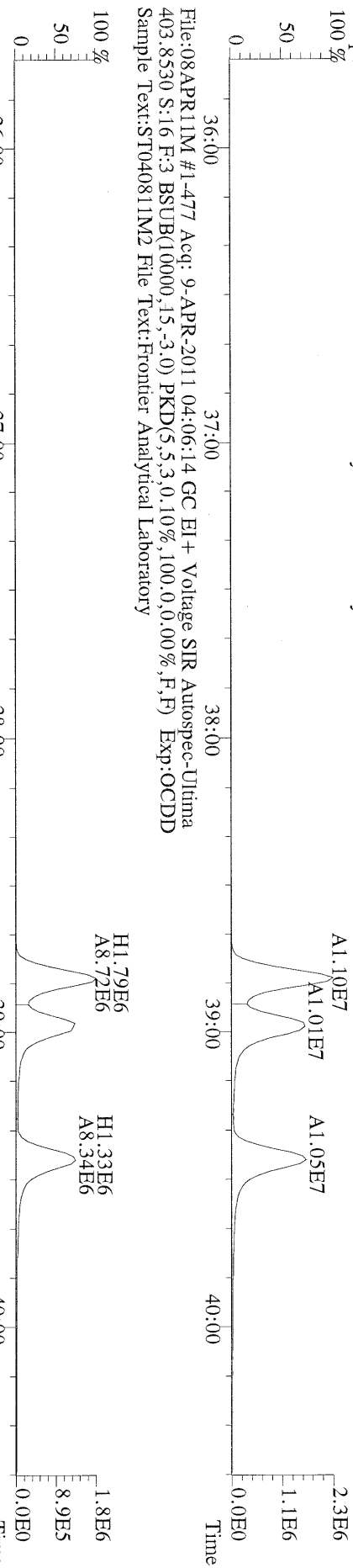
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 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



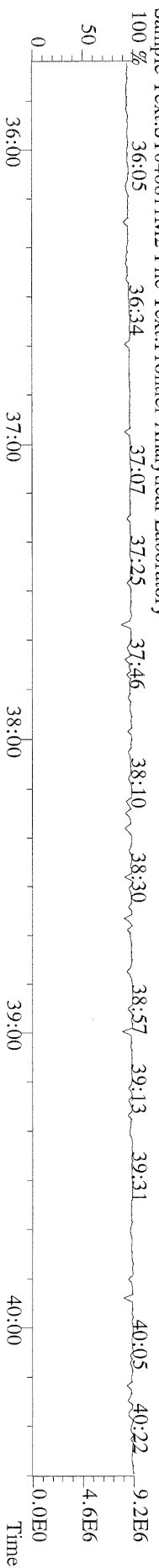
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389.8156 S:16 F:3 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
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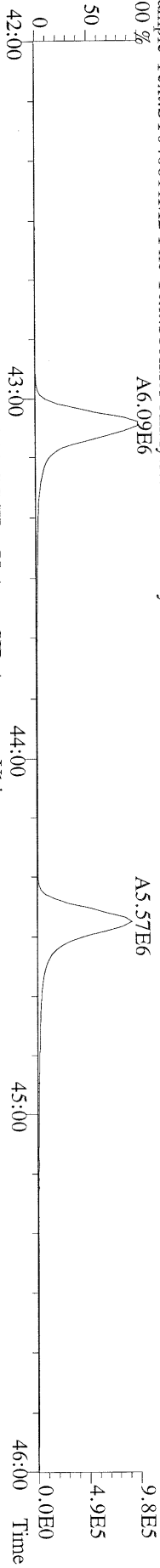
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401.8559 S:16 F:3 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



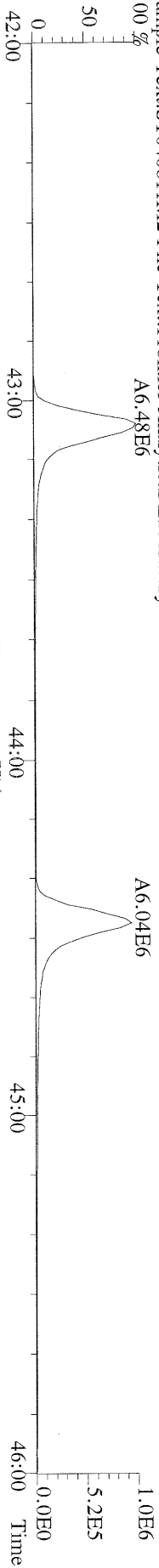
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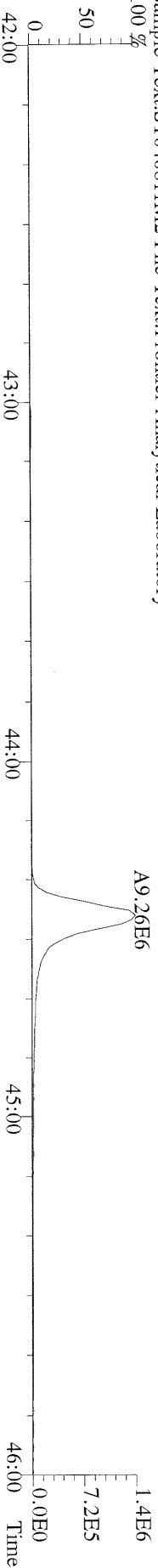
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423.7767 S:16 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



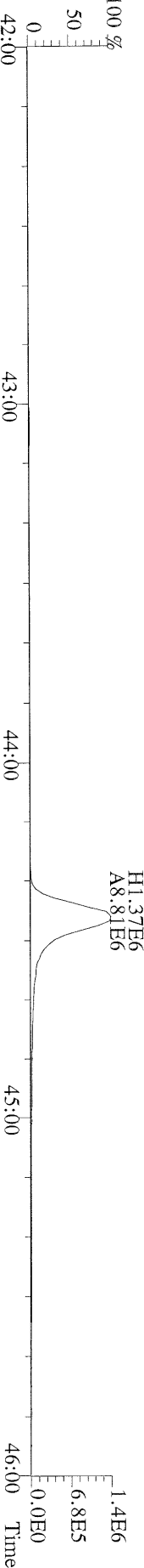
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425.7737 S:16 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



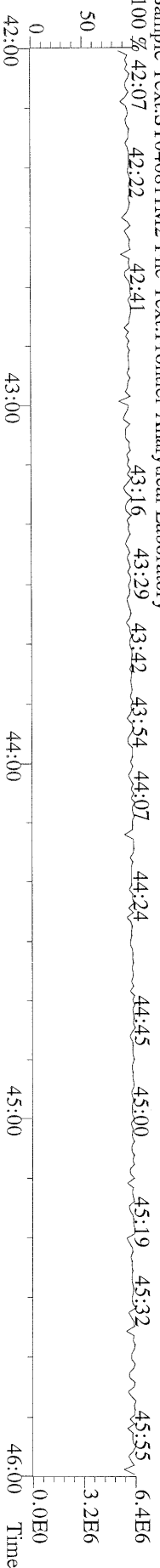
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



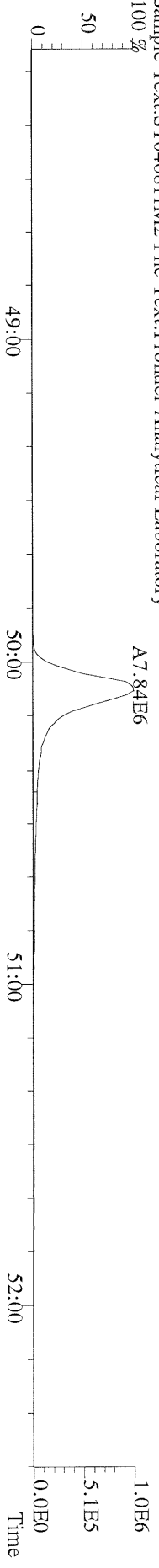
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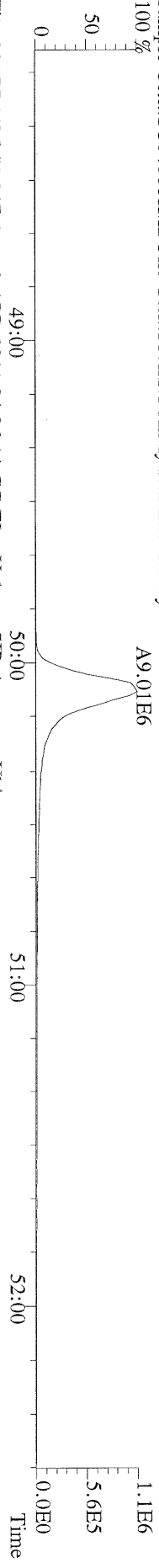
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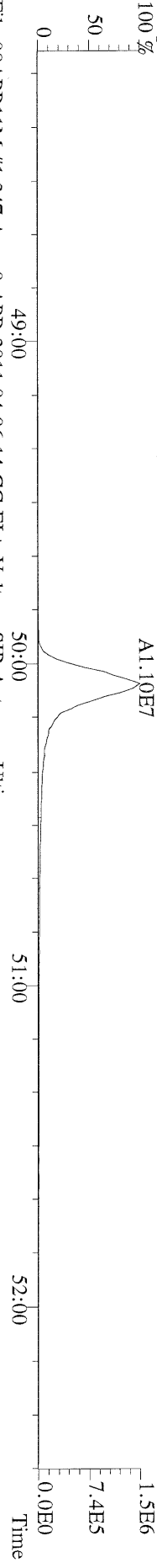
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457.7377 S:16 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



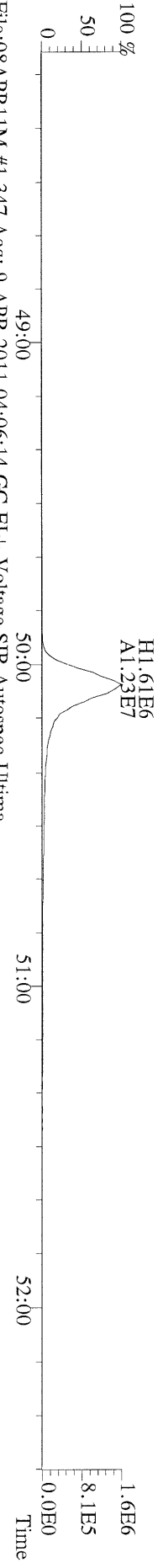
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



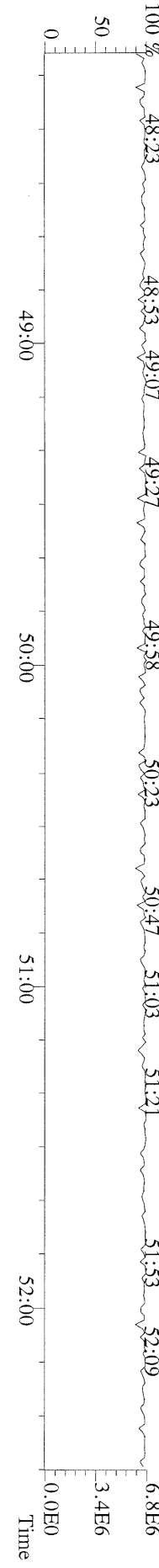
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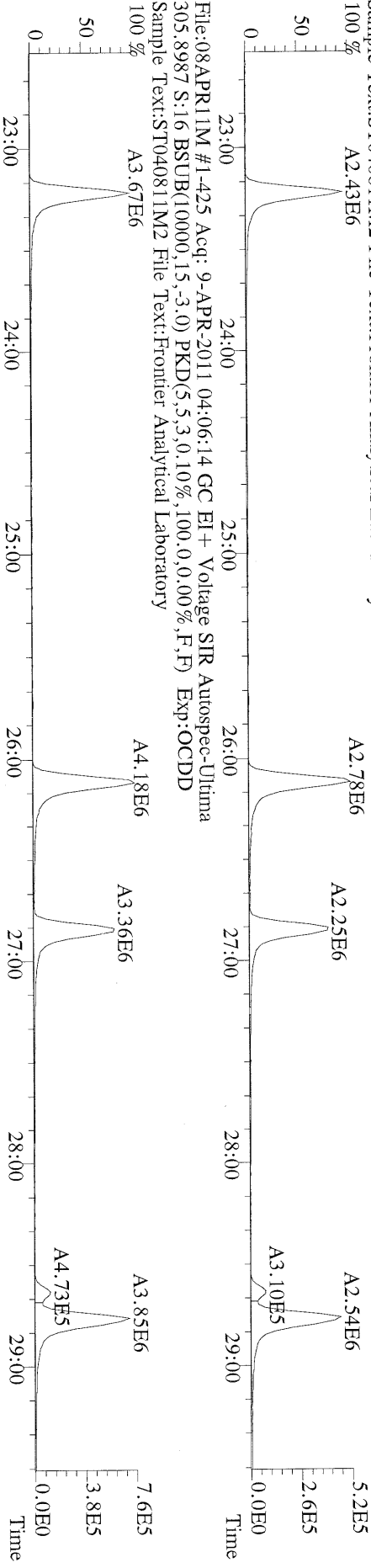
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



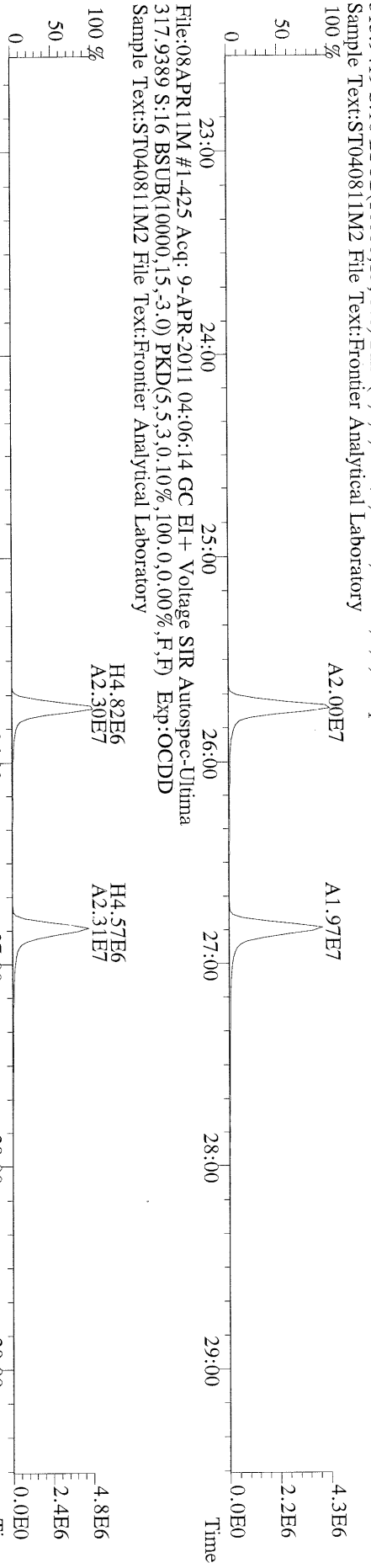
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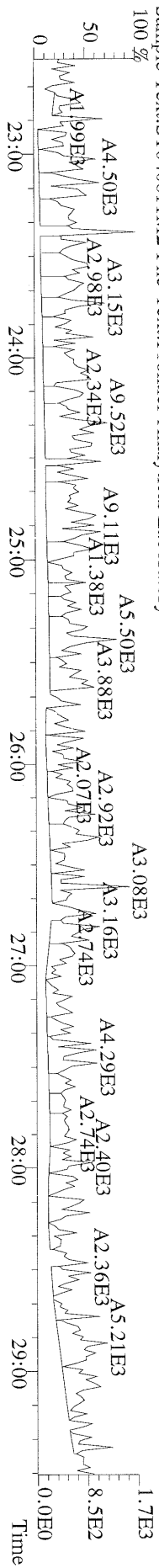
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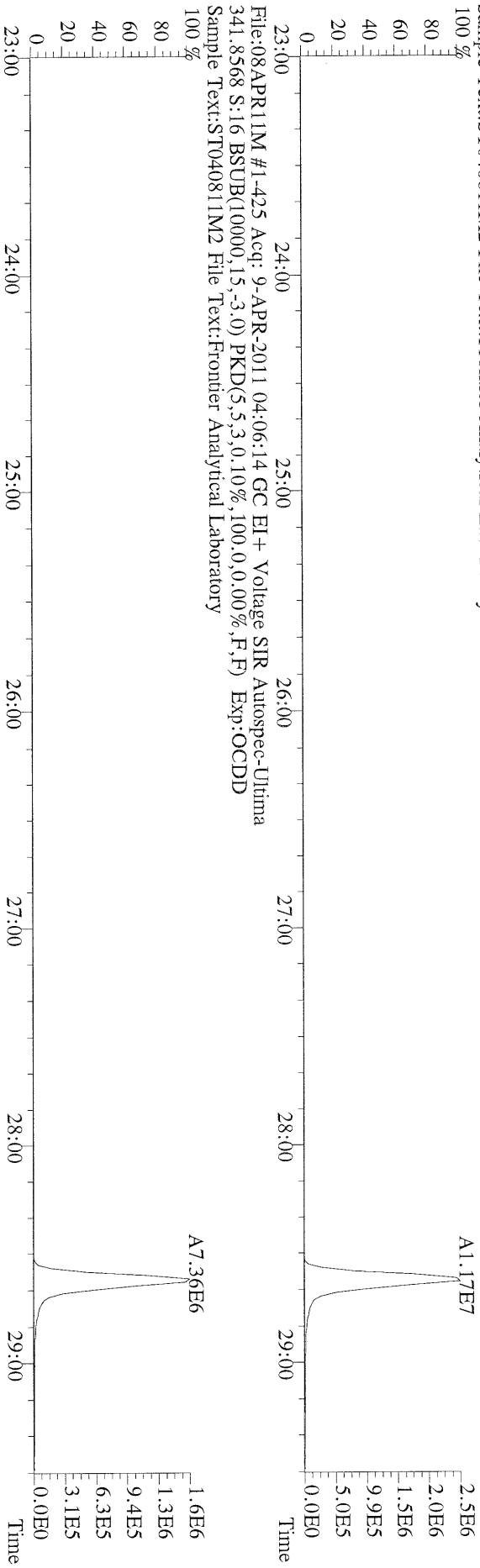
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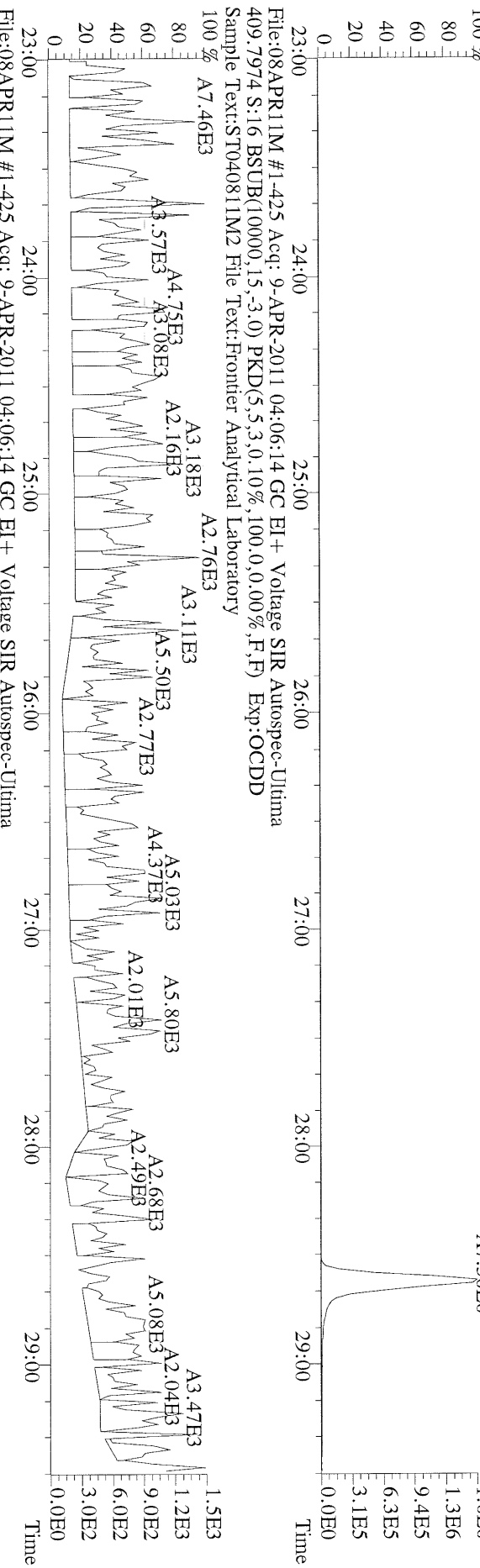
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 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



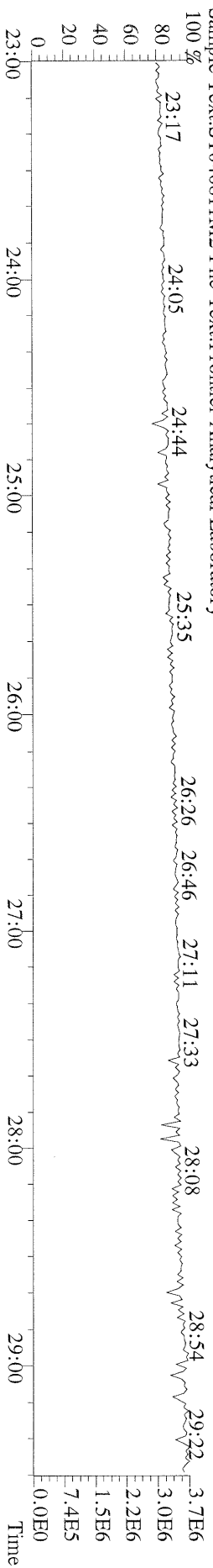
File:08APR11M #1-425 Acq: 9-APR-2011 04:06:14 GC EI+ Voltage SIR Autospec-Ultima
 339.8597 S:16 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0,0%,F,F) Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



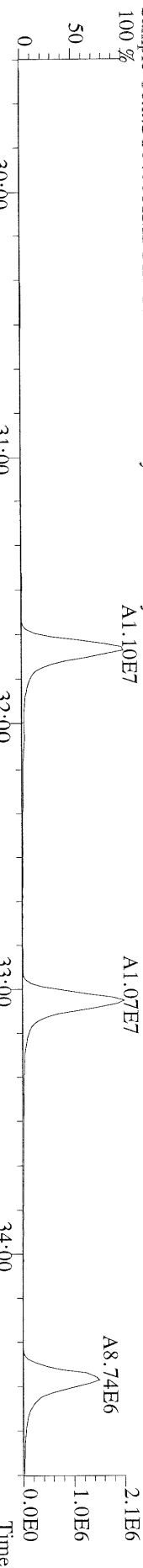
File:08APR11M #1-425 Acq: 9-APR-2011 04:06:14 GC EI+ Voltage SIR Autospec-Ultima
 341.8568 S:16 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0,0%,F,F) Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



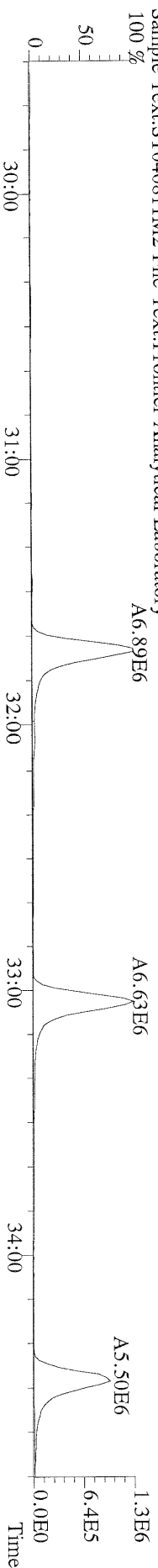
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 316.9824 S:16 Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



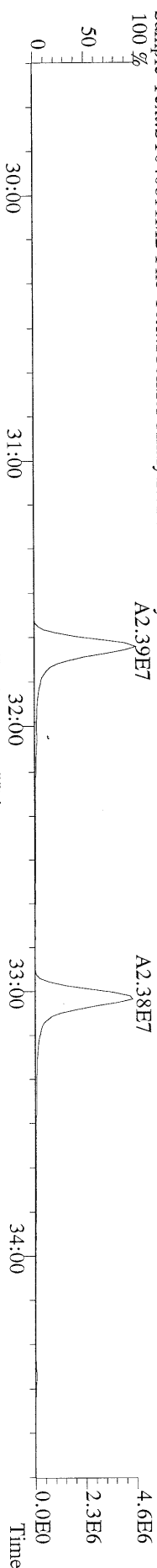
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Sample Text:ST04081IM2 File Text:Frontier Analytical Laboratory



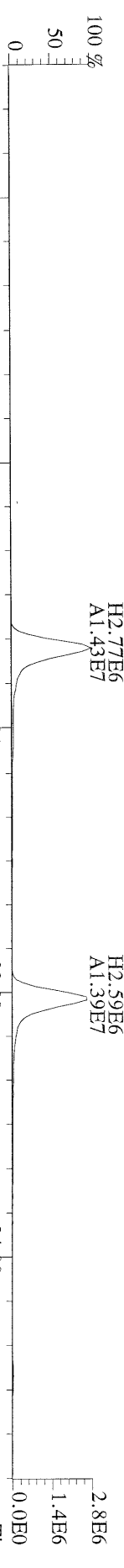
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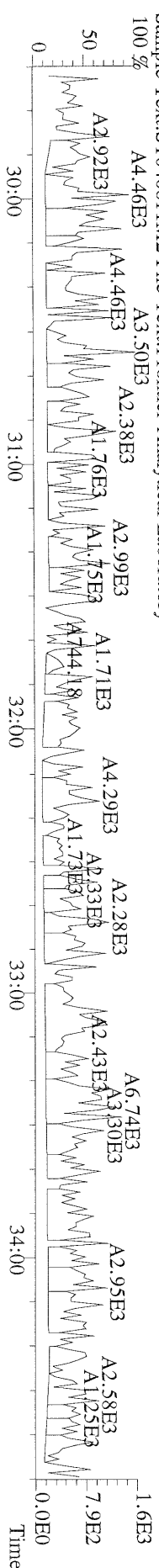
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351.9000 S:16 F:2 B SUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04081IM2 File Text:Frontier Analytical Laboratory



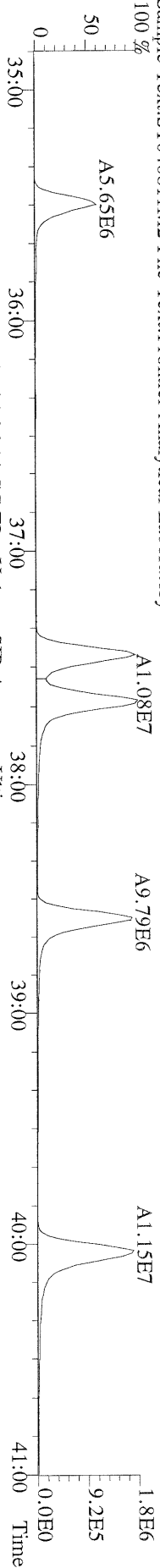
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353.8970 S:16 F:2 B SUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04081IM2 File Text:Frontier Analytical Laboratory



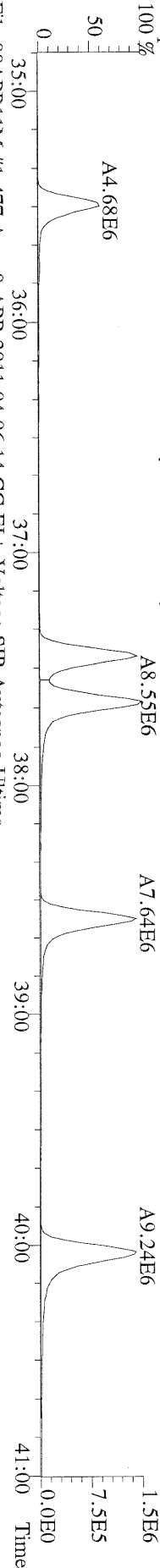
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409.7974 S:16 F:2 B SUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04081IM2 File Text:Frontier Analytical Laboratory



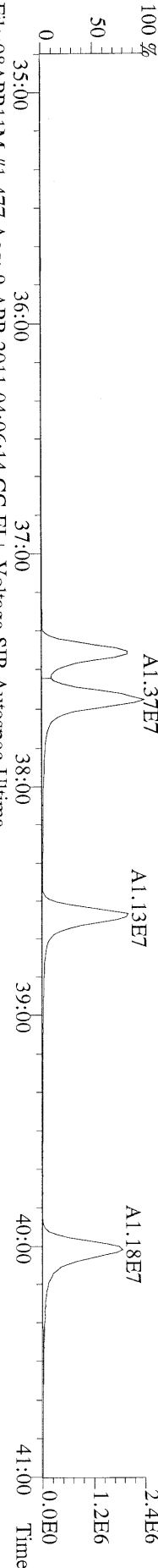
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 373.8207 S:16 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



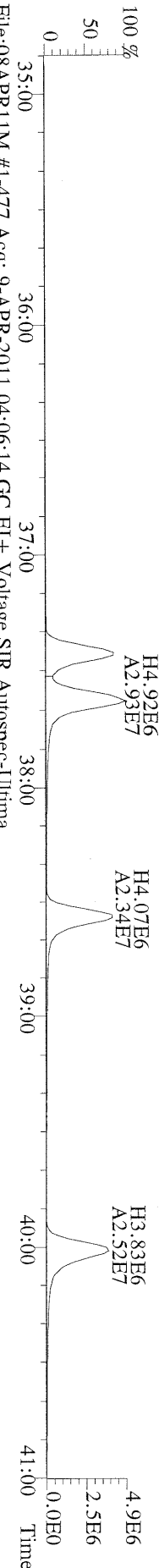
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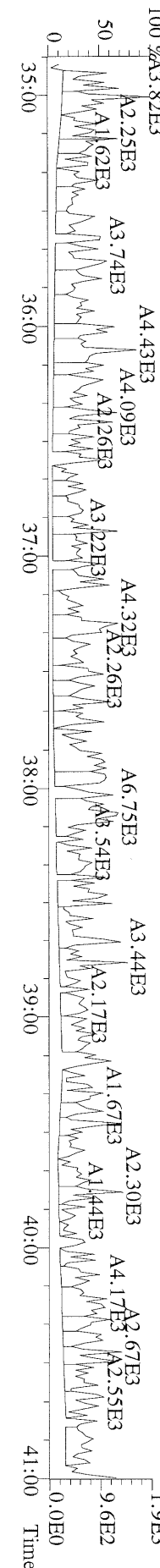
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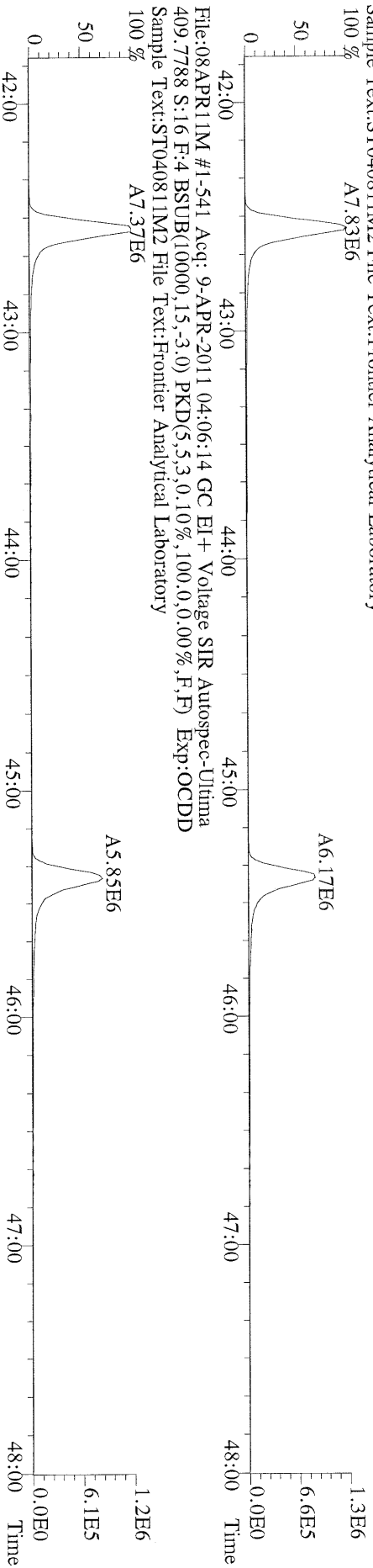
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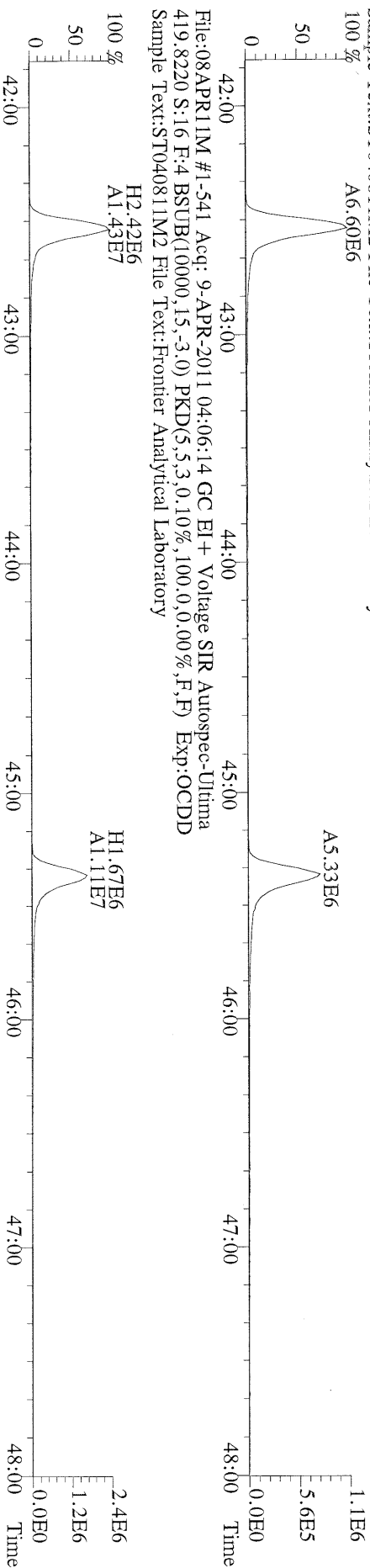
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 445.7555 S:16 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
 Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



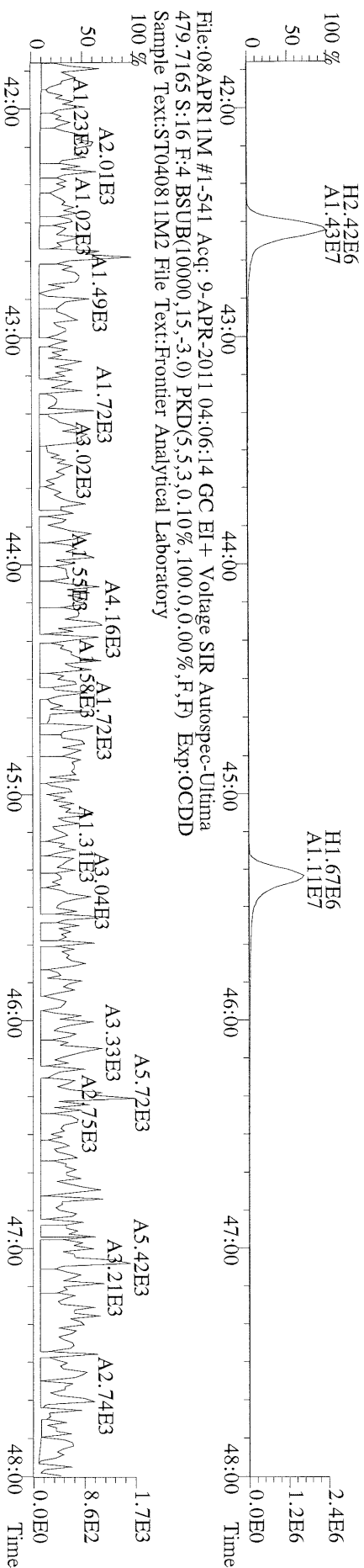
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407.7818 S:16 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



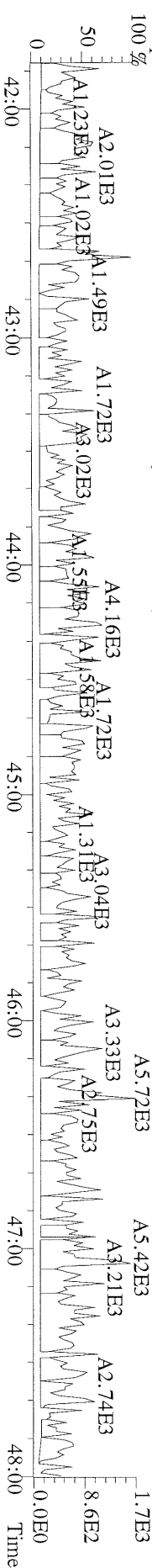
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



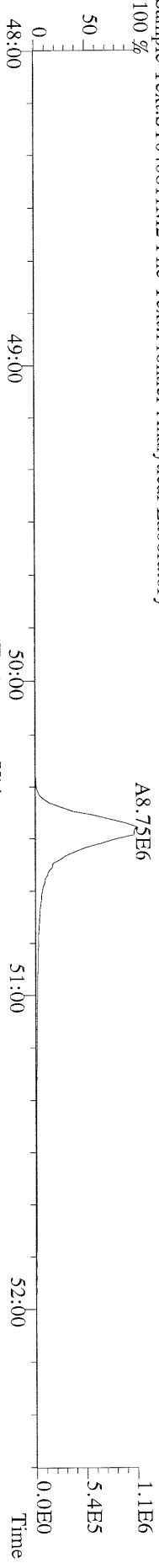
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



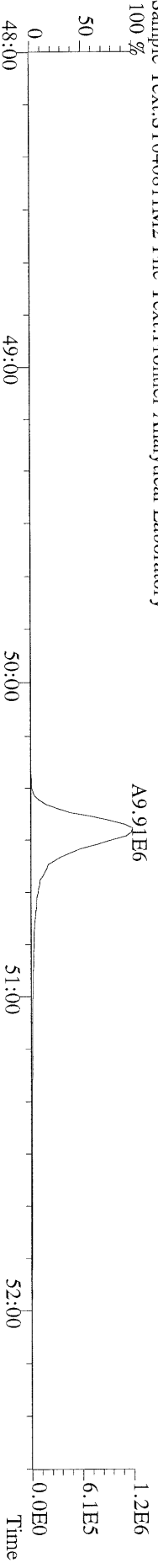
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479.7165 S:16 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



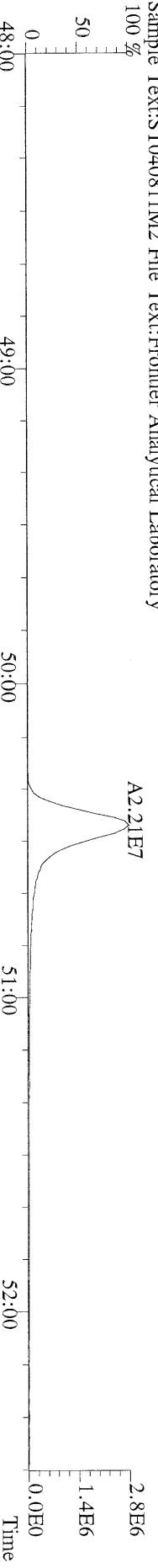
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



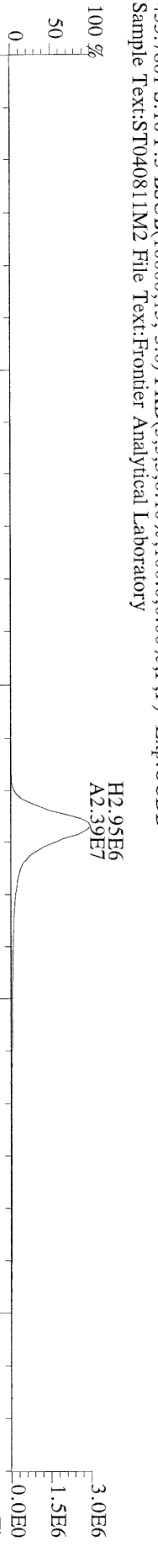
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



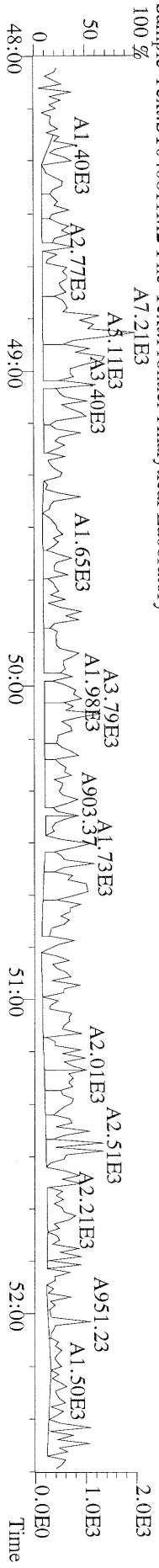
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory



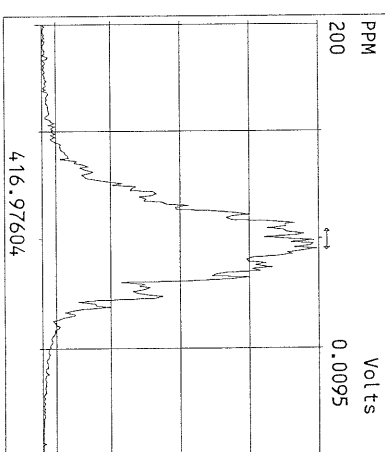
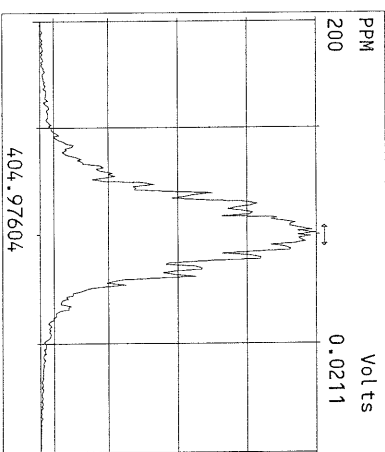
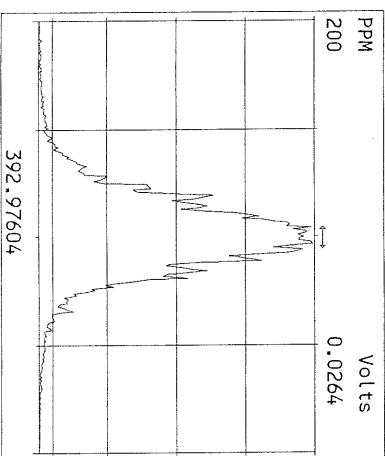
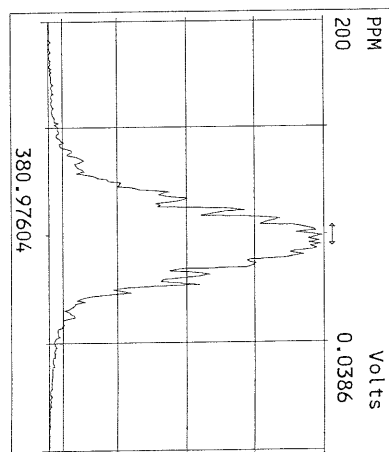
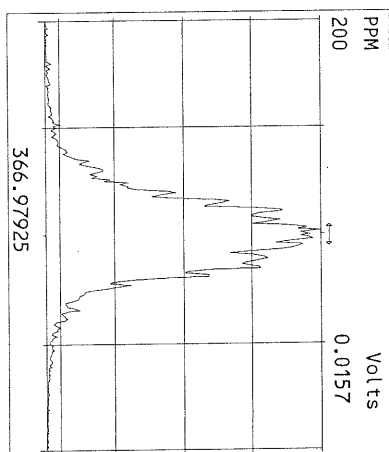
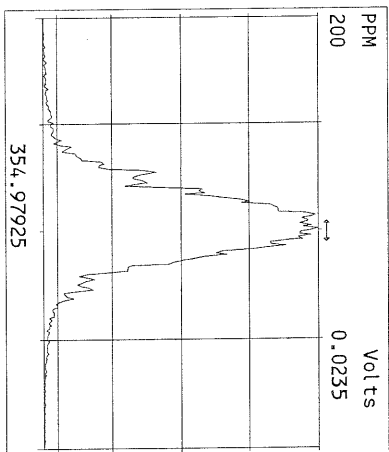
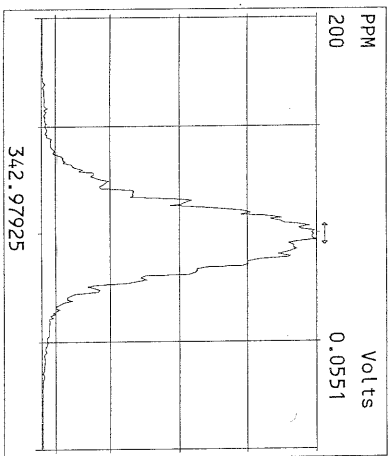
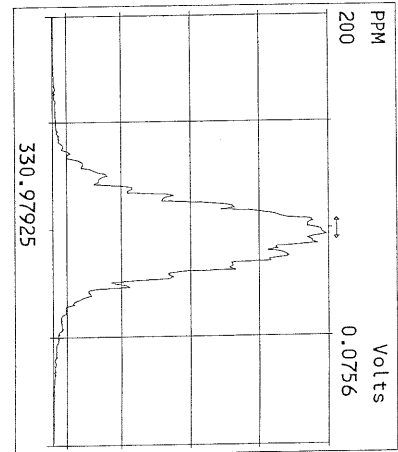
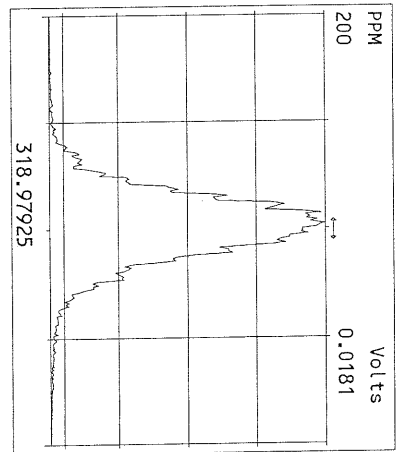
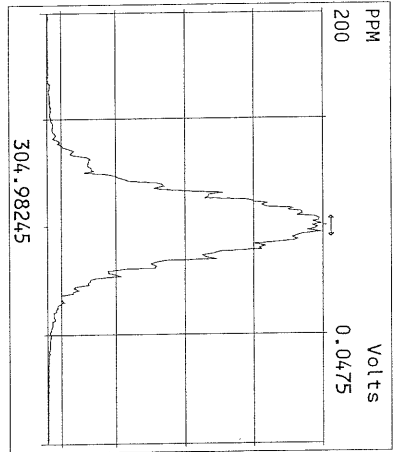
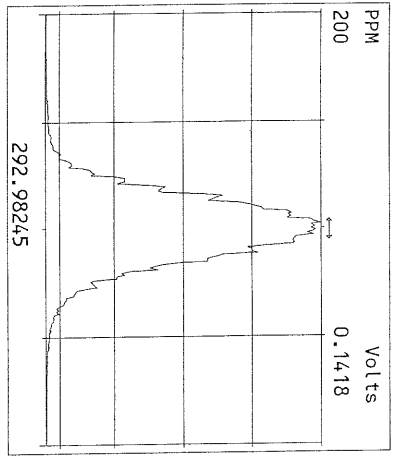
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory

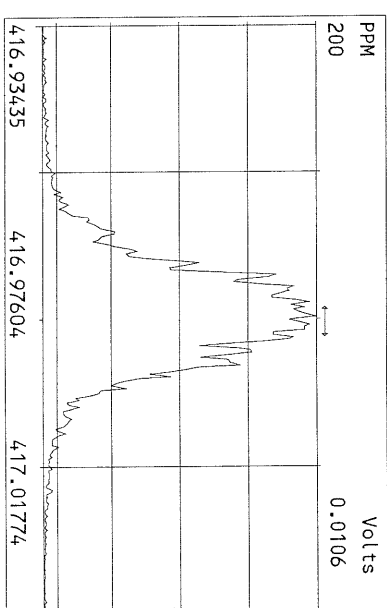
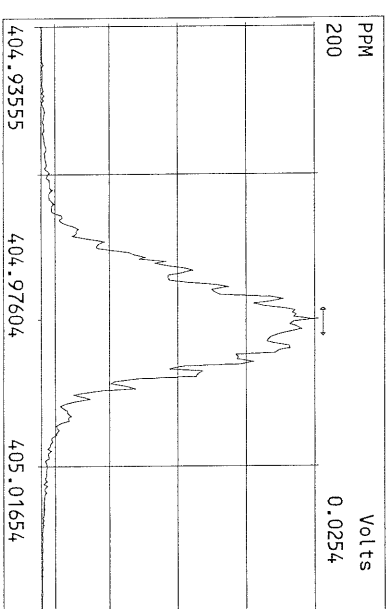
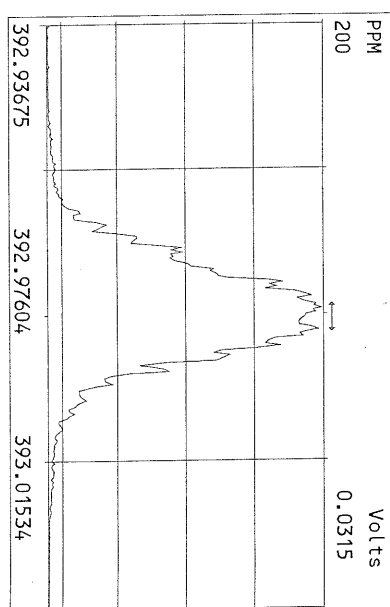
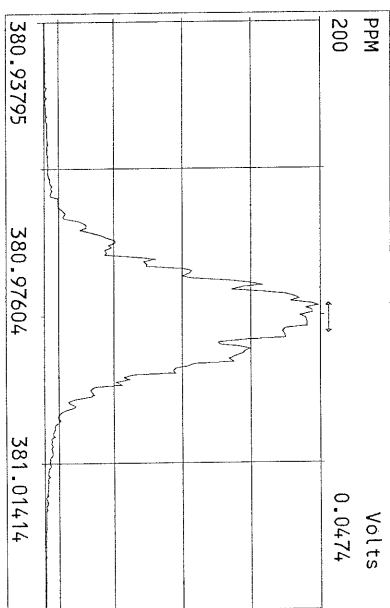
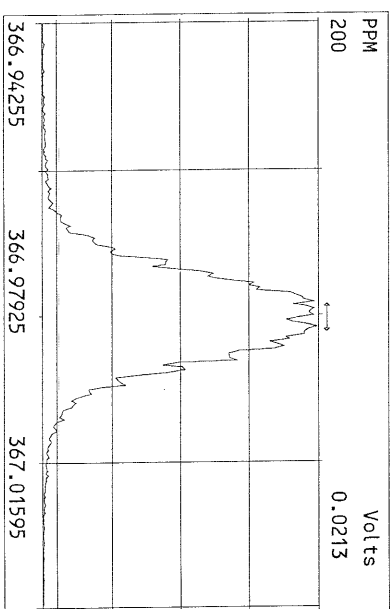
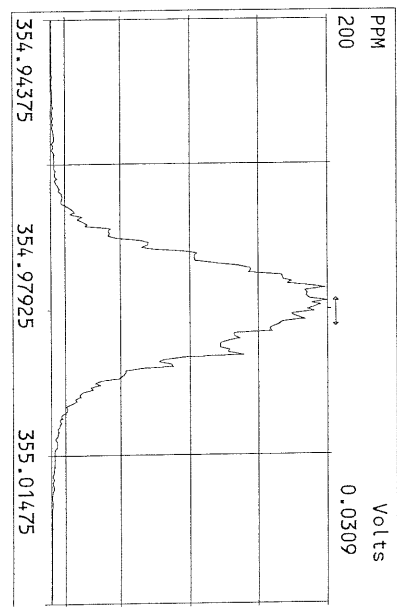
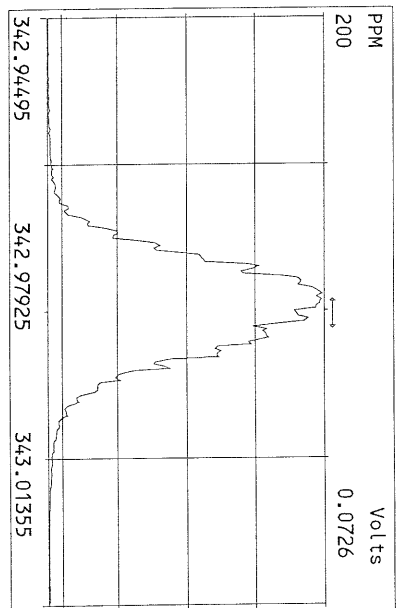
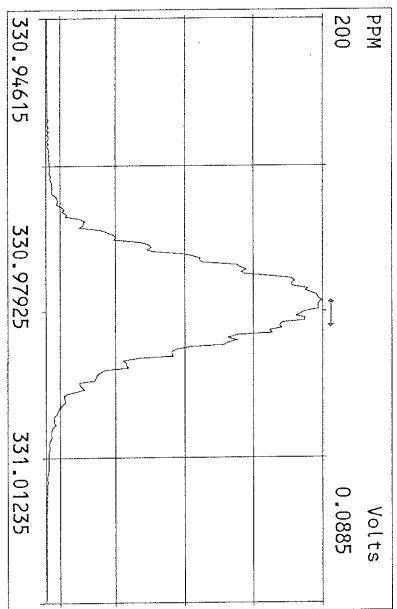


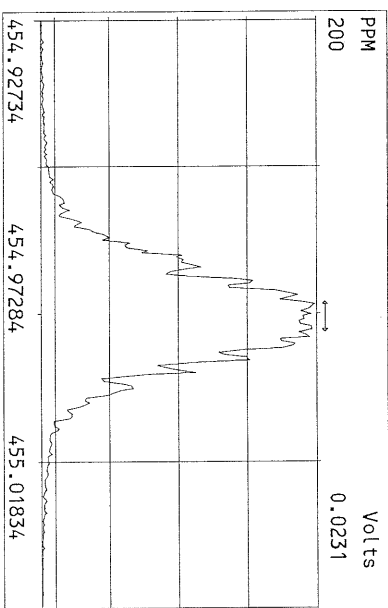
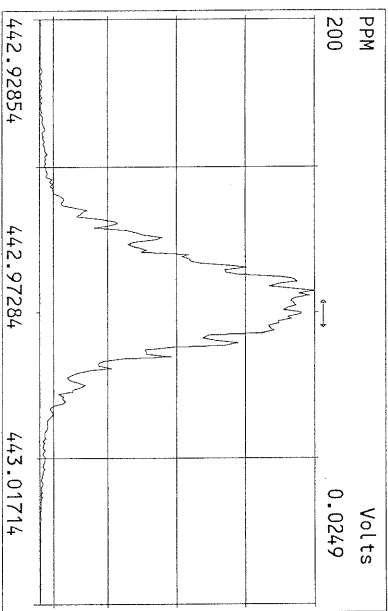
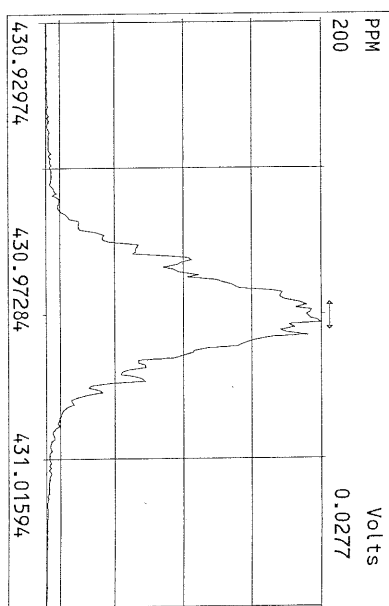
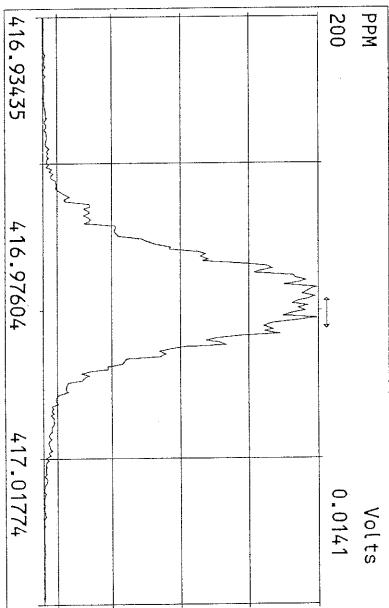
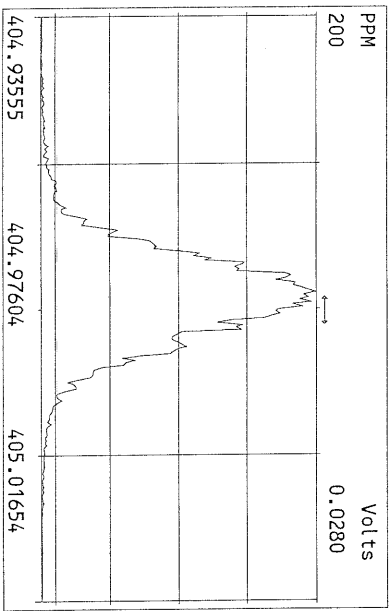
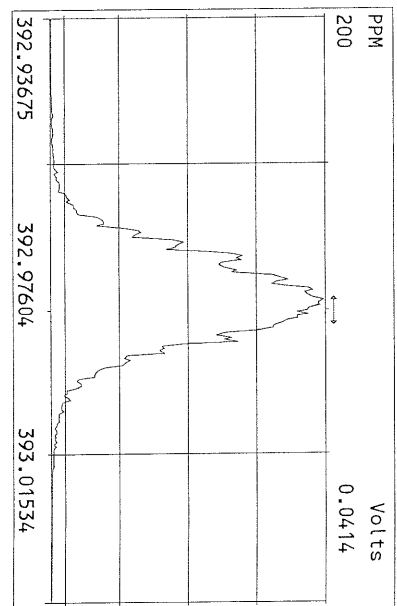
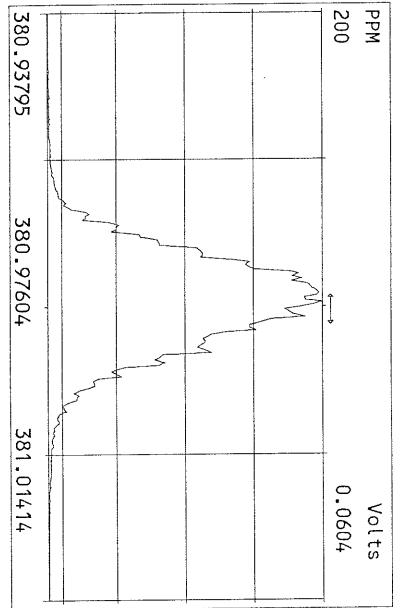
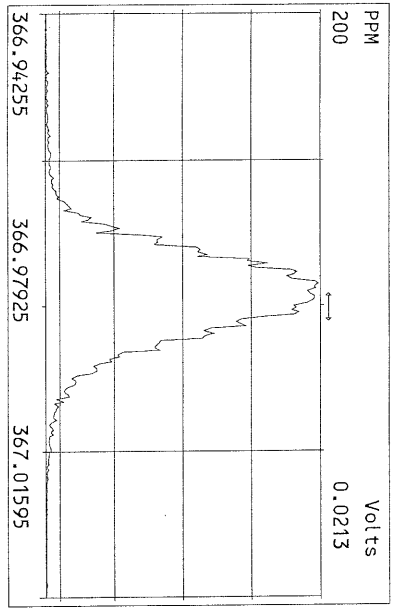
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Sample Text:ST040811M2 File Text:Frontier Analytical Laboratory

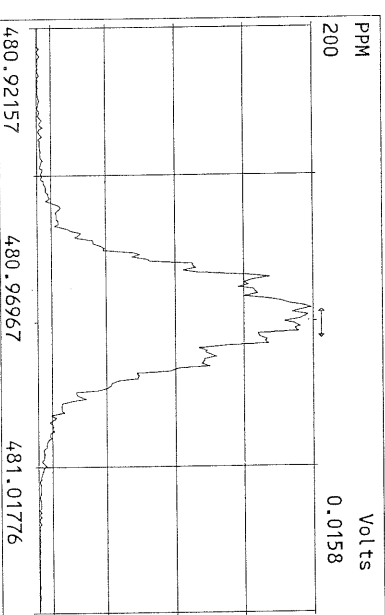
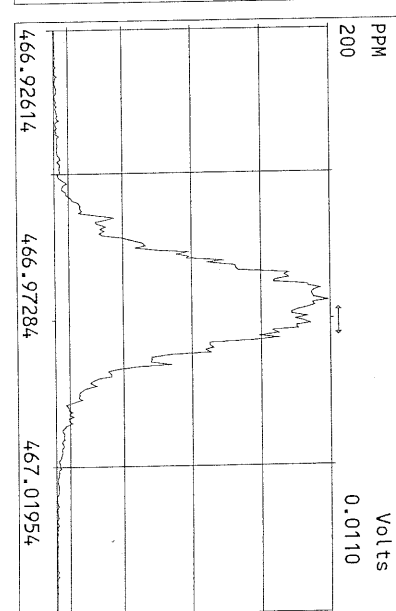
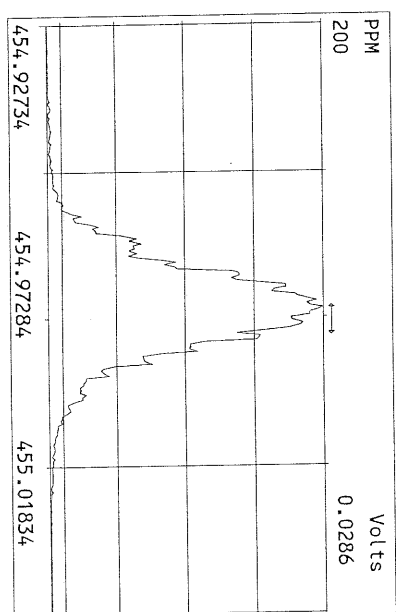
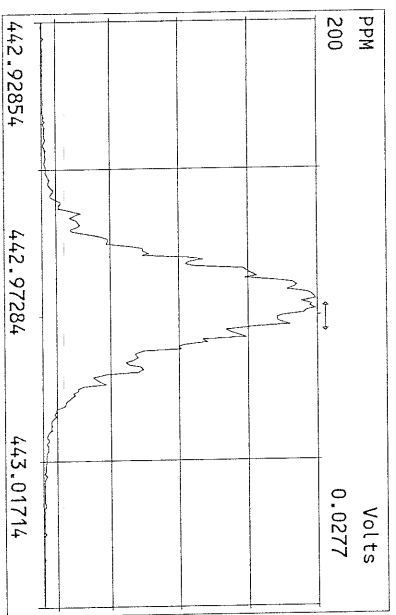
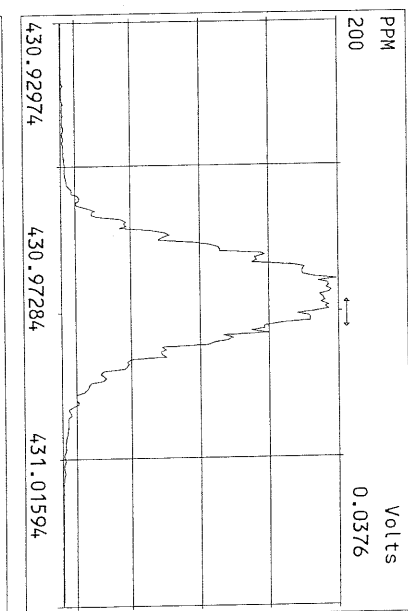
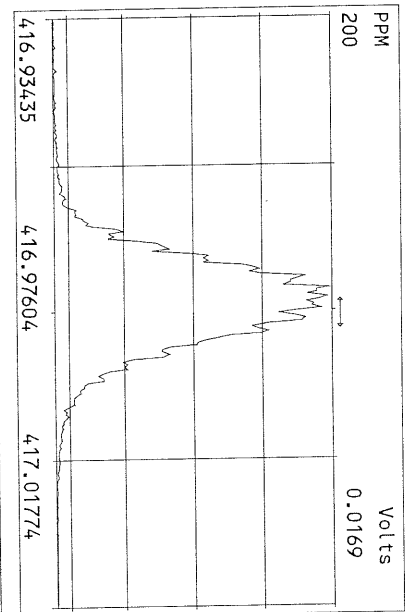
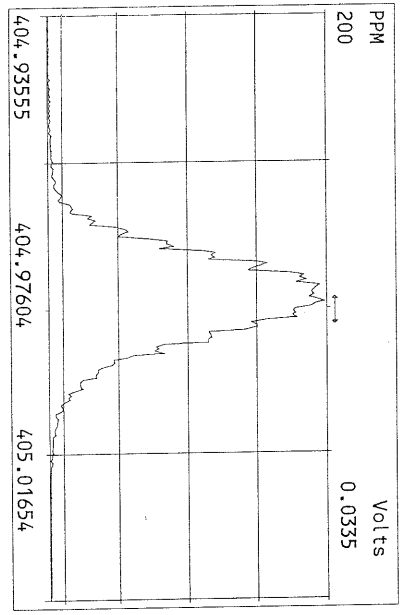


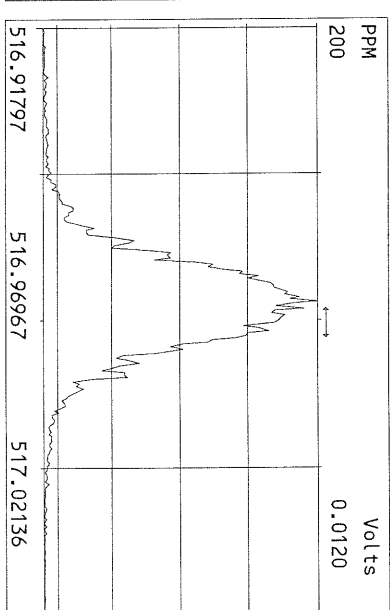
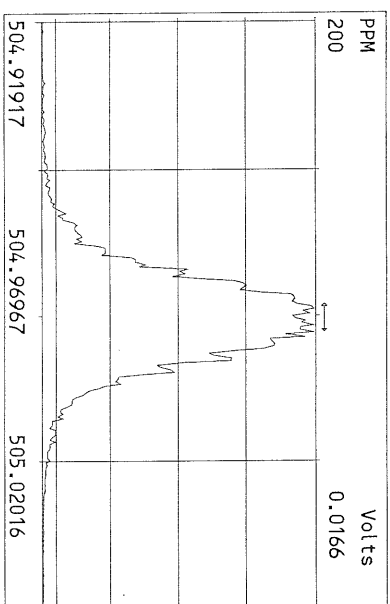
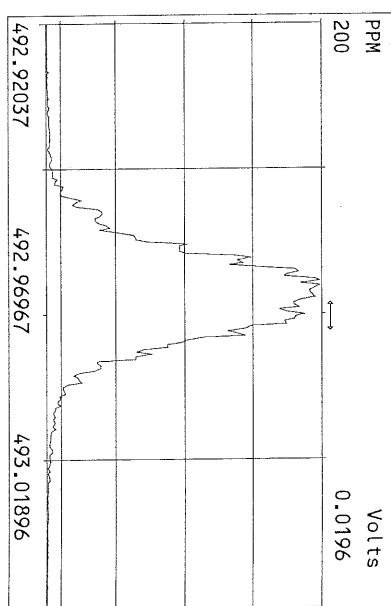
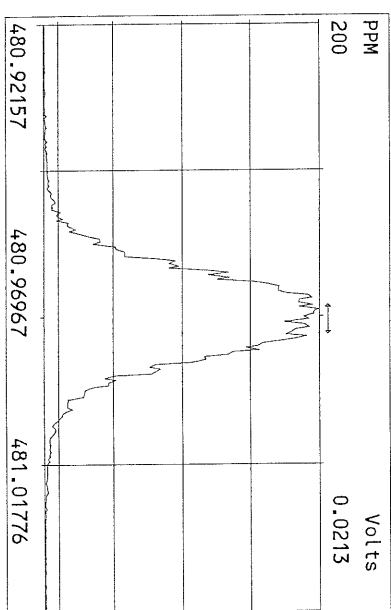
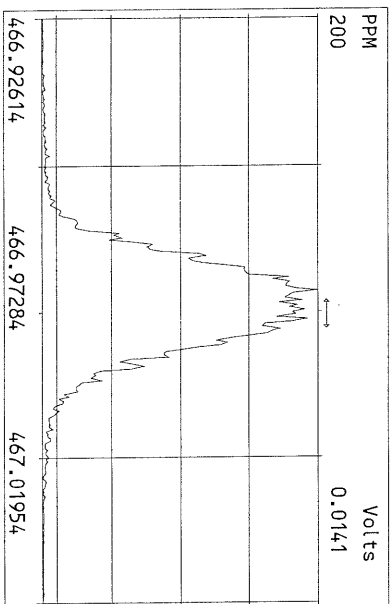
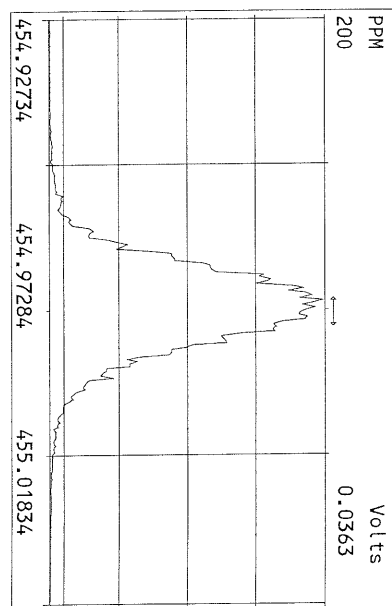
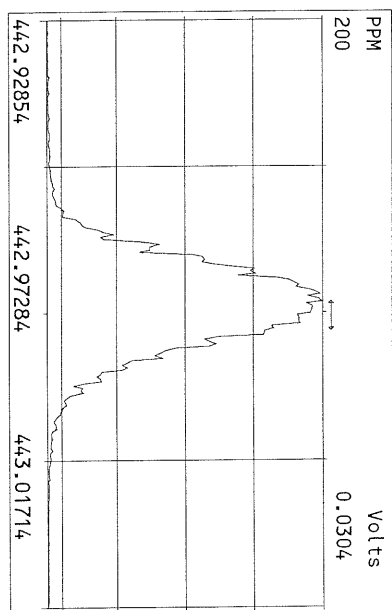
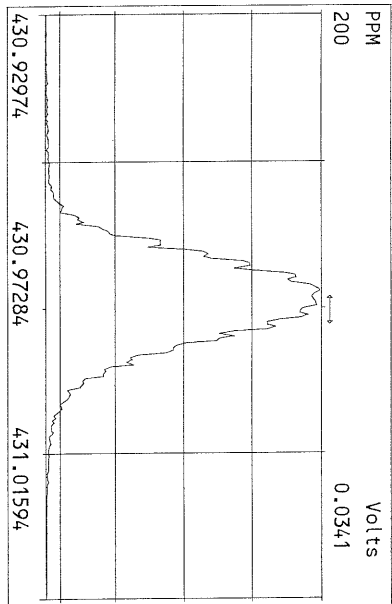
Peak Locate Examination: 9-APR-2011:08:31 File:08APR11M_RES_CHECK
Experiment:OCD Function:1 Reference:PFK











USEPA - ITD

FORM 4A

PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 20APR11M Sam:1

Analysis Date: 20-APR-11 10:16:17

NATIVE ANALYTES	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
2,3,7,8-TCDD	M/M+2	0.77	0.65-0.89	y	9.22	7.80 - 12.9 ✓
1,2,3,7,8-PeCDD	M+2/M+4	1.64	1.32-1.78	y	51.1	39.0 - 65.0 ✓
1,2,3,4,7,8-HxCDD	M+2/M+4	1.33	1.05-1.43	y	50.6	39.0 - 64.0 ✓
1,2,3,6,7,8-HxCDD	M+2/M+4	1.32	1.05-1.43	y	49.1	39.0 - 64.0 ✓
1,2,3,7,8,9-HxCDD	M+2/M+4	1.33	1.05-1.43	y	51.7	41.0 - 61.0 ✓
1,2,3,4,6,7,8-HpCDD	M+2/M+4	0.94	0.88-1.20	y	47.7	43.0 - 58.0 ✓
OCDD	M+2/M+4	0.93	0.76-1.02	y	98.0	79.0 - 126 ✓
2,3,7,8-TCDF	M/M+2	0.67	0.65-0.89	y	10.7	8.40 - 12.0 ✓
1,2,3,7,8-PeCDF	M+2/M+4	1.66	1.32-1.78	y	52.4	41.0 - 60.0 ✓
2,3,4,7,8-PeCDF	M+2/M+4	1.62	1.32-1.78	y	51.0	41.0 - 60.0 ✓
1,2,3,4,7,8-HxCDF	M+2/M+4	1.24	1.05-1.43	y	48.9	45.0 - 56.0 ✓
1,2,3,6,7,8-HxCDF	M+2/M+4	1.27	1.05-1.43	y	49.5	44.0 - 57.0 ✓
2,3,4,6,7,8-HxCDF	M+2/M+4	1.24	1.05-1.43	y	49.1	44.0 - 57.0 ✓
1,2,3,7,8,9-HxCDF	M+2/M+4	1.28	1.05-1.43	y	49.5	45.0 - 56.0 ✓
1,2,3,4,6,7,8-HpCDF	M+2/M+4	1.05	0.88-1.20	y	47.6	45.0 - 55.0 ✓
1,2,3,4,7,8,9-HpCDF	M+2/M+4	1.06	0.88-1.20	y	48.6	43.0 - 58.0 ✓
OCDF	M+2/M+4	0.87	0.76-1.02	y	90.0	63.0 - 159 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

Analyst: JDate: 4/21/11

USEPA - ITD

FORM 4B
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 20APR11M Sam:1

Analysis Date: 20-APR-11 10:16:17

LABELLED COMPOUNDS	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
13C-2,3,7,8-TCDD	M/M+2	0.80	0.65-0.89	y	94.3	82.0 - 121 ✓
13C-1,2,3,7,8-PeCDD	M+2/M+4	1.77	1.32-1.78	y	80.3	62.0 - 160 ✓
13C-1,2,3,4,7,8-HxCDD	M+2/M+4	1.28	1.05-1.43	y	96.0	85.0 - 117 ✓
13C-1,2,3,6,7,8-HxCDD	M+2/M+4	1.27	1.05-1.43	y	104	85.0 - 118 ✓
13C-1,2,3,4,6,7,8-HpCDD	M+2/M+4	1.04	0.88-1.20	y	123	72.0 - 138 ✓
13C-OCDD	M+2/M+4	0.96	0.76-1.02	y	230	96.0 - 415 ✓
13C-2,3,7,8-TCDF	M/M+2	0.88	0.65-0.89	y	98.2	71.0 - 140 ✓
13C-1,2,3,7,8-PeCDF	M+2/M+4	1.73	1.32-1.78	y	94.7	76.0 - 130 ✓
13C-2,3,4,7,8-PeCDF	M+2/M+4	1.71	1.32-1.78	y	95.2	77.0 - 130 ✓
13C-1,2,3,4,7,8-HxCDF	M/M+2	0.47	0.43-0.59	y	99.5	76.0 - 131 ✓
13C-1,2,3,6,7,8-HxCDF	M/M+2	0.49	0.43-0.59	y	102	70.0 - 143 ✓
13C-2,3,4,6,7,8-HxCDF	M/M+2	0.49	0.43-0.59	y	103	73.0 - 137 ✓
13C-1,2,3,7,8,9-HxCDF	M/M+2	0.48	0.43-0.59	y	107	74.0 - 135 ✓
13C-1,2,3,4,6,7,8-HpCDF	M/M+2	0.44	0.37-0.51	y	116	78.0 - 129 ✓
13C-1,2,3,4,7,8,9-HpCDF	M/M+2	0.46	0.37-0.51	y	116	77.0 - 129 ✓
13C-OCDF	M+2/M+4	0.93	0.76-1.02	y	227	96.0 - 415 ✓
CLEANUP STANDARD (4)						
37Cl-2,3,7,8-TCDD					7.92	7.80 - 12.8 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

(4) No ion abundance ratio; report concentration found.

Analyst: *[Signature]*

Date: *4/21/11*

USEPA - ITD

FORM 6A

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 20-APR-11 10:16:17

CS3 or VER Data Filename: 20APR11M

Sam:1

NATIVE ANALYTES	RETENTION TIME	RRT	RRT
	REFERENCE		QC LIMITS (1)
2,3,7,8-TCDD	13C-2,3,7,8-TCDD	1.001	0.999-1.002 ✓
2,3,7,8-TCDF	13C-2,3,7,8-TCDF	1.001	0.999-1.003 ✓
1,2,3,7,8-PeCDD	13C-1,2,3,7,8-PeCDD	1.001	0.999-1.002 ✓
1,2,3,7,8-PeCDF	13C-1,2,3,7,8-PeCDF	1.001	0.999-1.002 ✓
2,3,4,7,8-PeCDF	13C-2,3,4,7,8-PeCDF	1.000	0.999-1.002 ✓
LABELED COMPOUNDS			
37Cl-2,3,7,8-TCDD	13C-1,2,3,4-TCDD	1.022	0.989-1.052 ✓
13C-2,3,7,8-TCDD		1.022	0.976-1.043 ✓
13C-2,3,7,8-TCDF		0.993	0.923-1.103 ✓
13C-1,2,3,7,8-PeCDD		1.239	1.000-1.567 ✓
13C-1,2,3,7,8-PeCDF		1.174	0.923-1.203 ✓
13C-2,3,4,7,8-PeCDF		1.224	0.923-1.303 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified in Table 2, Method 1613.

Analyst: _____

Date: _____

FORM 6B
PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 20-APR-11 10:16:17 CS3 or VER Data Filename: 20APR11M Sam:1

NATIVE ANALYTES	RETENTION TIME REFERENCE	RRT	RRT QC LIMITS (1)
1,2,3,4,7,8-HxCDD	13C-1,2,3,4,7,8-HxCDD	1.001	0.999-1.001 ✓
1,2,3,6,7,8-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.000	0.998-1.004 ✓
1,2,3,7,8,9-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.012	1.000-1.019 ✓
1,2,3,4,7,8-HxCDF	13C-1,2,3,4,7,8-HxCDF	1.000	0.999-1.001 ✓
1,2,3,6,7,8-HxCDF	13C-1,2,3,6,7,8-HxCDF	1.001	0.997-1.005 ✓
2,3,4,6,7,8-HxCDF	13C-2,3,4,6,7,8-HxCDF	1.001	0.999-1.001 ✓
1,2,3,7,8,9-HxCDF	13C-1,2,3,7,8,9-HxCDF	1.001	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDD	13C-1,2,3,4,6,7,8-HpCDD	1.000	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDF	13C-1,2,3,4,6,7,8-HpCDF	1.001	0.999-1.001 ✓
1,2,3,4,7,8,9-HpCDF	13C-1,2,3,4,7,8,9-HpCDF	1.000	0.999-1.001 ✓
OCDD	13C-OCDD	1.001	0.999-1.001 ✓
OCDF	13C-OCDF	1.001	0.999-1.001 ✓
LABELED COMPOUNDS			
13C-1,2,3,4,7,8-HxCDD	13C-1,2,3,7,8,9-HxCDD	0.984	0.977-1.000 ✓
13C-1,2,3,6,7,8-HxCDD		0.989	0.981-1.003 ✓
13C-1,2,3,4,7,8-HxCDF		0.949	0.944-0.970 ✓
13C-1,2,3,6,7,8-HxCDF		0.954	0.949-0.975 ✓
13C-2,3,4,6,7,8-HxCDF		0.978	0.959-1.021 ✓
13C-1,2,3,7,8,9-HxCDF		1.015	0.977-1.047 ✓
13C-1,2,3,4,6,7,8-HpCDD		1.128	1.086-1.130 ✓
13C-1,2,3,4,6,7,8-HpCDF		1.079	1.043-1.085 ✓
13C-1,2,3,4,7,8,9-HpCDF		1.151	1.057-1.154 ✓
13C-OCDD		1.269	1.032-1.311 ✓
13C-OCDF		1.279	1.000-1.311 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified
in Table 2, Method 1613.

Analyst: Date: 4/21/11

FAL ID: ST042011M1 Filename: 20APR11M Sam:1 Acquired: 20-APR-11 10:16:17 ICal: PCDDFAL3-3-7-11

Client ID: 1613 CS3 100511J ConCal: ST042011M1 EndCal: ST042011M2

Results: GC Column: DB5 Amount: 1.000 NATO 1989 Tox: 100 WHO 1998 Tox: 126 WHO 2005 Tox: 115

Name	Resp	RA	RT	RRF	Conc	Qual	Fac	Noise-1	Noise-2	DL
2,3,7,8-TCDD	1.94e+06	0.77 y	27:20	1.13	9.22		2.50	-	-	*
1,2,3,7,8-PeCDD	8.04e+06	1.64 y	33:10	1.02	51.1		2.50	-	-	*
1,2,3,4,7,8-HxCDD	9.25e+06	1.33 y	38:31	1.45	50.6		2.50	-	-	*
1,2,3,6,7,8-HxCDD	7.70e+06	1.32 y	38:41	1.45	49.1		2.50	-	-	*
1,2,3,7,8,9-HxCDD	8.89e+06	1.33 y	39:08	1.47	51.7		2.50	-	-	*
1,2,3,4,6,7,8-HpCDD	6.97e+06	0.94 y	44:07	1.30	47.7		2.50	-	-	*
OCDD	1.10e+07	0.93 y	49:39	1.45	98.0		2.50	-	-	*
2,3,7,8-TCDF	3.78e+06	0.67 y	26:35	1.15	10.7		2.50	-	-	*
1,2,3,7,8-PeCDF	1.17e+07	1.66 y	31:26	0.89	52.4		2.50	-	-	*
2,3,4,7,8-PeCDF	1.12e+07	1.62 y	32:44	0.89	51.0		2.50	-	-	*
1,2,3,4,7,8-HxCDF	9.97e+06	1.24 y	37:08	1.01	48.9		2.50	-	-	*
1,2,3,6,7,8-HxCDF	1.14e+07	1.27 y	37:20	0.89	49.5		2.50	-	-	*
2,3,4,6,7,8-HxCDF	1.06e+07	1.24 y	38:16	1.02	49.1		2.50	-	-	*
1,2,3,7,8,9-HxCDF	1.27e+07	1.28 y	39:42	1.10	49.5		2.50	-	-	*
1,2,3,4,6,7,8-HpCDF	8.90e+06	1.05 y	42:13	1.48	47.6		2.50	-	-	*
1,2,3,4,7,8,9-HpCDF	6.77e+06	1.06 y	45:01	1.43	48.6		2.50	-	-	*
OCDF	1.11e+07	0.87 y	50:02	0.84	90.0		2.50	-	-	*
13C-2,3,7,8-TCDD	1.85e+07	0.80 y	27:19	1.03	94.3					Rec 94.3
13C-1,2,3,7,8-PeCDD	1.55e+07	1.77 y	33:08	1.01	80.3					80.3
13C-1,2,3,4,7,8-HxCDD	1.26e+07	1.28 y	38:30	1.19	96.0					96.0
13C-1,2,3,6,7,8-HxCDD	1.08e+07	1.27 y	38:40	0.94	104					104
13C-1,2,3,4,6,7,8-HpCDD	1.12e+07	1.04 y	44:06	0.83	123					123
13C-OCDD	1.54e+07	0.96 y	49:38	0.61	230					115
13C-2,3,7,8-TCDF	3.09e+07	0.88 y	26:34	0.98	98.2					98.2
13C-1,2,3,7,8-PeCDF	2.53e+07	1.73 y	31:24	0.83	94.7					94.7
13C-2,3,4,7,8-PeCDF	2.46e+07	1.71 y	32:43	0.80	95.2					95.2
13C-1,2,3,4,7,8-HxCDF	2.02e+07	0.47 y	37:07	1.84	99.5					99.5
13C-1,2,3,6,7,8-HxCDF	2.59e+07	0.49 y	37:18	2.29	102					102
13C-2,3,4,6,7,8-HxCDF	2.12e+07	0.49 y	38:14	1.86	103					103
13C-1,2,3,7,8,9-HxCDF	2.33e+07	0.48 y	39:40	1.98	107					107
13C-1,2,3,4,6,7,8-HpCDF	1.26e+07	0.44 y	42:11	0.99	116					116
13C-1,2,3,4,7,8,9-HpCDF	9.77e+06	0.46 y	45:00	0.77	116					116
13C-OCDF	2.91e+07	0.93 y	50:00	1.17	227					113
37Cl-2,3,7,8-TCDD	1.10e+06		27:20	0.73	7.92					79.2
13C-1,2,3,4-TCDD	1.91e+07	0.81 y	26:45	-	50.2					
13C-1,2,3,4-TCDF	3.22e+07	0.87 y	25:30	-	44.7					
13C-1,2,3,7,8,9-HxCDD	1.10e+07	1.30 y	39:06	-	44.5					
Total Tetra-Dioxins	1.00e+07		22:47	1.13	47.8		2.50	-	-	* 19
Total Penta-Dioxins	1.71e+07		30:12	1.02	108		2.50	-	-	* 10
Total Hexa-Dioxins	2.88e+07		36:04	1.46	169		2.50	-	-	* 6
Total Hepta-Dioxins	1.48e+07		42:45	1.30	101		2.50	-	-	* 6
Total Tetra-Furans	1.81e+07		23:01	1.15	51.1		2.50	-	-	* 20
1st Fn. Tot Penta-Furans	1.31e+07		28:22	0.89	59.2		2.50	-	-	* PeCDF 2
Total Penta-Furans	3.29e+07		30:08	0.89	148		2.50	-	-	* 208 9
Total Hexa-Furans	5.11e+07		35:11	1.00	225		2.50	-	-	* 10
Total Hepta-Furans	1.58e+07		42:13	1.46	97.2		2.50	-	-	* 13

Analyst:

Date: 4/21/11

Frontier Analytical Laboratory - Acquisition Log

Run Name:20APR11M

Instrument: FAL3

GC: DB5

Experiment:OCDD

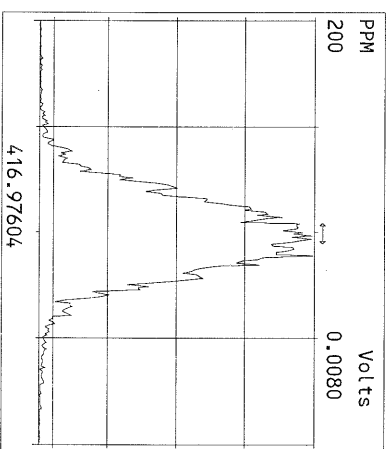
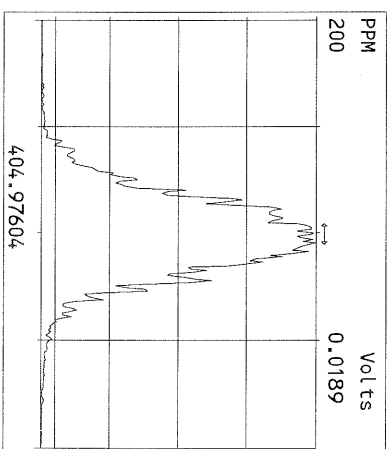
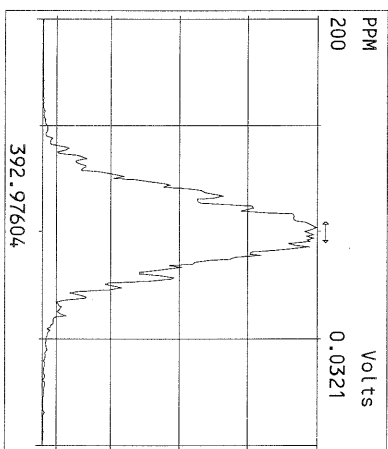
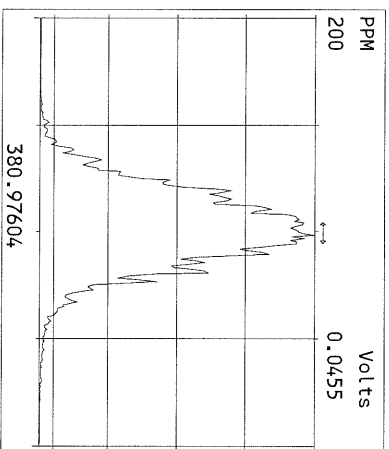
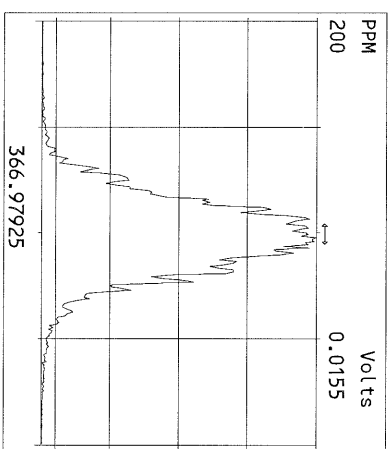
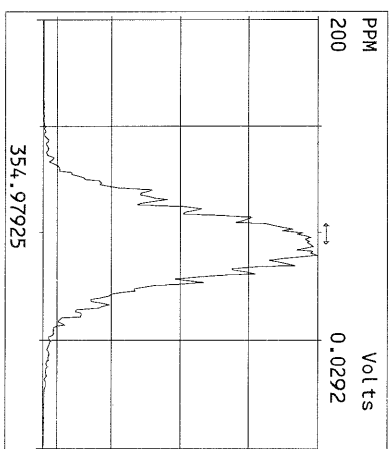
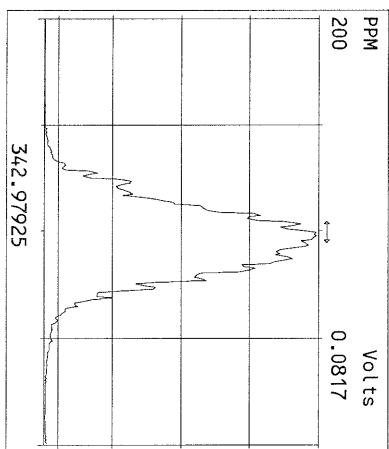
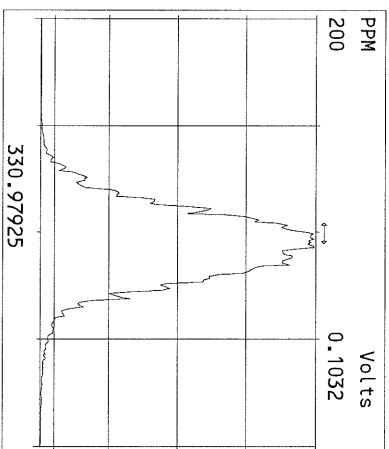
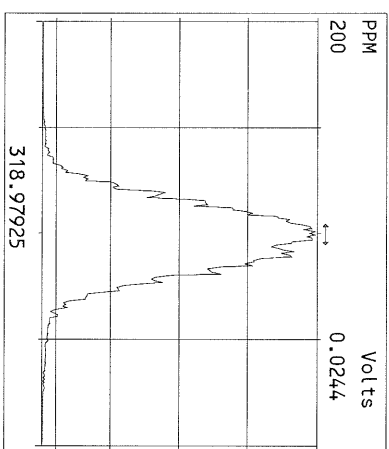
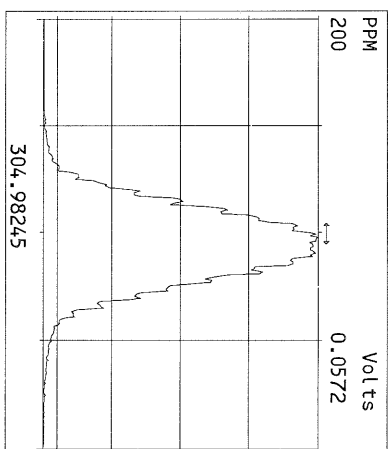
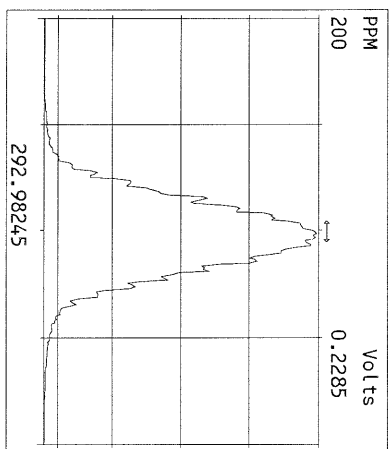
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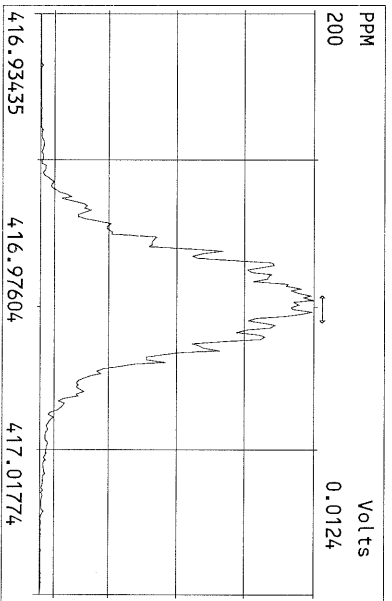
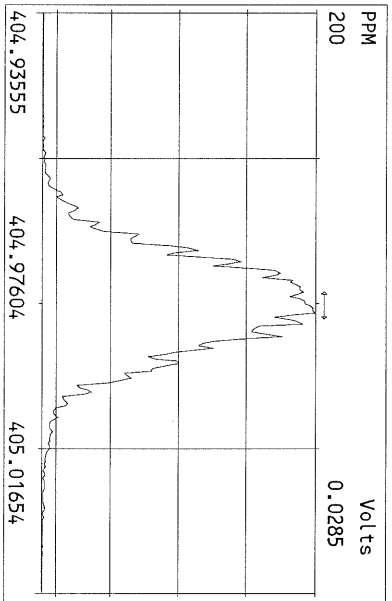
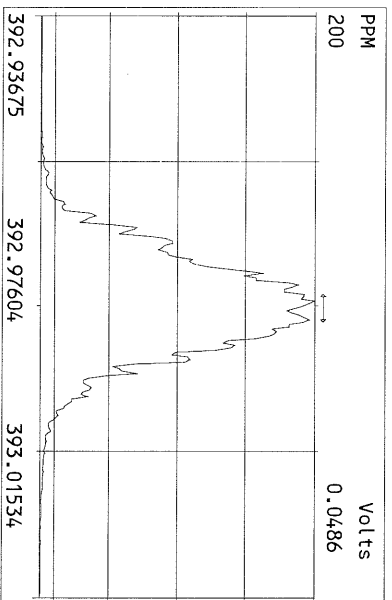
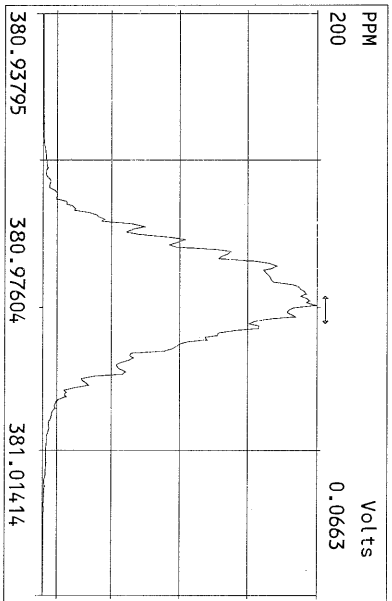
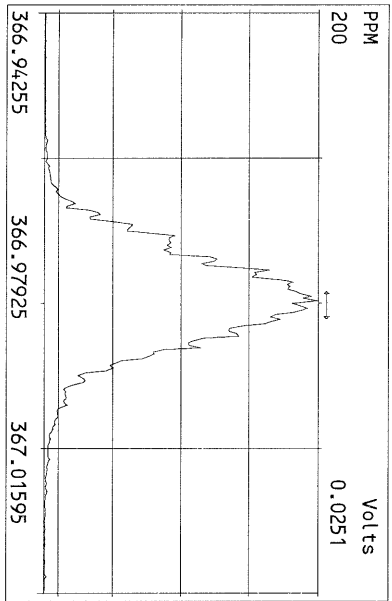
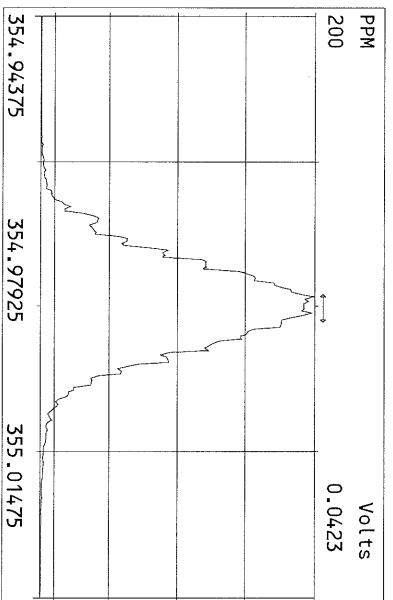
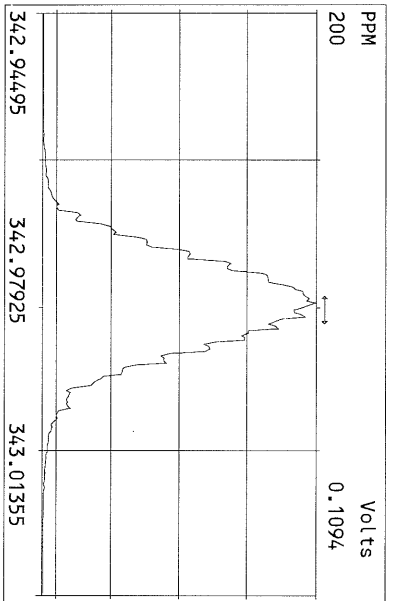
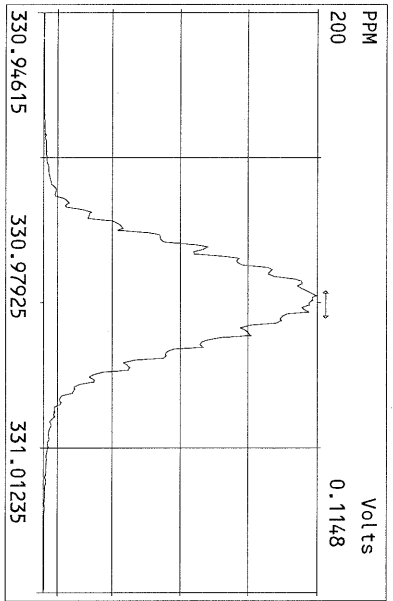
8/4/2011

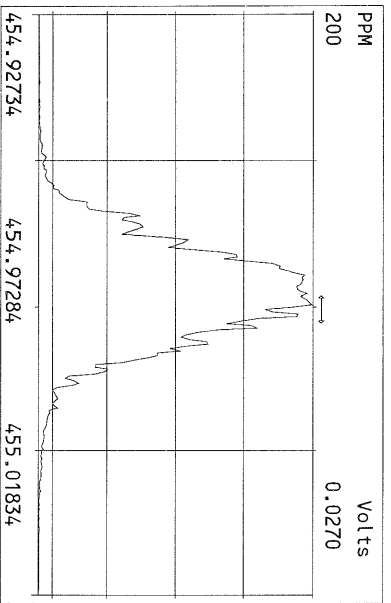
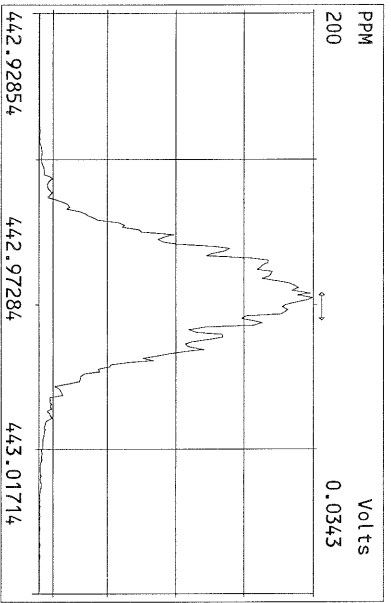
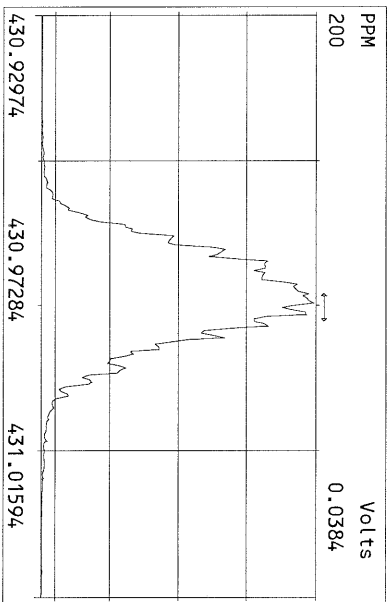
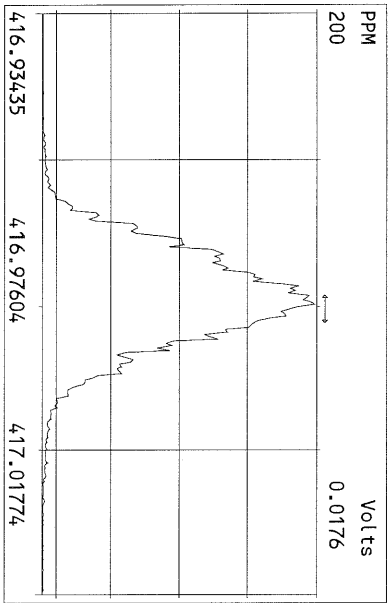
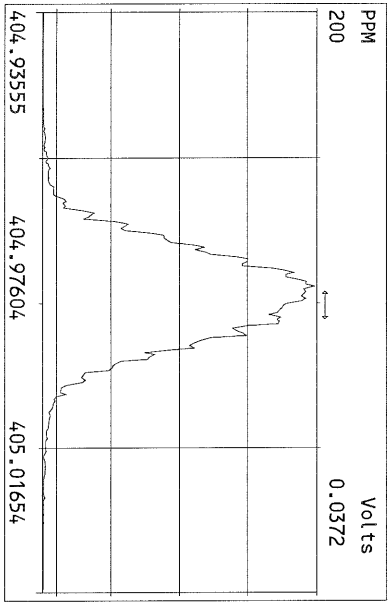
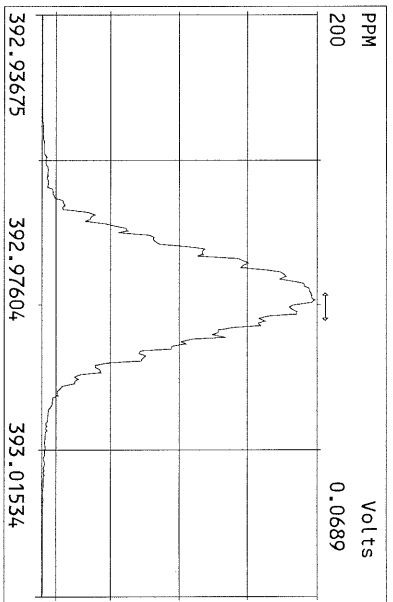
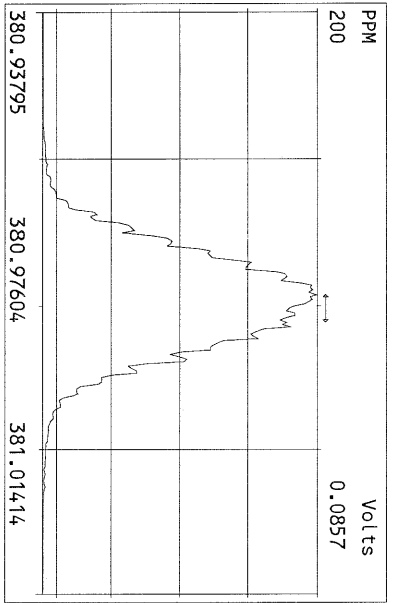
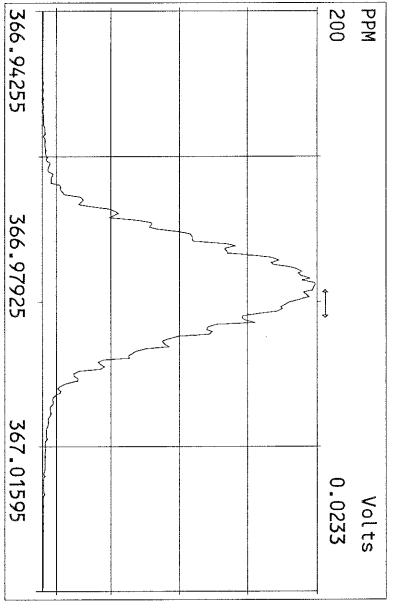
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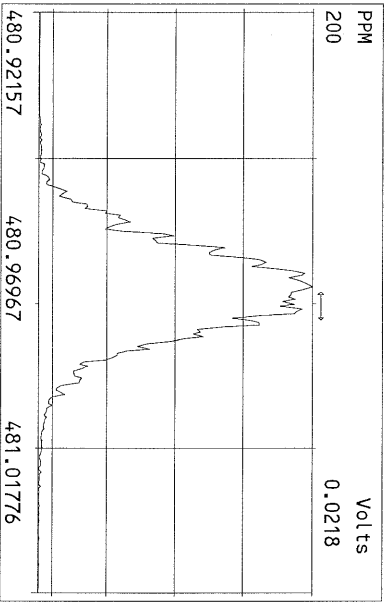
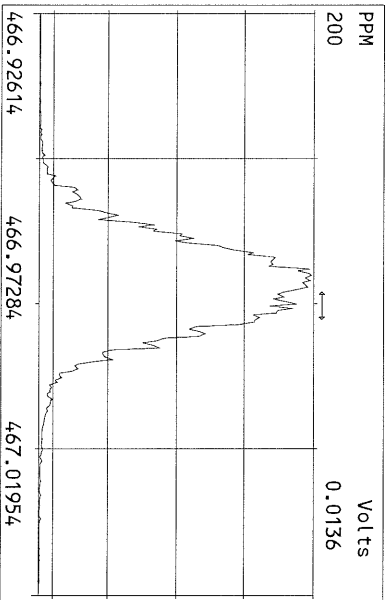
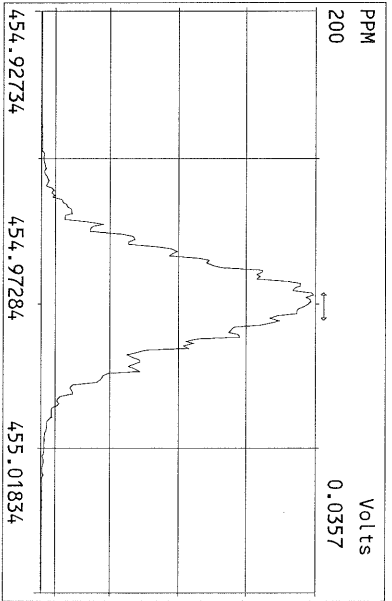
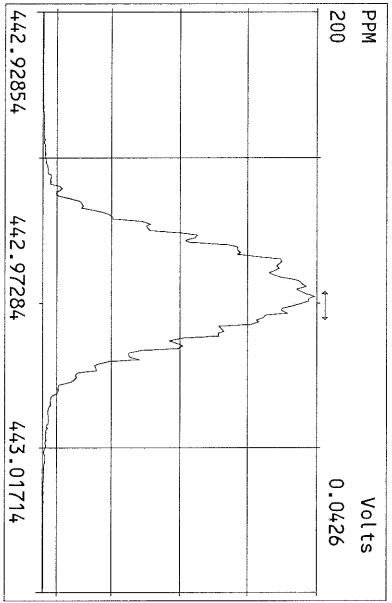
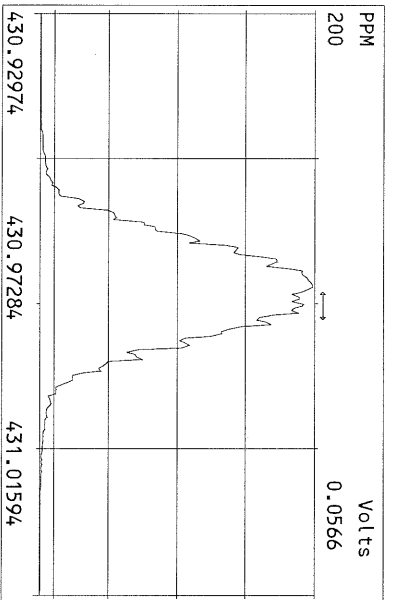
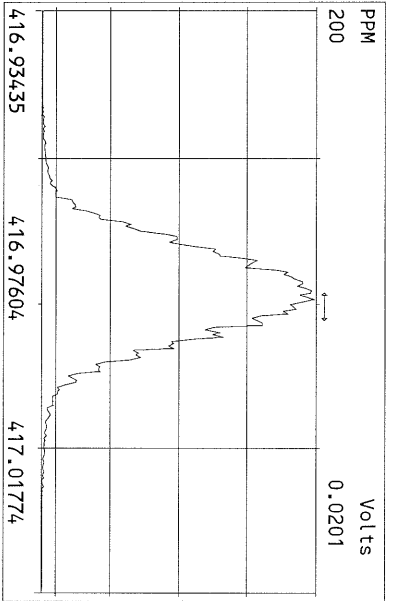
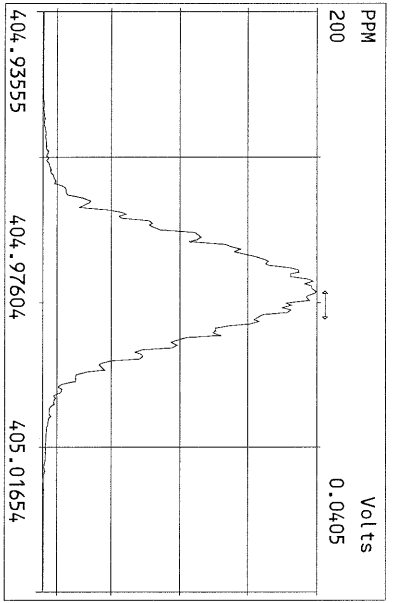
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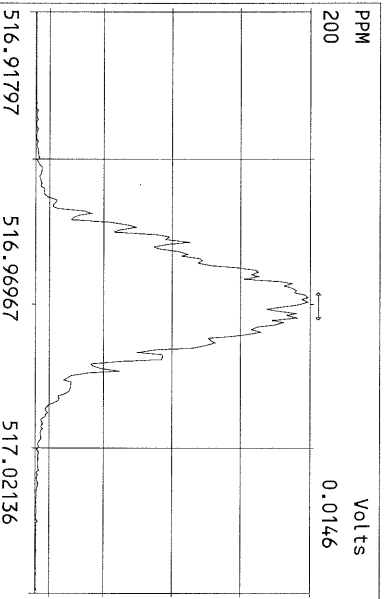
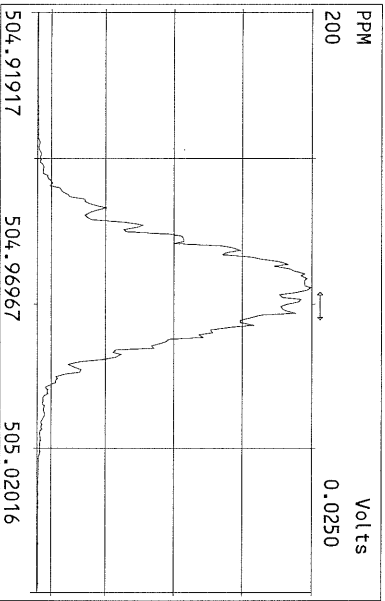
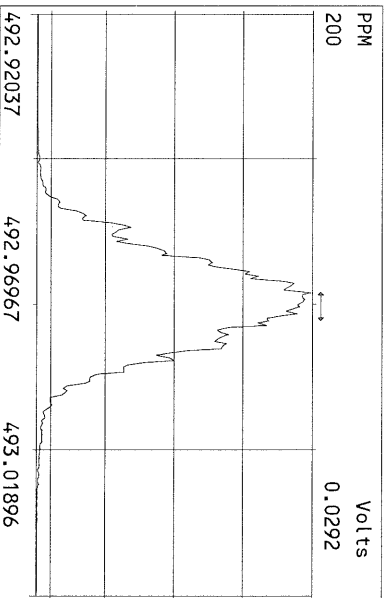
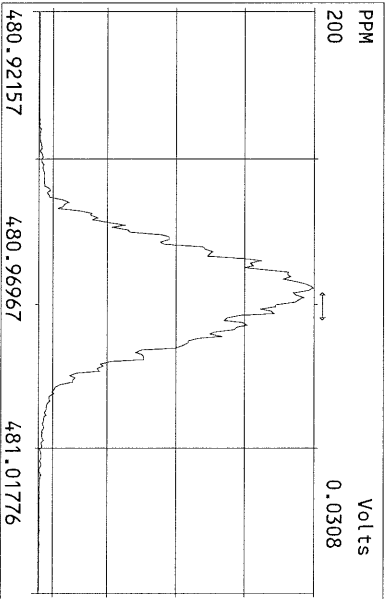
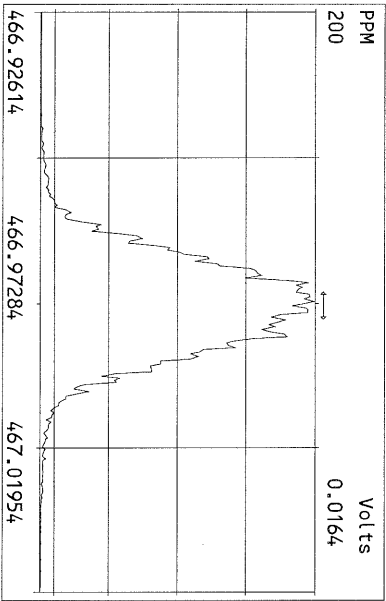
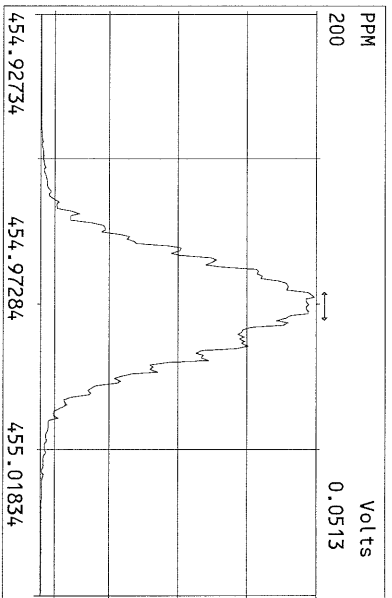
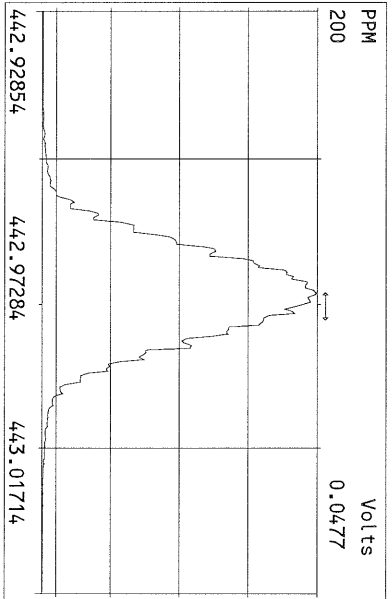
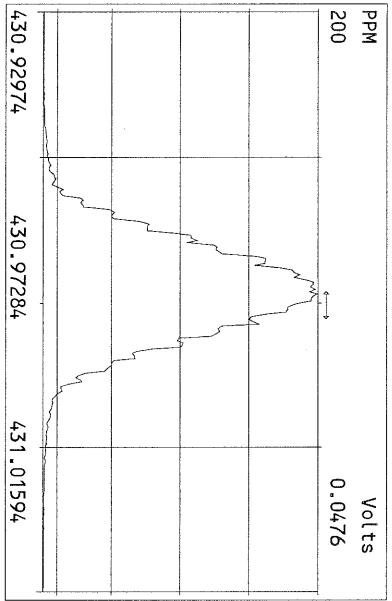
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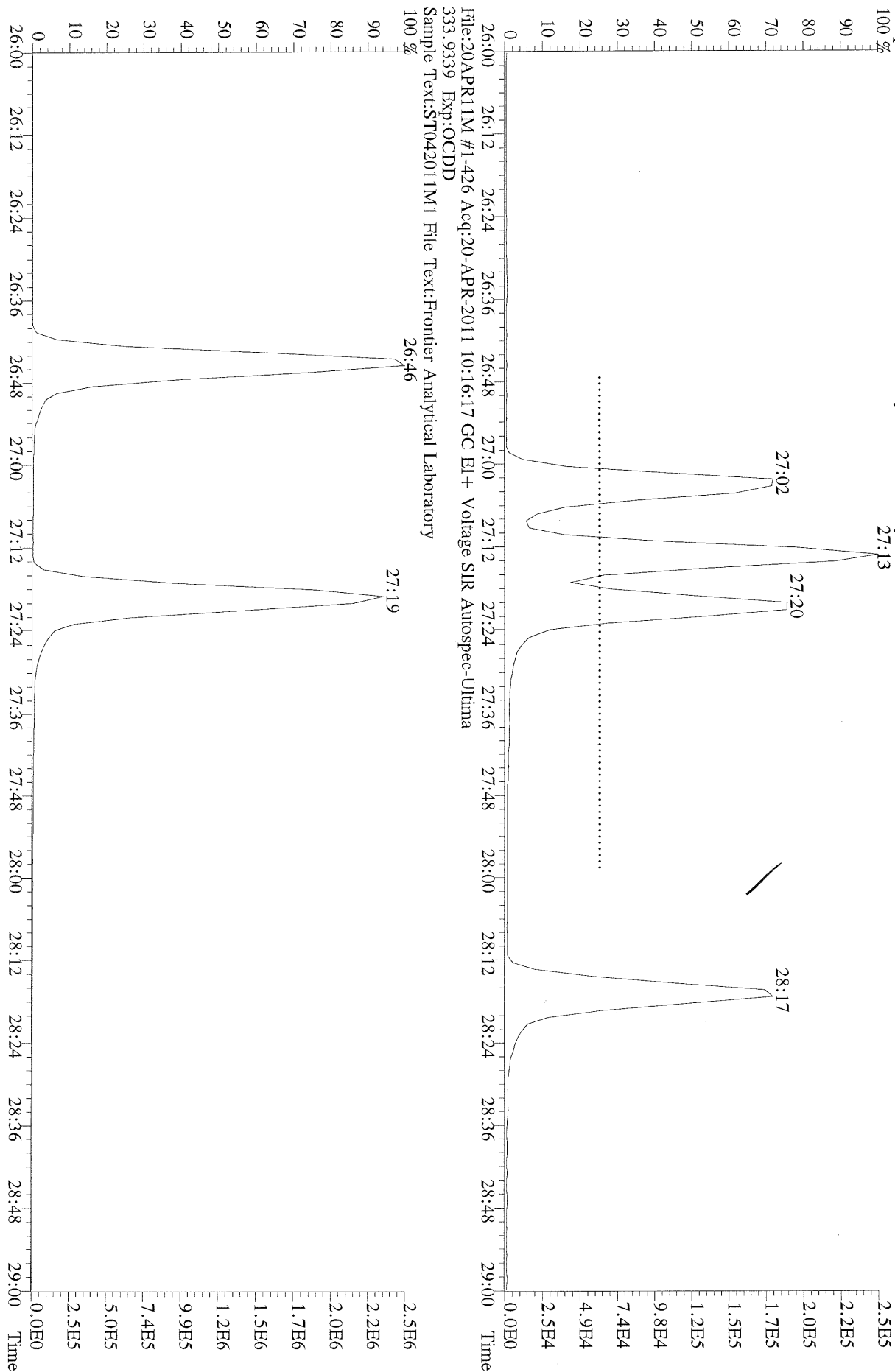




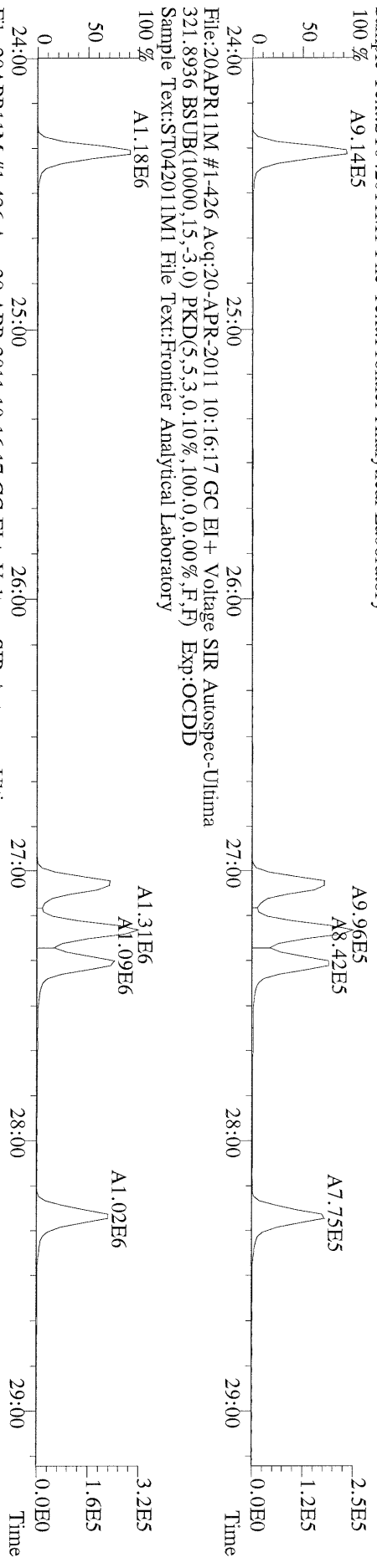




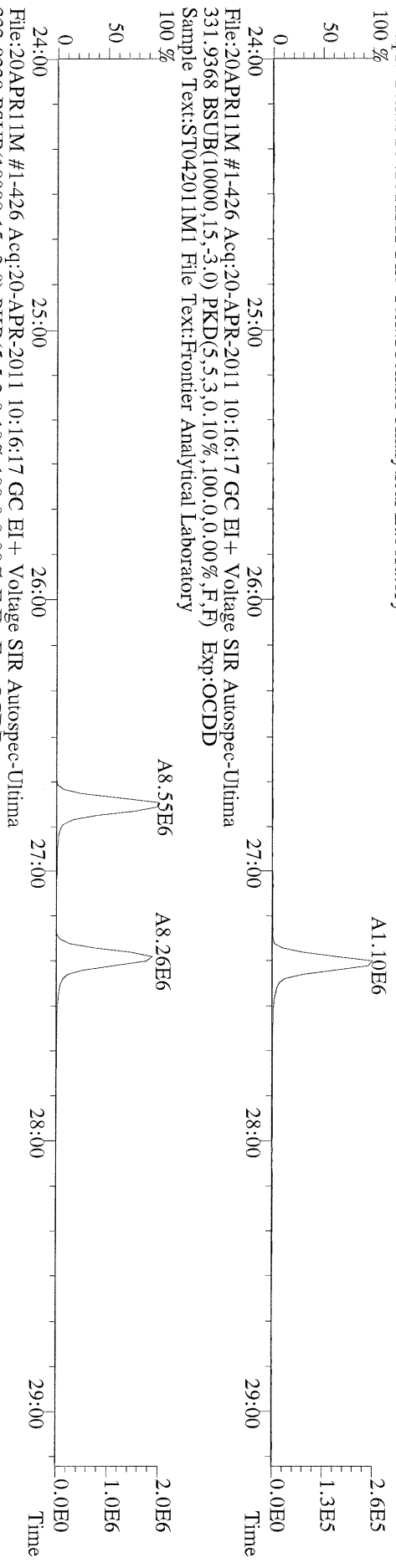
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319.8965 Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



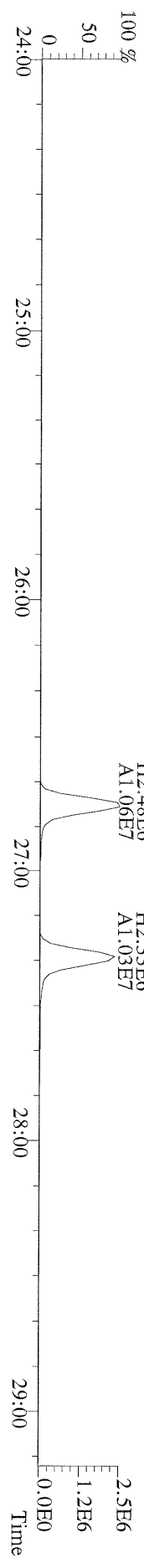
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319.8965 BSUB(10000,15,-3,0) PKD(5.5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



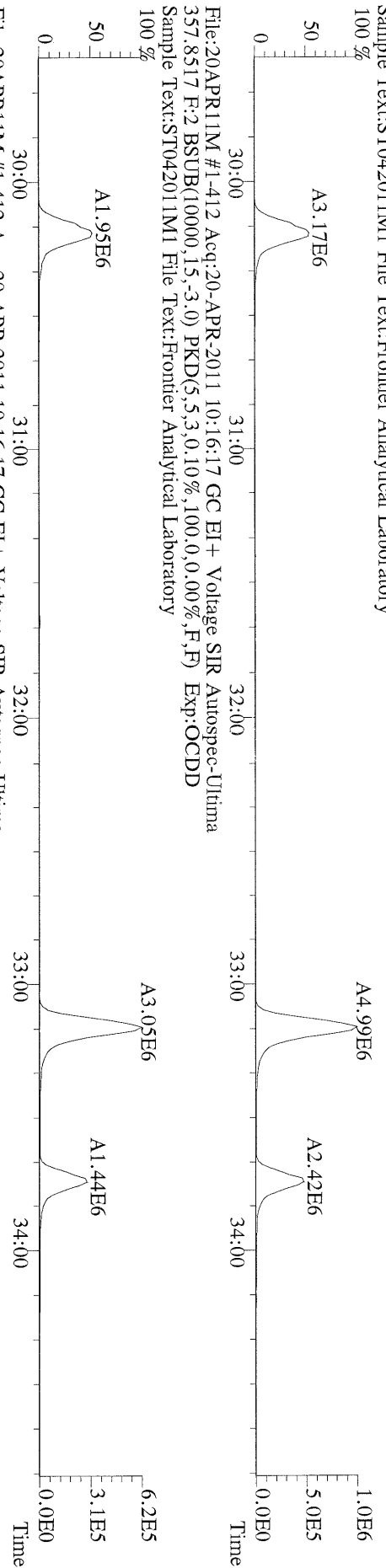
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327.8847 BSUB(10000,15,-3,0) PKD(5.5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



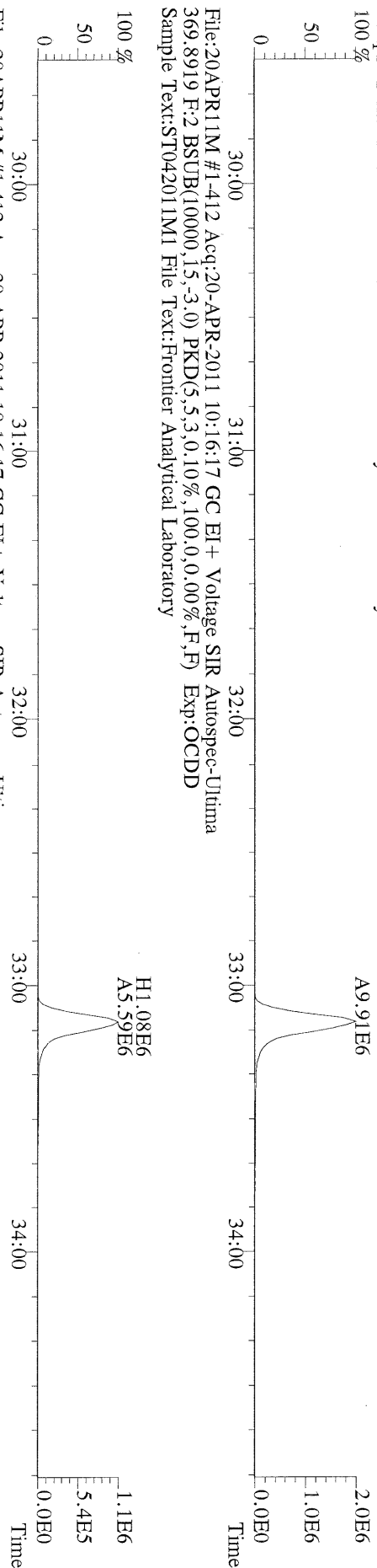
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333.9339 BSUB(10000,15,-3,0) PKD(5.5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



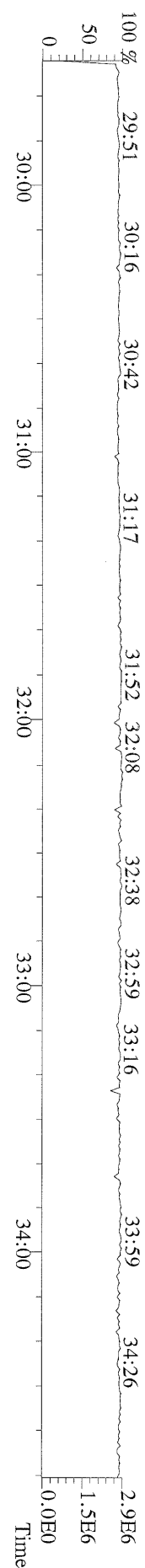
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 Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



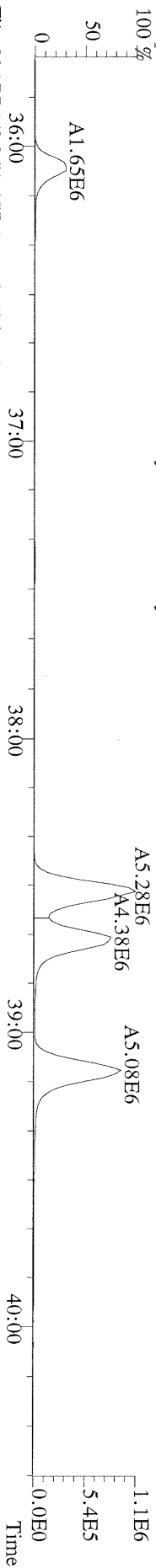
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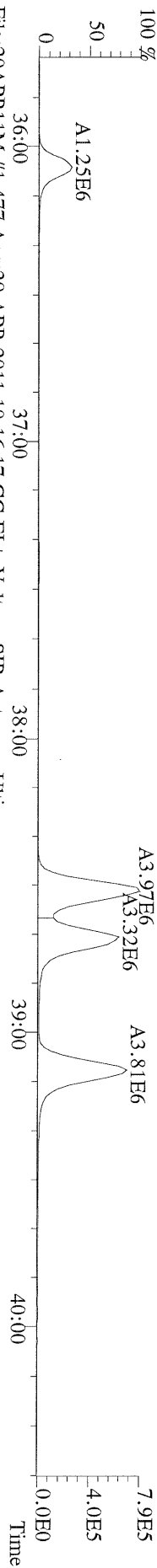
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 366.9792 F:2 Exp:OCDD
 Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



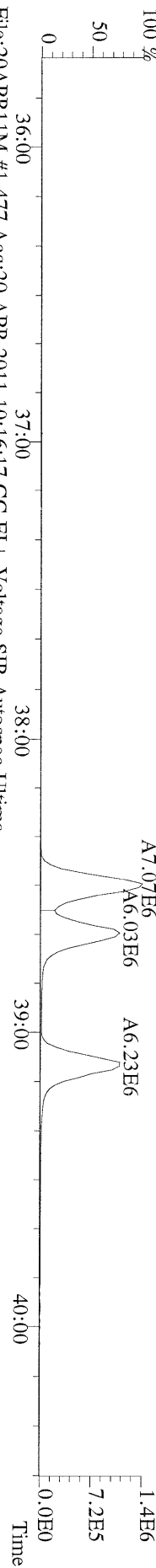
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389.8156 F:3 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



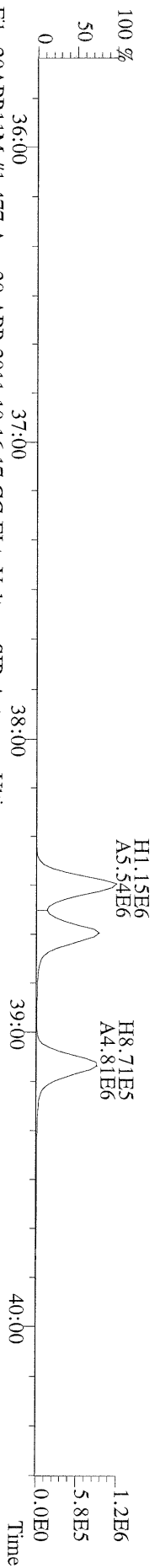
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391.8127 F:3 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
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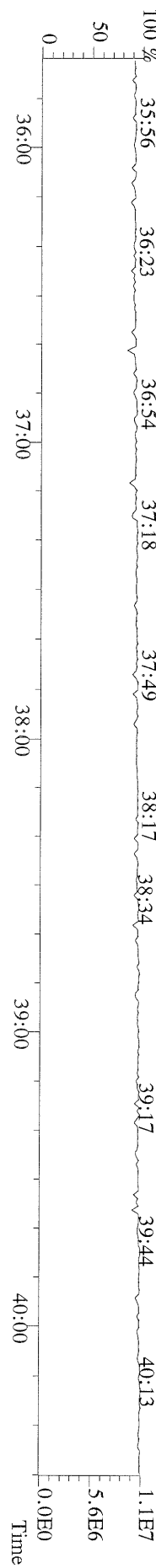
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



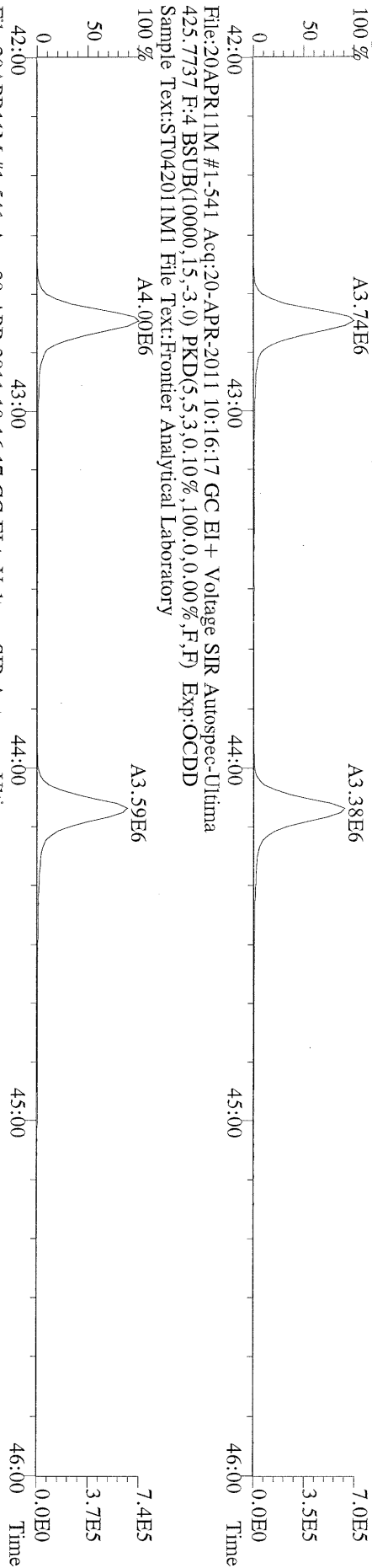
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



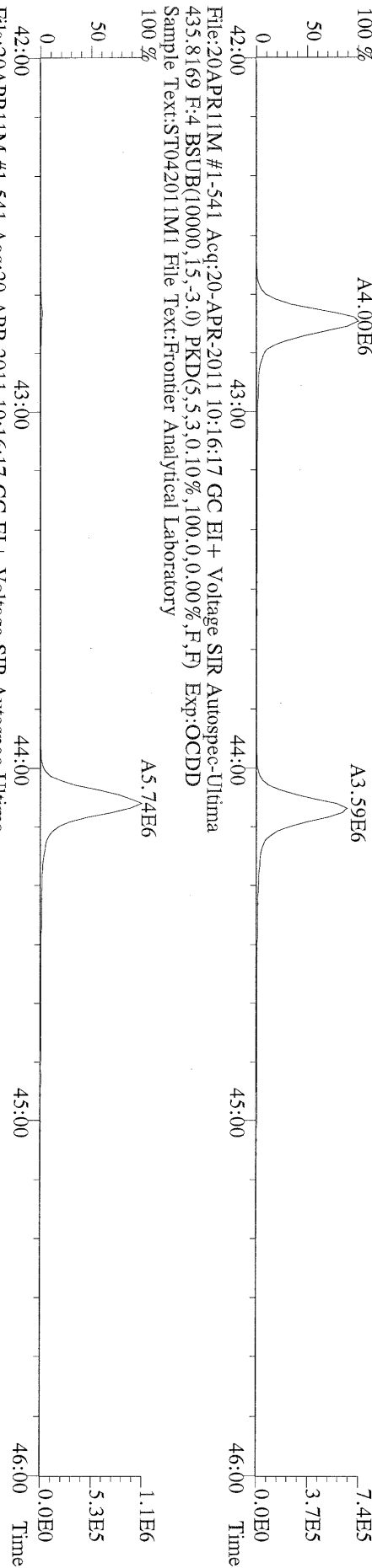
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



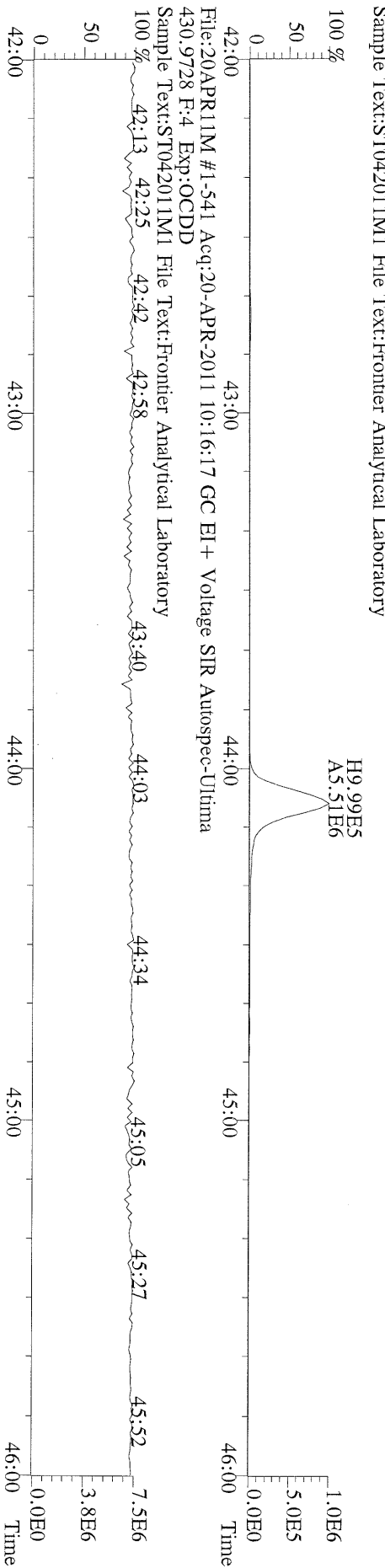
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423.7767 F:4 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



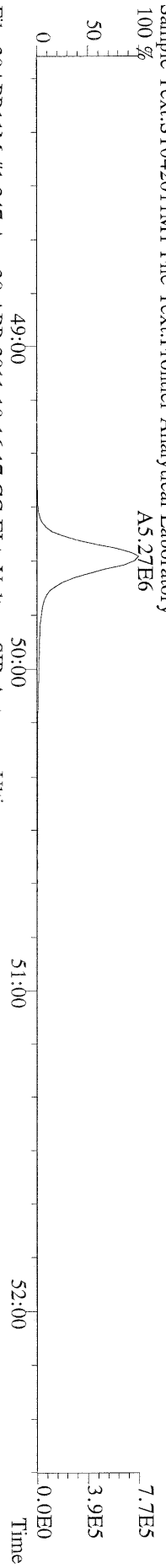
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425.7737 F:4 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



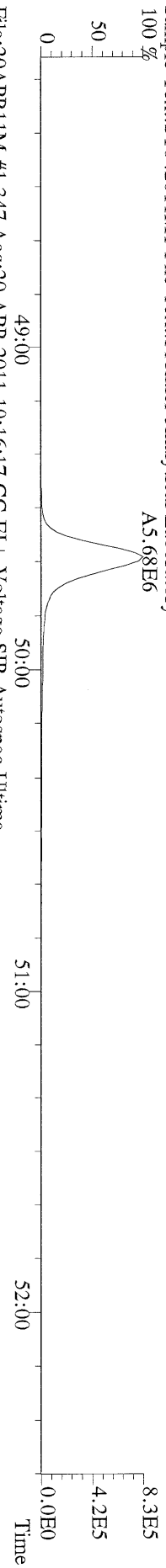
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437.8140 F:4 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



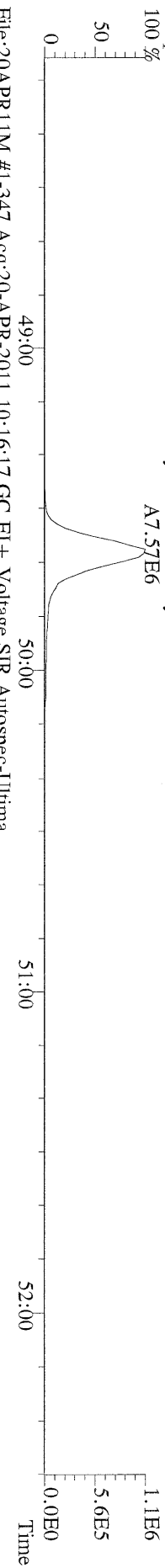
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457.7377 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



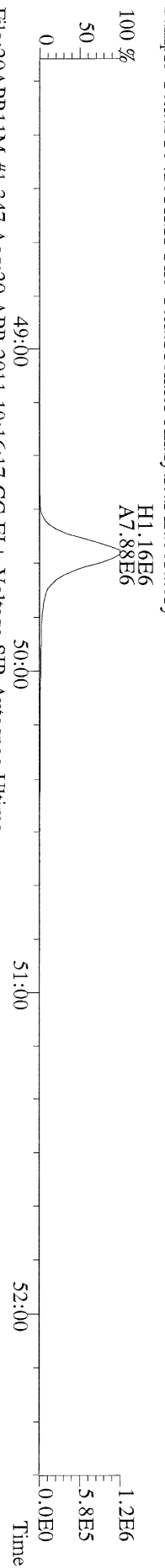
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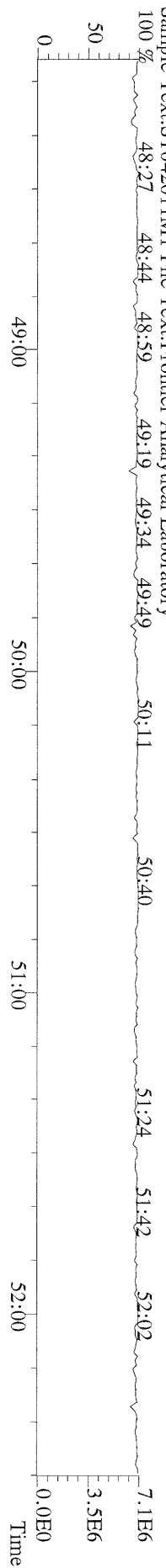
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



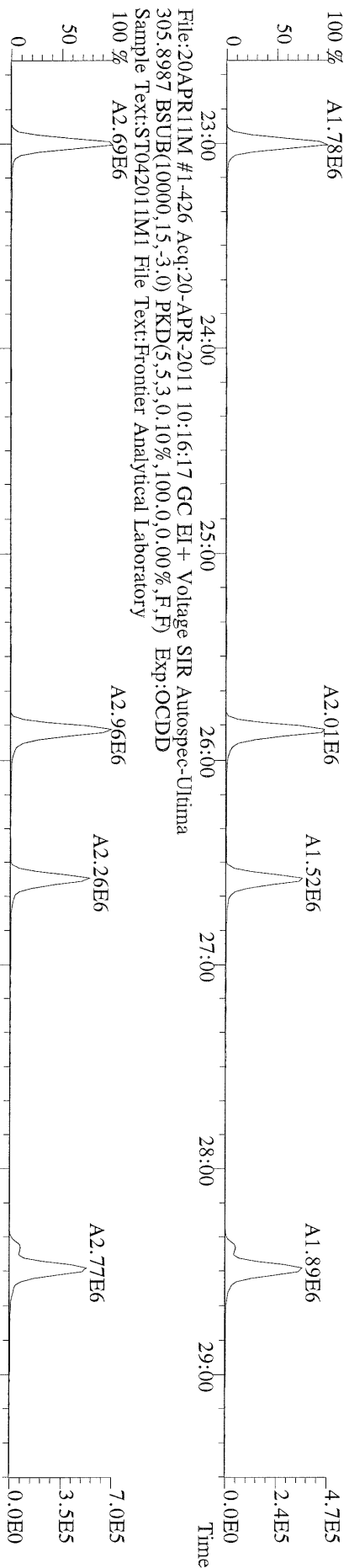
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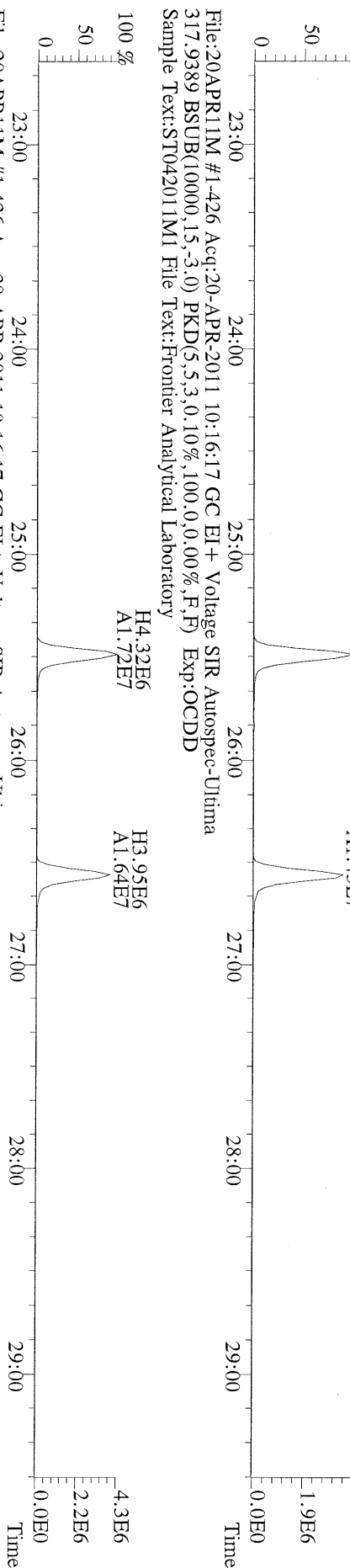
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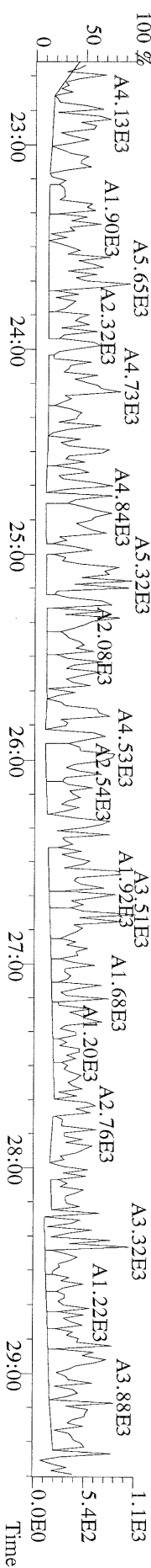
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303.9016 BSUB(10000,15,-3.0) PKD(5.5,3.0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



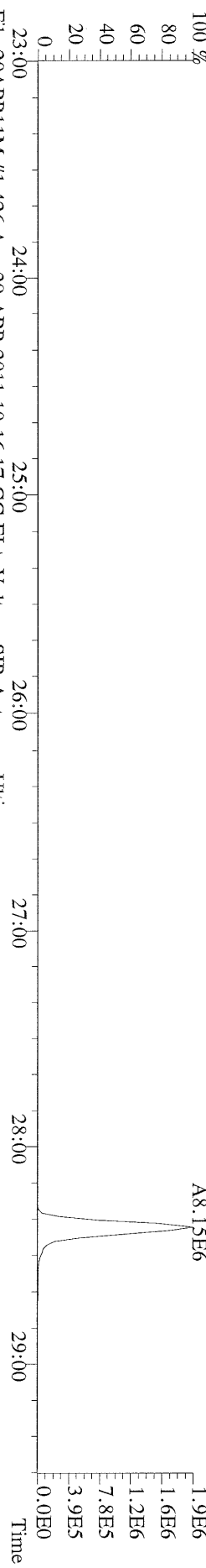
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



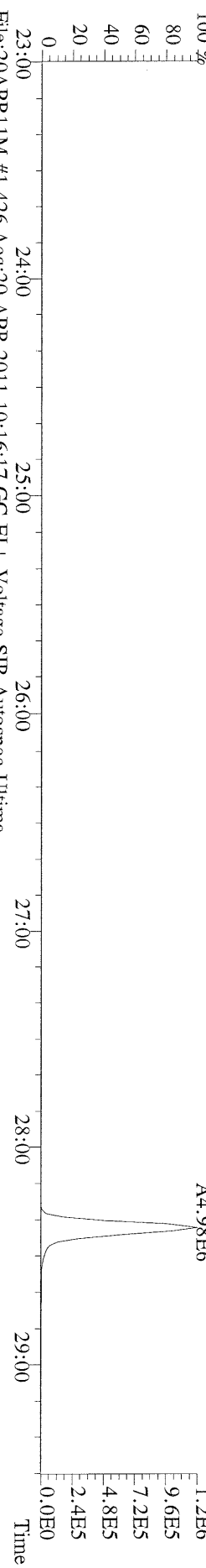
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375.8364 BSUB(10000,15,-3.0) PKD(5.5,3.0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



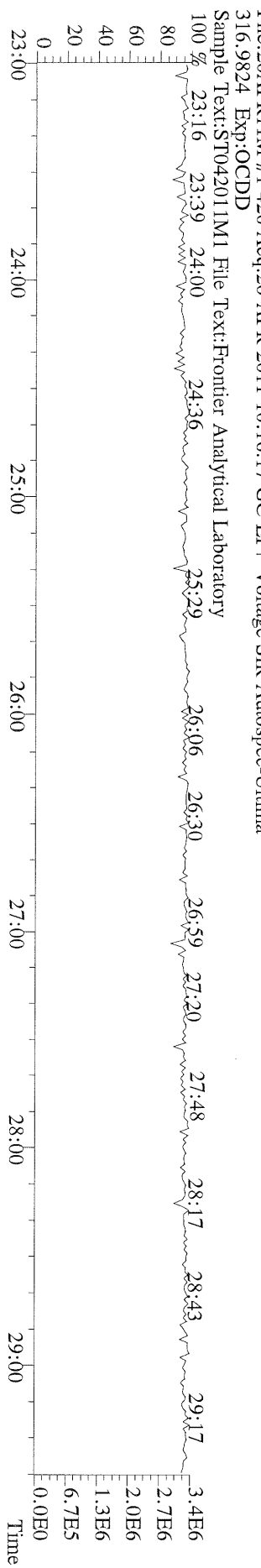
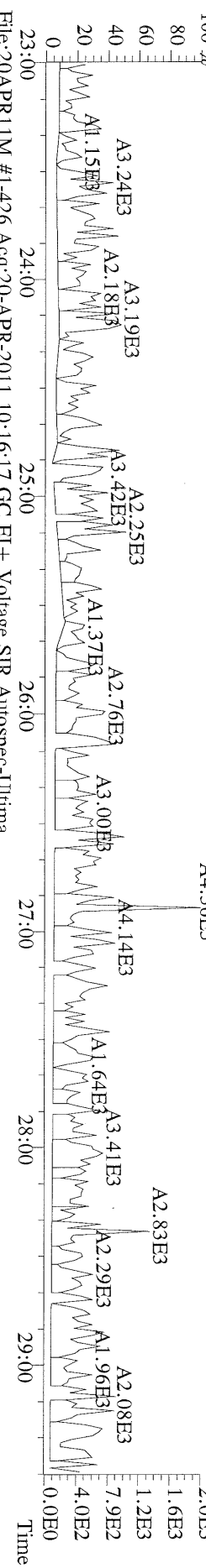
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 339.8597 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



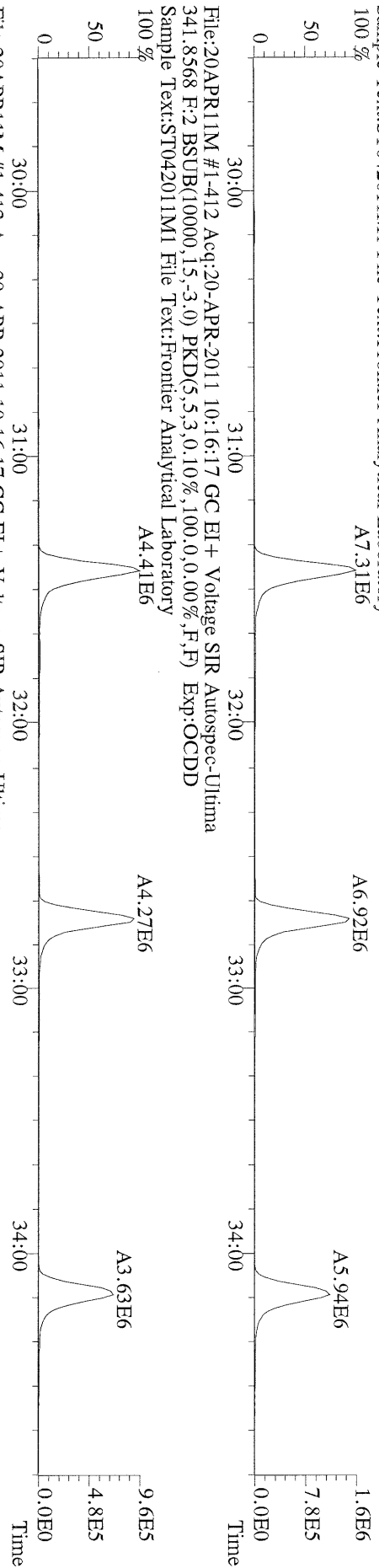
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 Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



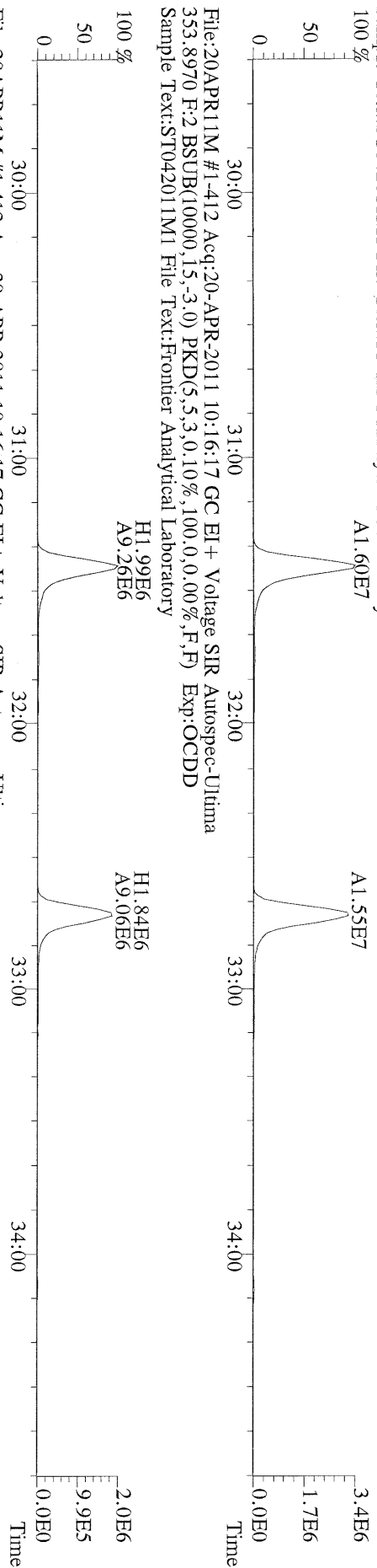
File:20APR11M #1-426 Acq:20-APR-2011 10:16:17 GC EI+ Voltage SIR Autospec-Ultima
 409.7974 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



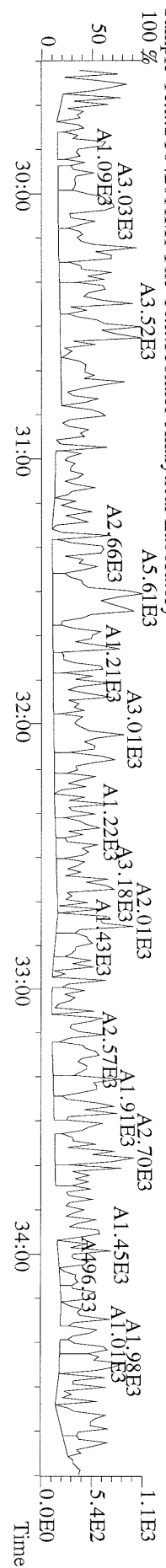
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339.8597 F:2 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



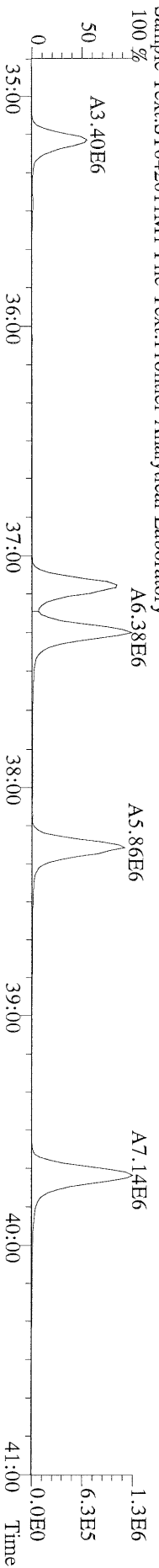
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351.9000 F:2 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



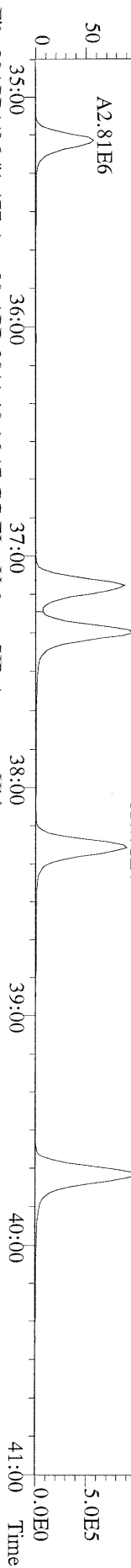
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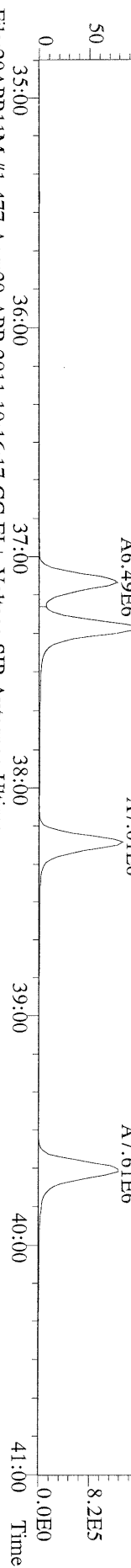
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373.8207 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



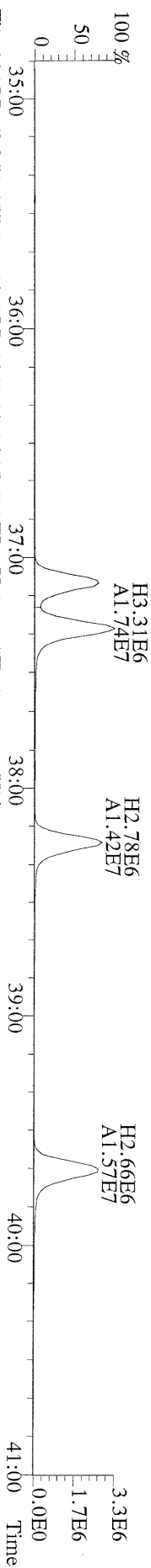
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



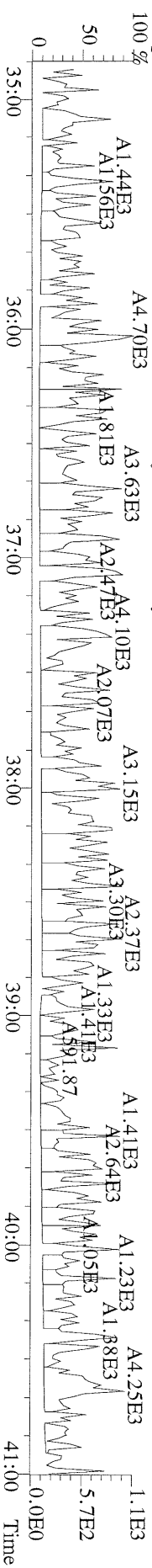
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383.8639 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



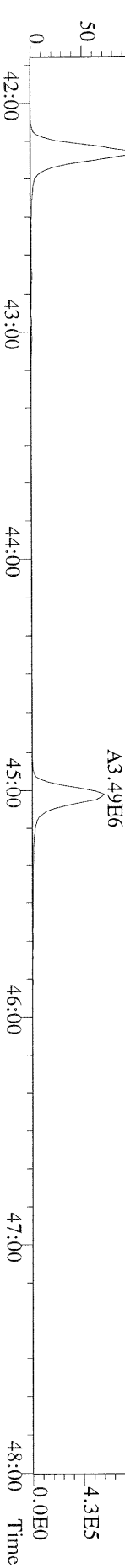
File:20APR11M #1-477 Acq:20-APR-2011 10:16:17 GC EI+ Voltage SIR Autospec-Utima
385.8610 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



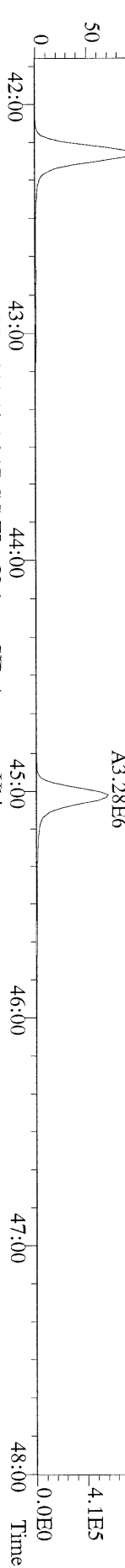
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445.7555 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory



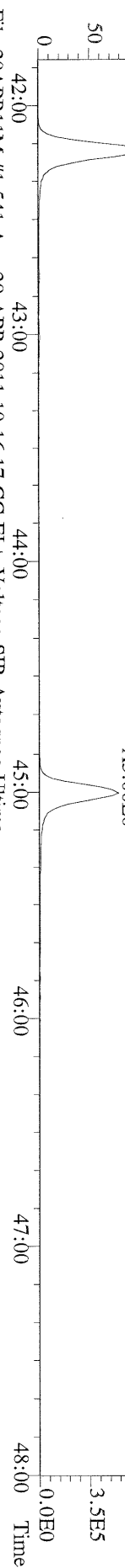
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407.7818 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



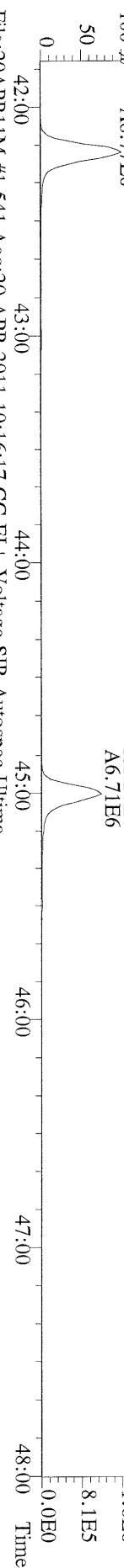
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409.7788 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



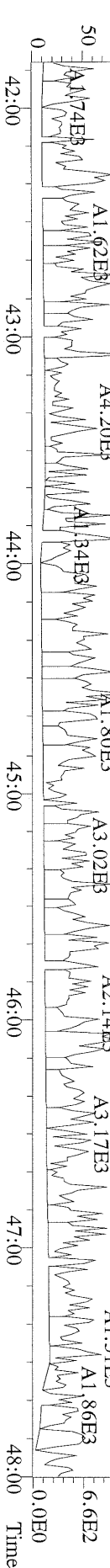
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417.8253 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



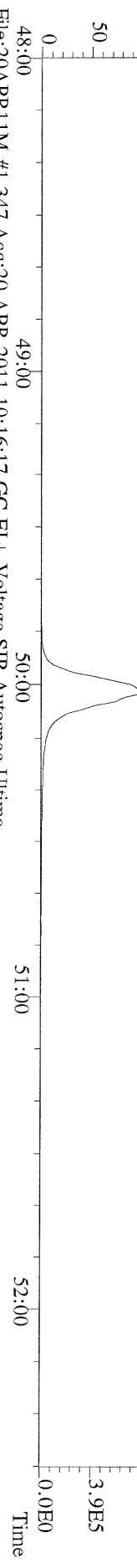
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419.8220 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



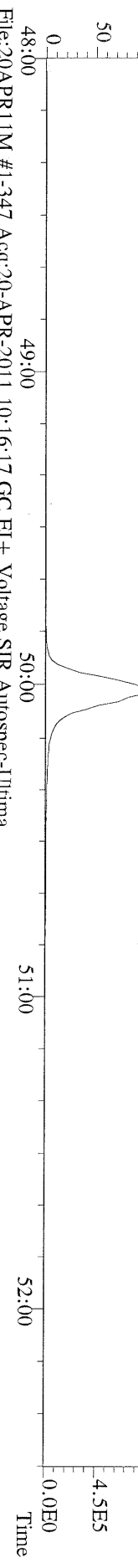
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Sample Text:ST042011M1 File Text:Frontier Analytical Laboratory
100 %



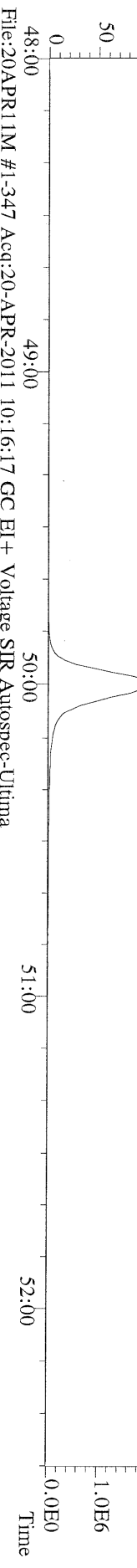
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441.7428 F:5 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM1 File Text:Frontier Analytical Laboratory



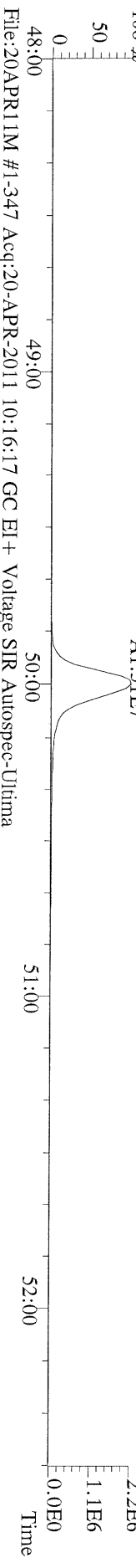
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Sample Text:ST04201IM1 File Text:Frontier Analytical Laboratory



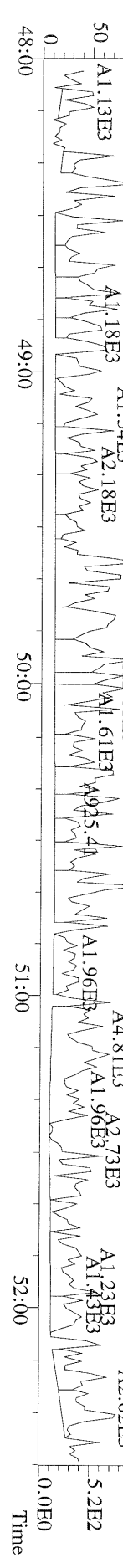
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453.7831 F:5 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM1 File Text:Frontier Analytical Laboratory

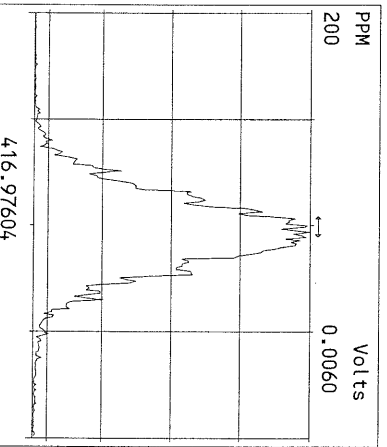
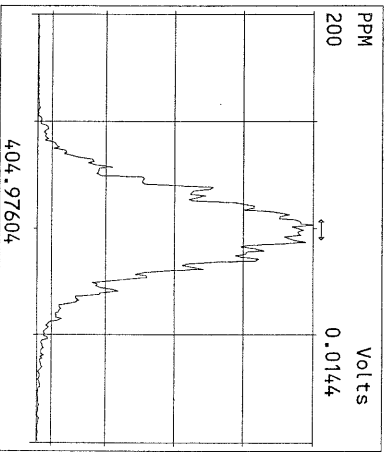
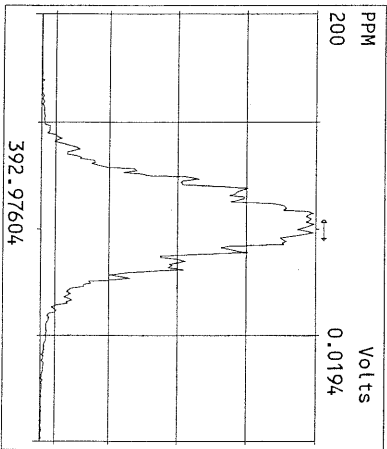
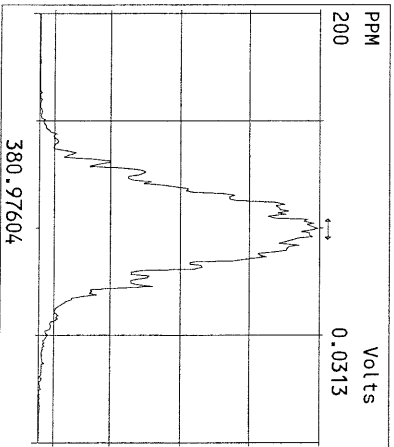
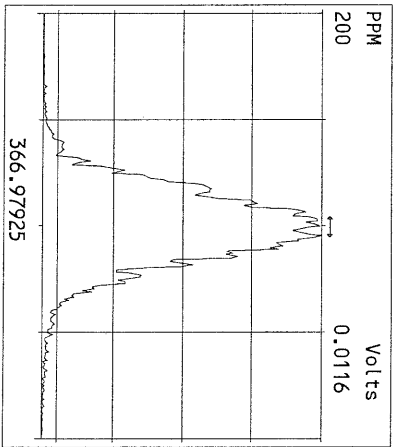
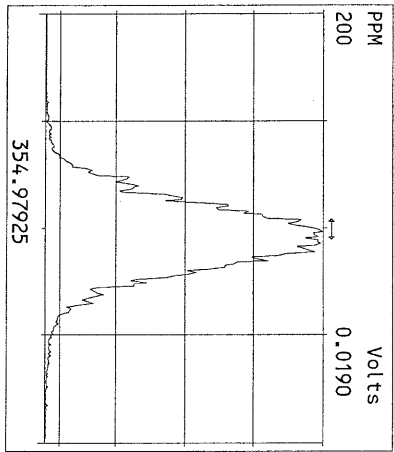
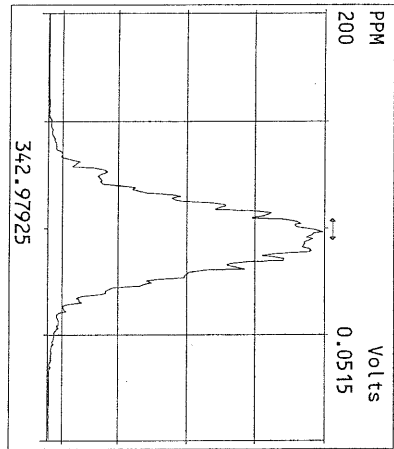
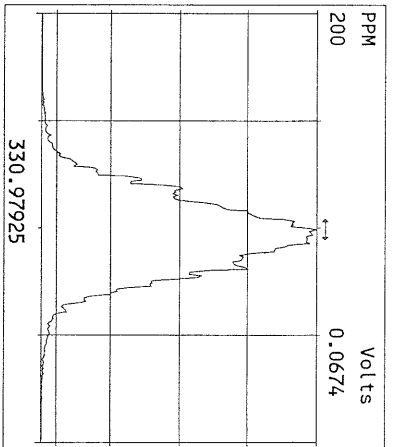
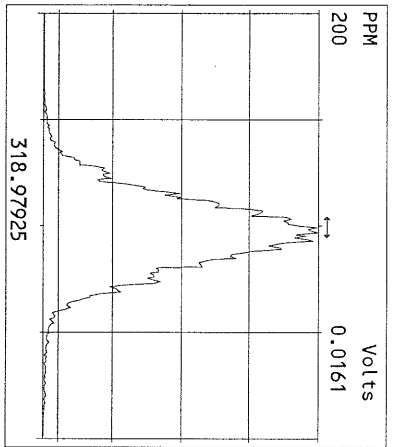
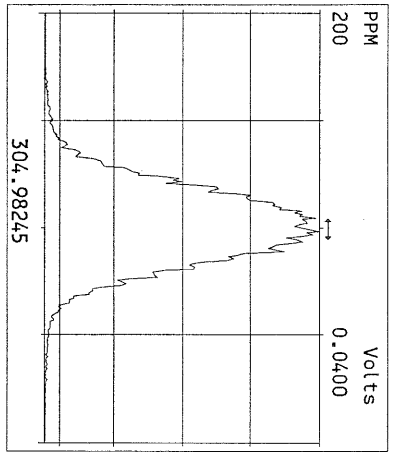
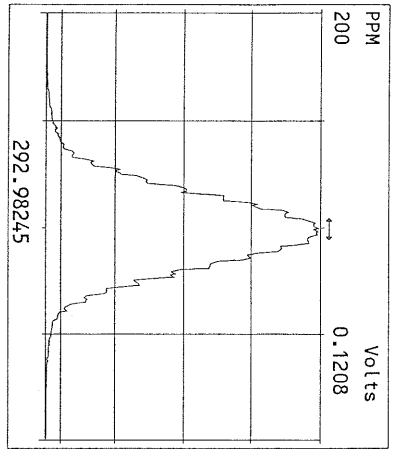


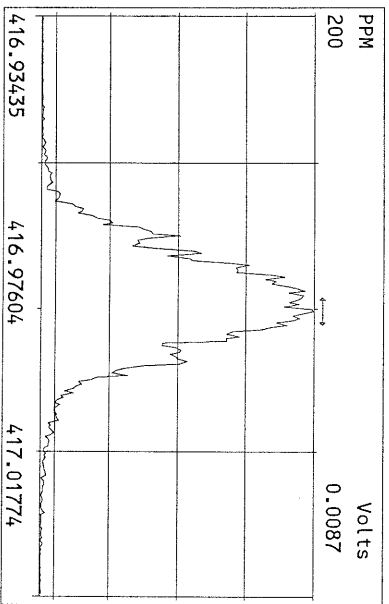
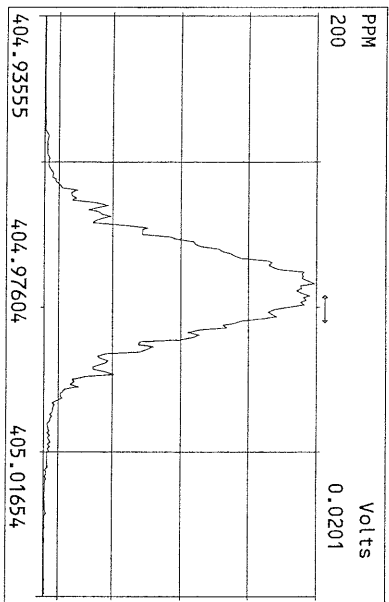
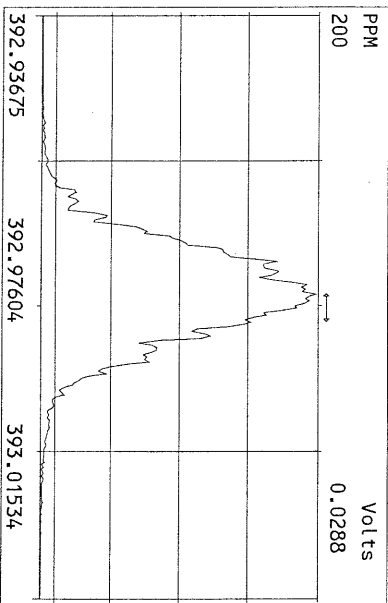
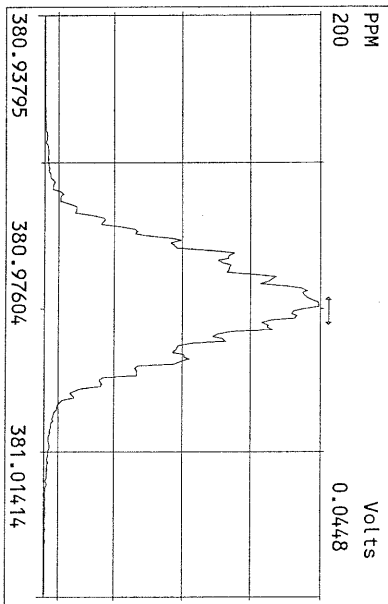
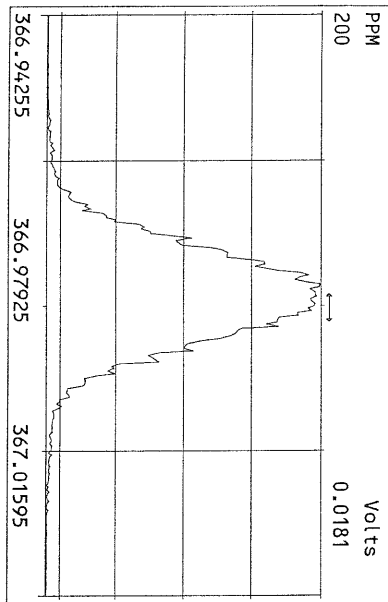
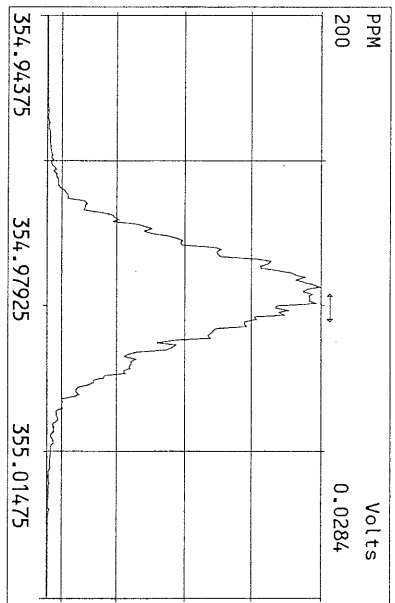
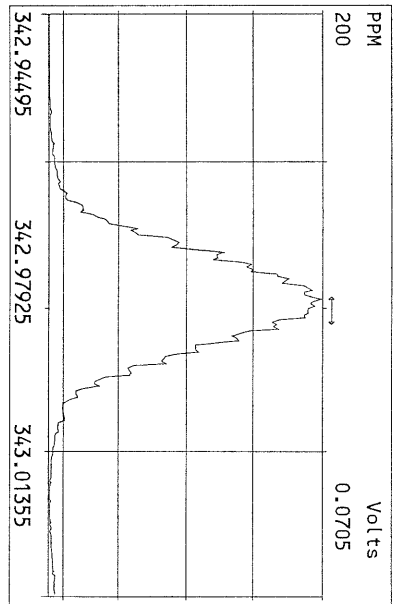
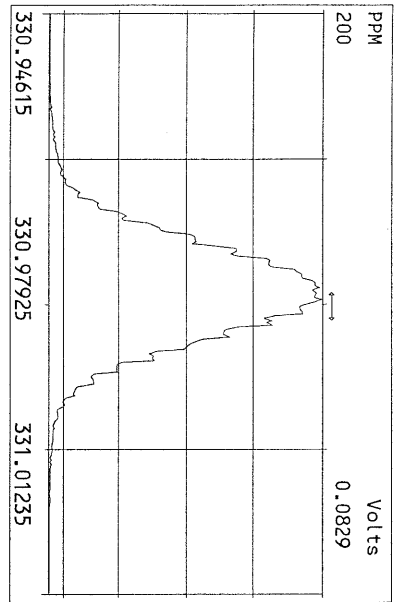
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455.7801 F:5 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM1 File Text:Frontier Analytical Laboratory

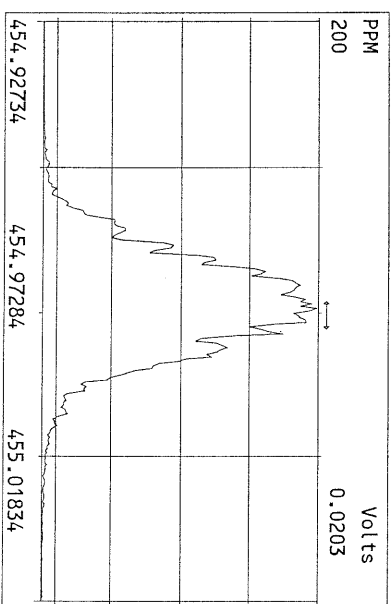
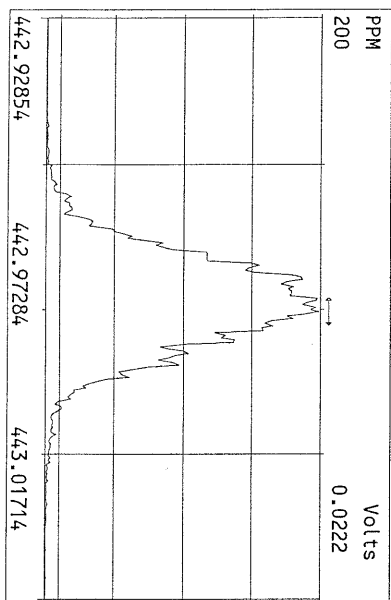
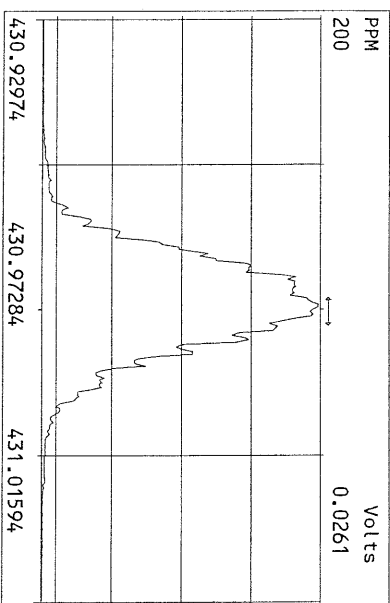
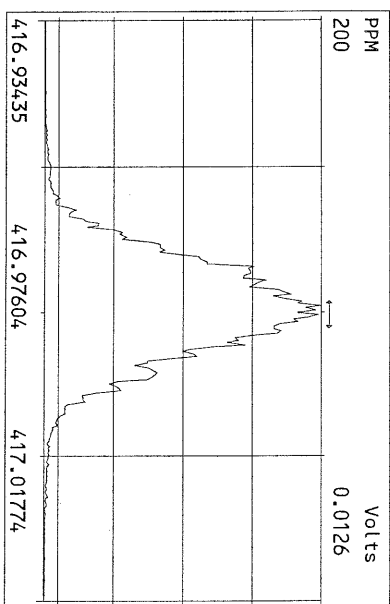
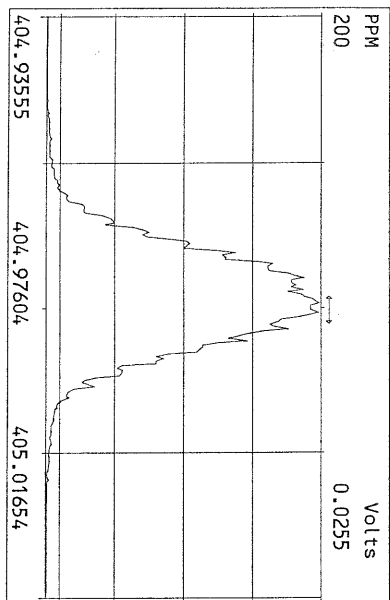
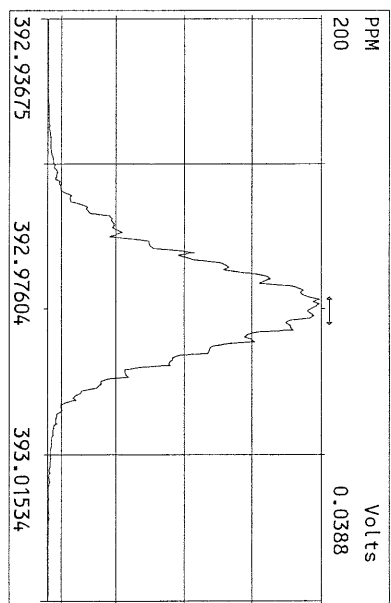
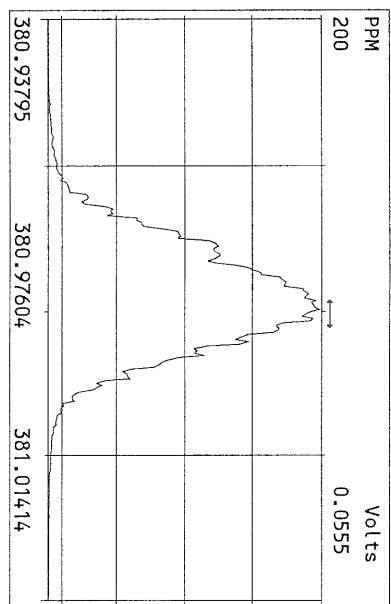
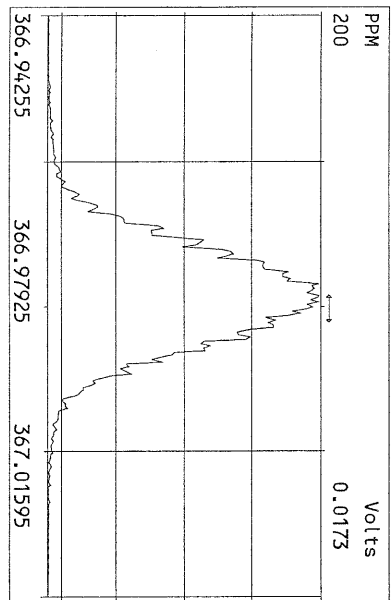


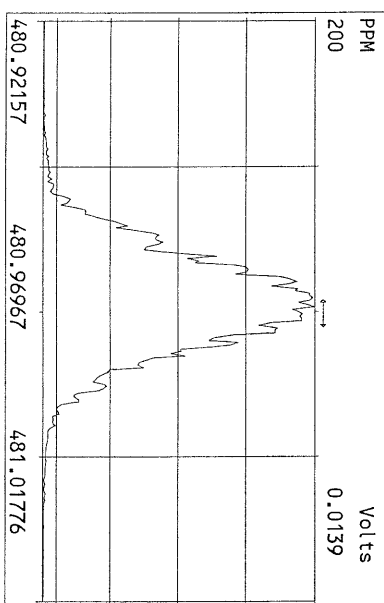
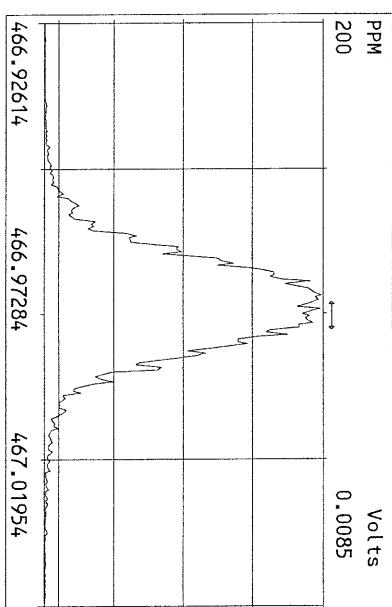
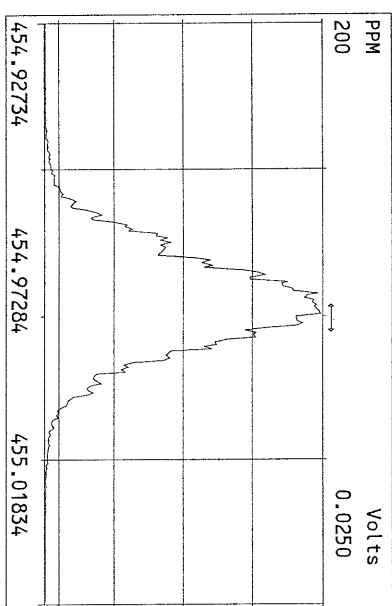
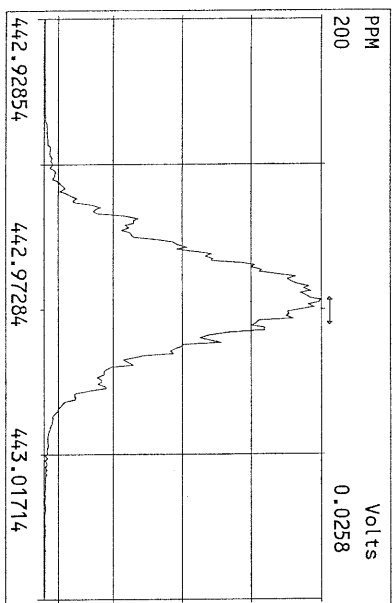
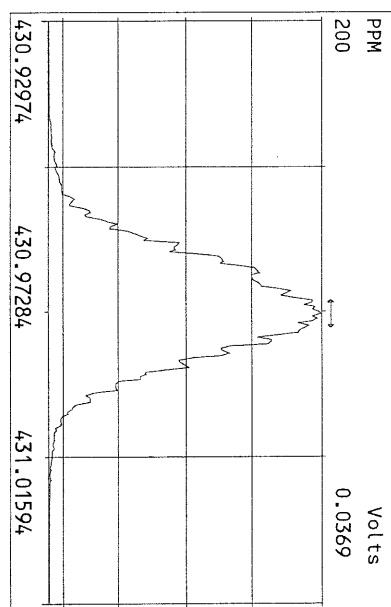
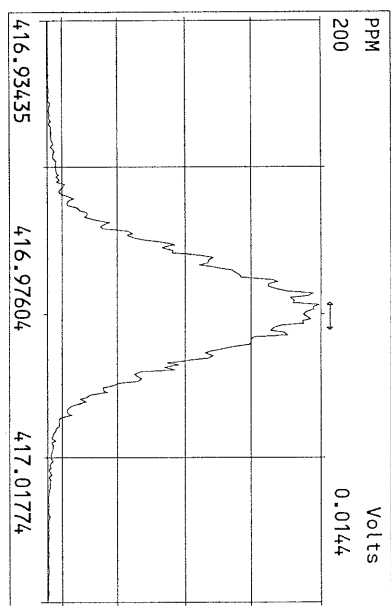
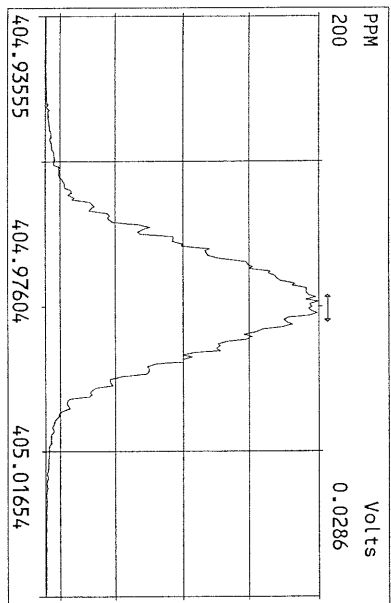
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513.6775 F:5 BSUB(10000,15,-3.0) PKD(5,5.3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM1 File Text:Frontier Analytical Laboratory

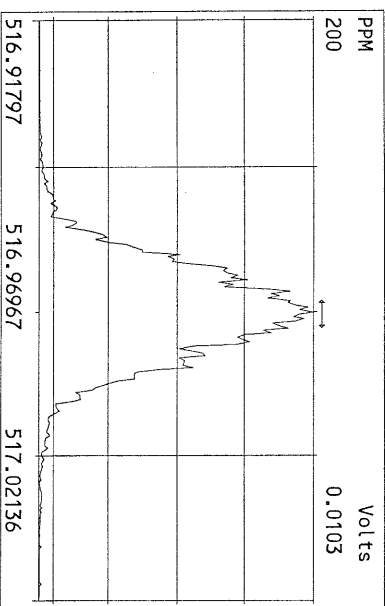
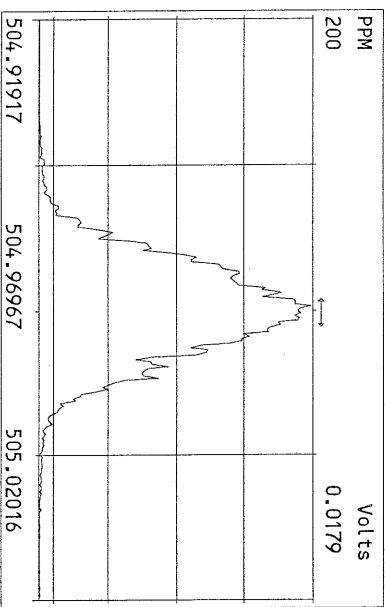
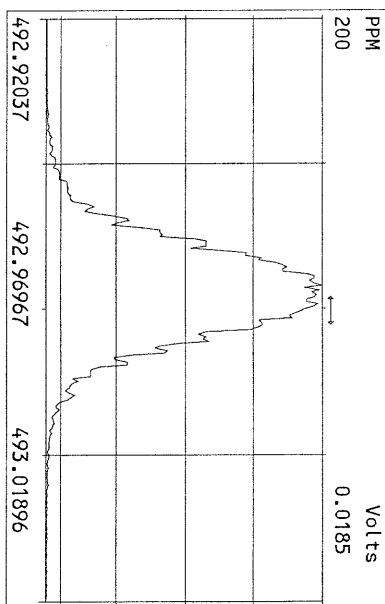
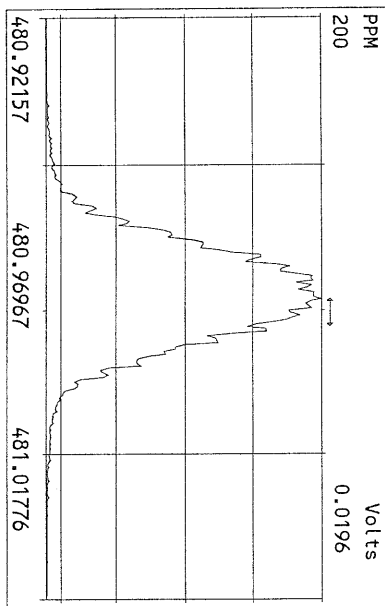
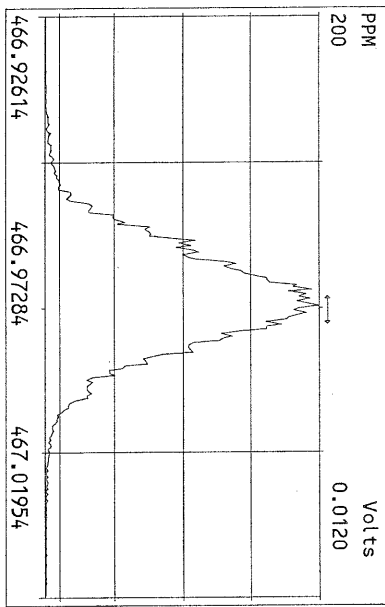
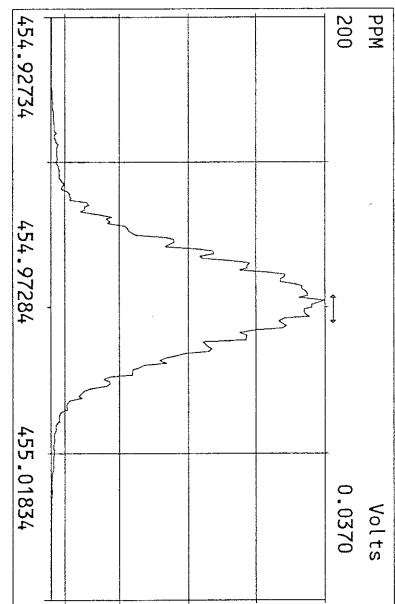
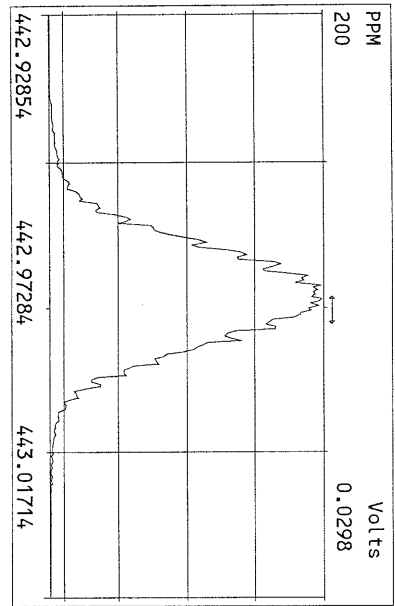
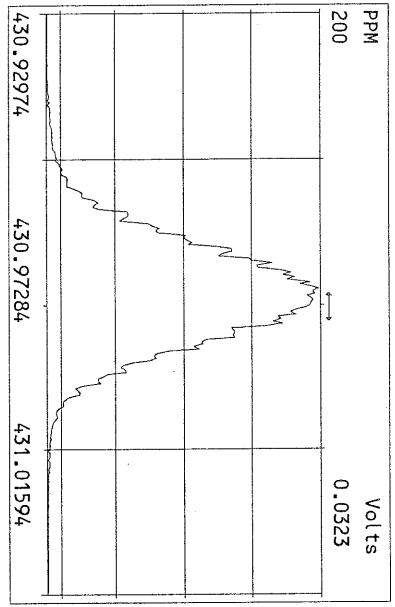












USEPA - ITD

FORM 4A
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3 GC Column ID: DB5

VER Data Filename: 20APR11M Sam:17 Analysis Date: 21-APR-11 01:02:13

NATIVE ANALYTES	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
2,3,7,8-TCDD	M/M+2	0.77	0.65-0.89	y	8.99	7.80 - 12.9 ✓
1,2,3,7,8-PeCDD	M+2/M+4	1.62	1.32-1.78	y	53.0	39.0 - 65.0 ✓
1,2,3,4,7,8-HxCDD	M+2/M+4	1.33	1.05-1.43	y	51.3	39.0 - 64.0 ✓
1,2,3,6,7,8-HxCDD	M+2/M+4	1.32	1.05-1.43	y	50.9	39.0 - 64.0 ✓
1,2,3,7,8,9-HxCDD	M+2/M+4	1.32	1.05-1.43	y	52.9	41.0 - 61.0 ✓
1,2,3,4,6,7,8-HpCDD	M+2/M+4	0.96	0.88-1.20	y	48.9	43.0 - 58.0 ✓
OCDD	M+2/M+4	0.93	0.76-1.02	y	98.2	79.0 - 126 ✓
2,3,7,8-TCDF	M/M+2	0.66	0.65-0.89	y	10.7	8.40 - 12.0 ✓
1,2,3,7,8-PeCDF	M+2/M+4	1.68	1.32-1.78	y	53.8	41.0 - 60.0 ✓
2,3,4,7,8-PeCDF	M+2/M+4	1.65	1.32-1.78	y	52.1	41.0 - 60.0 ✓
1,2,3,4,7,8-HxCDF	M+2/M+4	1.27	1.05-1.43	y	49.5	45.0 - 56.0 ✓
1,2,3,6,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	50.2	44.0 - 57.0 ✓
2,3,4,6,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	49.3	44.0 - 57.0 ✓
1,2,3,7,8,9-HxCDF	M+2/M+4	1.27	1.05-1.43	y	50.3	45.0 - 56.0 ✓
1,2,3,4,6,7,8-HpCDF	M+2/M+4	1.04	0.88-1.20	y	48.9	45.0 - 55.0 ✓
1,2,3,4,7,8,9-HpCDF	M+2/M+4	1.05	0.88-1.20	y	48.9	43.0 - 58.0 ✓
OCDF	M+2/M+4	0.89	0.76-1.02	y	96.4	63.0 - 159 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

Analyst: Date: 4/21/11

USEPA - ITD

FORM 4B
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 20APR11M Sam:17

Analysis Date: 21-APR-11 01:02:13

LABELLED COMPOUNDS	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
13C-2,3,7,8-TCDD	M/M+2	0.82	0.65-0.89	y	95.8	82.0 - 121 ✓
13C-1,2,3,7,8-PeCDD	M+2/M+4	1.76	1.32-1.78	y	86.0	62.0 - 160 ✓
13C-1,2,3,4,7,8-HxCDD	M+2/M+4	1.29	1.05-1.43	y	94.6	85.0 - 117 ✓
13C-1,2,3,6,7,8-HxCDD	M+2/M+4	1.27	1.05-1.43	y	98.4	85.0 - 118 ✓
13C-1,2,3,4,6,7,8-HpCDD	M+2/M+4	1.06	0.88-1.20	y	122	72.0 - 138 ✓
13C-OCDD	M+2/M+4	0.97	0.76-1.02	y	226	96.0 - 415 ✓
13C-2,3,7,8-TCDF	M/M+2	0.87	0.65-0.89	y	102	71.0 - 140 ✓
13C-1,2,3,7,8-PeCDF	M+2/M+4	1.73	1.32-1.78	y	99.6	76.0 - 130 ✓
13C-2,3,4,7,8-PeCDF	M+2/M+4	1.71	1.32-1.78	y	101	77.0 - 130 ✓
13C-1,2,3,4,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	102	76.0 - 131 ✓
13C-1,2,3,6,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	102	70.0 - 143 ✓
13C-2,3,4,6,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	104	73.0 - 137 ✓
13C-1,2,3,7,8,9-HxCDF	M/M+2	0.49	0.43-0.59	y	109	74.0 - 135 ✓
13C-1,2,3,4,6,7,8-HpCDF	M/M+2	0.45	0.37-0.51	y	118	78.0 - 129 ✓
13C-1,2,3,4,7,8,9-HpCDF	M/M+2	0.44	0.37-0.51	y	128	77.0 - 129 ✓
13C-OCDF	M+2/M+4	0.95	0.76-1.02	y	224	96.0 - 415 ✓
CLEANUP STANDARD (4)						
37Cl-2,3,7,8-TCDD					7.97	7.80 - 12.8 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

(4) No ion abundance ratio; report concentration found.

Analyst: _____

Date: 4/21/11

FORM 5
 PCDD/PCDF RT WINDOW AND ISOMER SPECIFICITY STANDARDS

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.:

Instrument ID: FAL3 Initial Calibration Date: 3/7/11

RT Window Data Filename: 20APR11M Sam:17 Analysis Date: 21-APR-11 Time: 01:02:13

DB-5 IS Data Filename: 20APR11M Sam:17 Analysis Date: 21-APR-11 Time: 01:02:13

DB-225 IS Date Filename: Analysis Date: Time:

DB-5 RT WINDOW DEFINING STANDARDS RESULTS

ISOMERS	ABSOLUTE RT	ISOMERS	ABSOLUTE RT
1,3,6,8-TCDD (F)	24:22 ✓	1,3,6,8-TCDF (F)	23:01 ✓
1,2,8,9-TCDD (L)	28:17 ✓	1,2,8,9-TCDF (L)	28:30 ✓
1,2,4,7,9-PeCDD (F)	30:12 ✓	1,3,4,6,8-PeCDF (F)	28:23 ✓
1,2,3,8,9-PeCDD (L)	33:45 ✓	1,2,3,8,9-PeCDF (L)	34:10 ✓
1,2,4,6,7,9-HxCDD (F)	36:06 ✓	1,2,3,4,6,8-HxCDF (F)	35:13 ✓
1,2,3,7,8,9-HxCDD (L)	39:10 ✓	1,2,3,7,8,9-HxCDF (L)	39:43 ✓
1,2,3,4,6,7,9-HpCDD (F)	42:46 ✓	1,2,3,4,6,7,8-HpCDF (F)	42:15 ✓
1,2,3,4,6,7,8-HpCDD (L)	44:09 ✓	1,2,3,4,7,8,9-HpCDF (L)	45:03 ✓

(F) = First eluting iosmer (DB-5); (L) = Last eluting isomer (DB-5)

=====

ISOMER SPECIFICITY (IS) TEST STANDARD RESULTS

% VALLEY HEIGHT
 BETWEEN
 COMPARED PEAKS (1)
 <25%

(1) To meet contract requirement, %Valley Height Between Compared Peaks shall not exceed 25% (section 15.4.2.2, Method 1613).

Analyst: 8

Date: 4/2/11

USEPA - ITD

FORM 6A

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 21-APR-11 01:02:13

CS3 or VER Data Filename: 20APR11M

Sam:17

NATIVE ANALYTES	RETENTION TIME REFERENCE	RRT	RRT QC LIMITS (1)
2,3,7,8-TCDD	13C-2,3,7,8-TCDD	1.001	0.999-1.002 ✓
2,3,7,8-TCDF	13C-2,3,7,8-TCDF	1.001	0.999-1.003 ✓
1,2,3,7,8-PeCDD	13C-1,2,3,7,8-PeCDD	1.001	0.999-1.002 ✓
1,2,3,7,8-PeCDF	13C-1,2,3,7,8-PeCDF	1.000	0.999-1.002 ✓
2,3,4,7,8-PeCDF	13C-2,3,4,7,8-PeCDF	1.001	0.999-1.002 ✓
LABELED COMPOUNDS			
37Cl-2,3,7,8-TCDD	13C-1,2,3,4-TCDD	1.021	0.989-1.052 ✓
13C-2,3,7,8-TCDD		1.021	0.976-1.043 ✓
13C-2,3,7,8-TCDF		0.993	0.923-1.103 ✓
13C-1,2,3,7,8-PeCDD		1.239	1.000-1.567 ✓
13C-1,2,3,7,8-PeCDF		1.174	0.923-1.203 ✓
13C-2,3,4,7,8-PeCDF		1.223	0.923-1.303 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified in Table 2, Method 1613.

Analyst: 

Date: 4/21/11

USEPA - ITD

FORM 6B

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11


Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 21-APR-11 01:02:13 CS3 or VER Data Filename: 20APR11M Sam:17

NATIVE ANALYTES	RETENTION TIME REFERENCE	RRT	RRT QC LIMITS (1)
1,2,3,4,7,8-HxCDD	13C-1,2,3,4,7,8-HxCDD	1.001	0.999-1.001 ✓
1,2,3,6,7,8-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.001	0.998-1.004 ✓
1,2,3,7,8,9-HxCDD	13C-1,2,3,6,7,8-HxCDD	1.012	1.000-1.019 ✓
1,2,3,4,7,8-HxCDF	13C-1,2,3,4,7,8-HxCDF	1.000	0.999-1.001 ✓
1,2,3,6,7,8-HxCDF	13C-1,2,3,6,7,8-HxCDF	1.001	0.997-1.005 ✓
2,3,4,6,7,8-HxCDF	13C-2,3,4,6,7,8-HxCDF	1.001	0.999-1.001 ✓
1,2,3,7,8,9-HxCDF	13C-1,2,3,7,8,9-HxCDF	1.001	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDD	13C-1,2,3,4,6,7,8-HpCDD	1.001	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDF	13C-1,2,3,4,6,7,8-HpCDF	1.000	0.999-1.001 ✓
1,2,3,4,7,8,9-HpCDF	13C-1,2,3,4,7,8,9-HpCDF	1.001	0.999-1.001 ✓
OCDD	13C-OCDD	1.001	0.999-1.001 ✓
OCDF	13C-OCDF	1.001	0.999-1.001 ✓
LABELED COMPOUNDS			
13C-1,2,3,4,7,8-HxCDD	13C-1,2,3,7,8,9-HxCDD	0.984	0.977-1.000 ✓
13C-1,2,3,6,7,8-HxCDD		0.989	0.981-1.003 ✓
13C-1,2,3,4,7,8-HxCDF		0.949	0.944-0.970 ✓
13C-1,2,3,6,7,8-HxCDF		0.954	0.949-0.975 ✓
13C-2,3,4,6,7,8-HxCDF		0.978	0.959-1.021 ✓
13C-1,2,3,7,8,9-HxCDF		1.015	0.977-1.047 ✓
13C-1,2,3,4,6,7,8-HpCDD		1.128	1.086-1.130 ✓
13C-1,2,3,4,6,7,8-HpCDF		1.080	1.043-1.085 ✓
13C-1,2,3,4,7,8,9-HpCDF		1.151	1.057-1.154 ✓
13C-OCDD		1.269	1.032-1.311 ✓
13C-OCDF		1.279	1.000-1.311 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified
in Table 2, Method 1613.

Analyst: Date: 

Name	Resp	RA	RT	RRF	NATO 1989 Tox:		WHO 1998 Tox:		WHO 2005 Tox:		DL
					Conc	Qual	Fac	Noise-1	Noise-2	DL	
2,3,7,8-TCDD	1.97e+06	0.77 y	27:21	1.13	8.99	2.50	-	-	-	*	
1,2,3,7,8-PeCDD	9.18e+06	1.62 y	33:11	1.02	53.0	2.50	-	-	-	*	
1,2,3,4,7,8-HxCDD	1.14e+07	1.33 y	38:33	1.45	51.3	2.50	-	-	-	*	
1,2,3,6,7,8-HxCDD	9.27e+06	1.32 y	38:43	1.45	50.9	2.50	-	-	-	*	
1,2,3,7,8,9-HxCDD	1.08e+07	1.32 y	39:10	1.47	52.9	2.50	-	-	-	*	
1,2,3,4,6,7,8-HpCDD	8.73e+06	0.96 y	44:09	1.30	48.9	2.50	-	-	-	*	
OCDD	1.33e+07	0.93 y	49:42	1.45	98.2	2.50	-	-	-	*	
2,3,7,8-TCDF	4.14e+06	0.66 y	26:36	1.15	10.7	2.50	-	-	-	*	
1,2,3,7,8-PeCDF	1.33e+07	1.68 y	31:27	0.89	53.8	2.50	-	-	-	*	
2,3,4,7,8-PeCDF	1.27e+07	1.65 y	32:46	0.89	52.1	2.50	-	-	-	*	
1,2,3,4,7,8-HxCDF	1.27e+07	1.27 y	37:09	1.01	49.5	2.50	-	-	-	*	
1,2,3,6,7,8-HxCDF	1.42e+07	1.26 y	37:21	0.89	50.2	2.50	-	-	-	*	
2,3,4,6,7,8-HxCDF	1.32e+07	1.26 y	38:17	1.02	49.3	2.50	-	-	-	*	
1,2,3,7,8,9-HxCDF	1.63e+07	1.27 y	39:43	1.10	50.3	2.50	-	-	-	*	
1,2,3,4,6,7,8-HpCDF	1.14e+07	1.04 y	42:15	1.48	48.9	2.50	-	-	-	*	
1,2,3,4,7,8,9-HpCDF	9.28e+06	1.05 y	45:03	1.43	48.9	2.50	-	-	-	*	
OCDF	1.45e+07	0.89 y	50:04	0.84	96.4	2.50	-	-	-	*	
										Rec	
13C-2,3,7,8-TCDD	1.93e+07	0.82 y	27:20	1.03	95.8					95.8	
13C-1,2,3,7,8-PeCDD	1.71e+07	1.76 y	33:10	1.01	86.0					86.0	
13C-1,2,3,4,7,8-HxCDD	1.53e+07	1.29 y	38:31	1.19	94.6					94.6	
13C-1,2,3,6,7,8-HxCDD	1.25e+07	1.27 y	38:41	0.94	98.4					98.4	
13C-1,2,3,4,6,7,8-HpCDD	1.37e+07	1.06 y	44:07	0.83	122					122	
13C-OCDD	1.87e+07	0.97 y	49:40	0.61	226					113	
13C-2,3,7,8-TCDF	3.36e+07	0.87 y	26:35	0.98	102					102	
13C-1,2,3,7,8-PeCDF	2.78e+07	1.73 y	31:26	0.83	99.6					99.6	
13C-2,3,4,7,8-PeCDF	2.73e+07	1.71 y	32:44	0.80	101					101	
13C-1,2,3,4,7,8-HxCDF	2.55e+07	0.48 y	37:08	1.84	102					102	
13C-1,2,3,6,7,8-HxCDF	3.16e+07	0.48 y	37:20	2.29	102					102	
13C-2,3,4,6,7,8-HxCDF	2.62e+07	0.48 y	38:16	1.86	104					104	
13C-1,2,3,7,8,9-HxCDF	2.93e+07	0.49 y	39:42	1.98	109					109	
13C-1,2,3,4,6,7,8-HpCDF	1.58e+07	0.45 y	42:14	0.99	118					118	
13C-1,2,3,4,7,8,9-HpCDF	1.33e+07	0.44 y	45:02	0.77	128					128	
13C-OCDF	3.56e+07	0.95 y	50:02	1.17	224					112	
37Cl-2,3,7,8-TCDD	1.14e+06		27:21	0.73	7.97					79.7	
13C-1,2,3,4-TCDD	1.96e+07	0.82 y	26:46	-	51.6						
13C-1,2,3,4-TCDF	3.37e+07	0.88 y	25:30	-	46.8						
13C-1,2,3,7,8,9-HxCDD	1.36e+07	1.28 y	39:08	-	54.8						
						Fac	Noise-1	Noise-2	DL	#Hom	
Total Tetra-Dioxins	1.06e+07		23:44	1.13	48.4	2.50	576	-	0.0276	22	
Total Penta-Dioxins	1.93e+07		30:12	1.02	112	2.50	1510	-	0.103	8	
Total Hexa-Dioxins	3.51e+07		36:06	1.46	173	2.50	1910	-	0.119	6	
Total Hepta-Dioxins	1.83e+07		42:46	1.30	103	2.50	608	-	0.0455	3	
Total Tetra-Furans	1.89e+07		23:01	1.15	49.0	2.50	884	-	0.0241	9	
1st Fn. Tot Penta-Furans	1.36e+07		28:23	0.89	55.5	2.50	432	-	0.0207	PeCDF 4	
Total Penta-Furans	3.78e+07		30:08	0.89	154	2.50	1990	-	0.0957	210 9	
Total Hexa-Furans	6.44e+07		35:13	1.00	228	2.50	1520	-	0.0701	11	
Total Hepta-Furans	2.09e+07		42:15	1.46	98.6	2.50	-	-	*	6	

Analyst: J

Date: 4/21/11

Frontier Analytical Laboratory - Acquisition Log

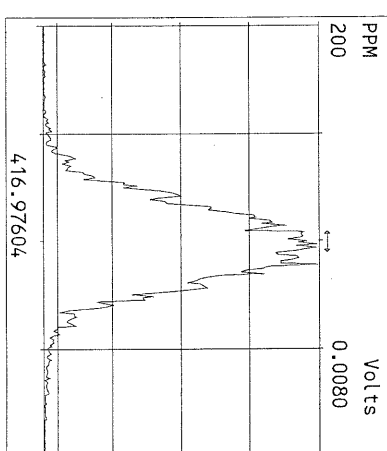
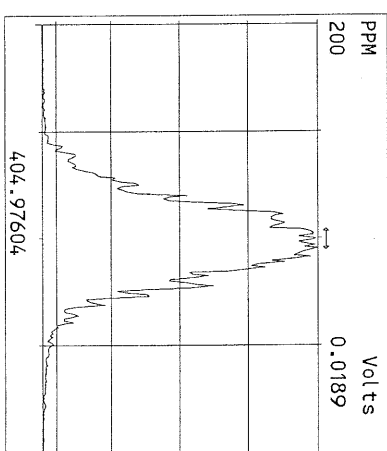
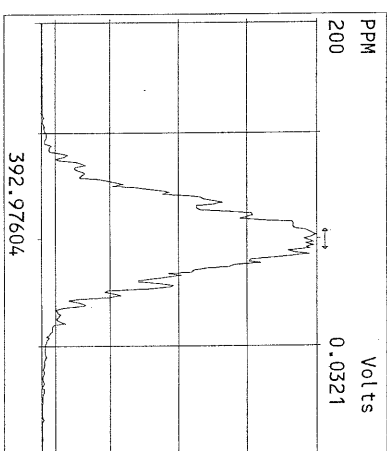
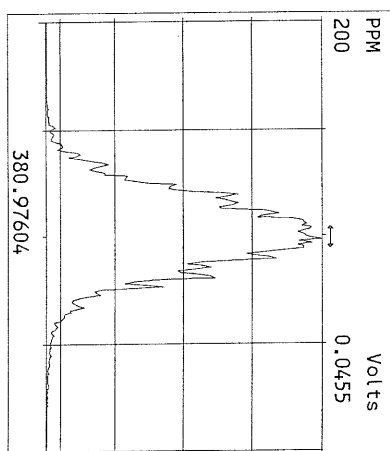
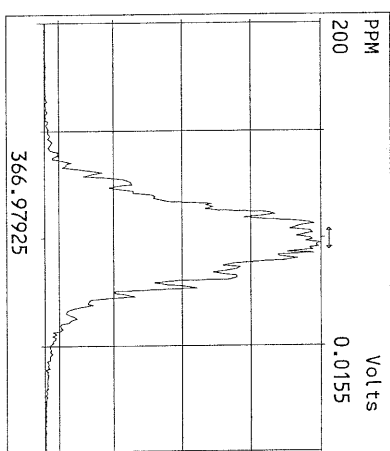
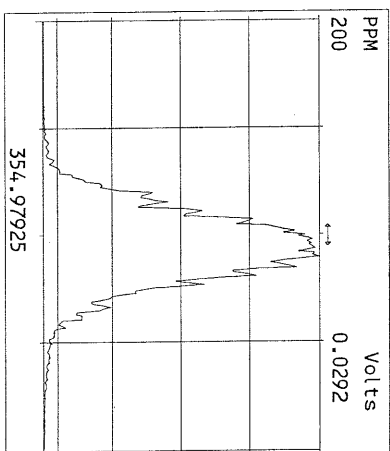
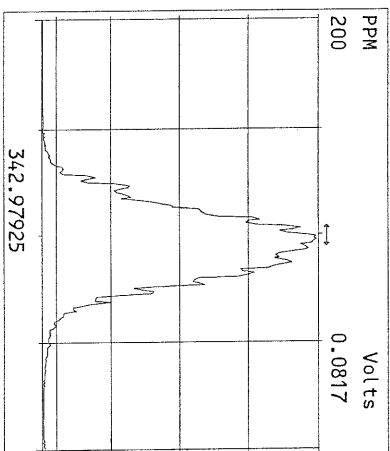
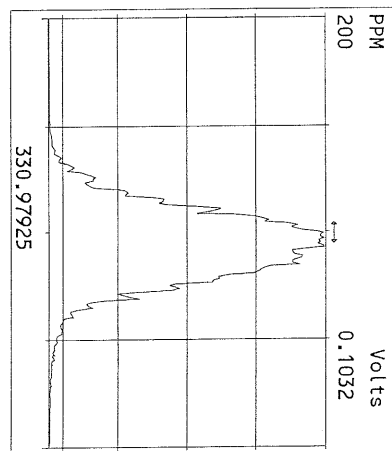
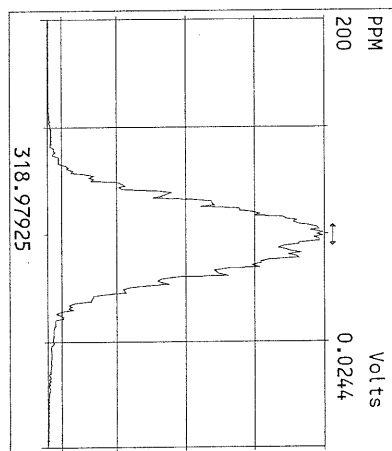
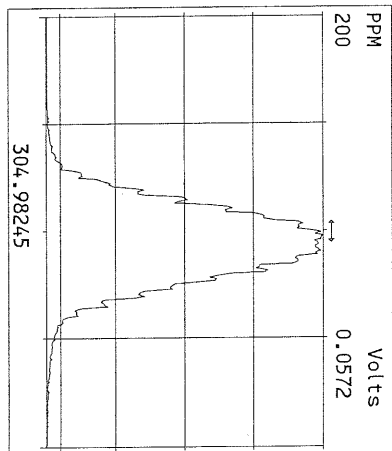
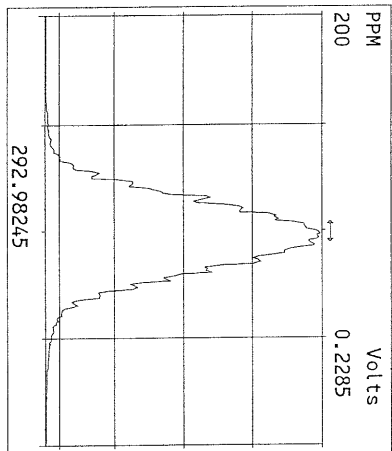
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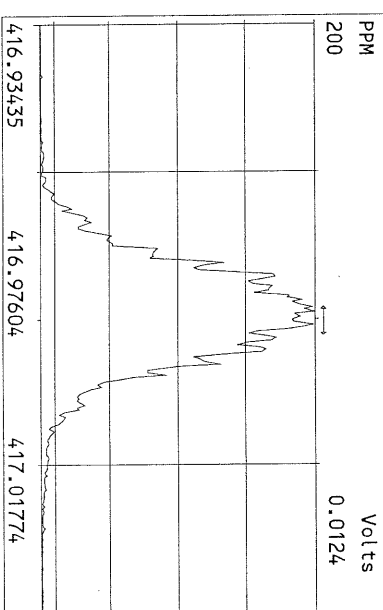
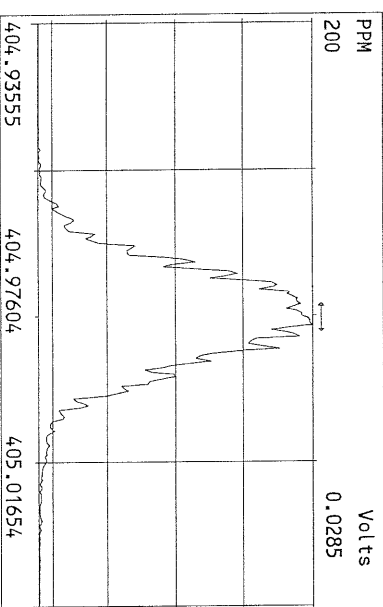
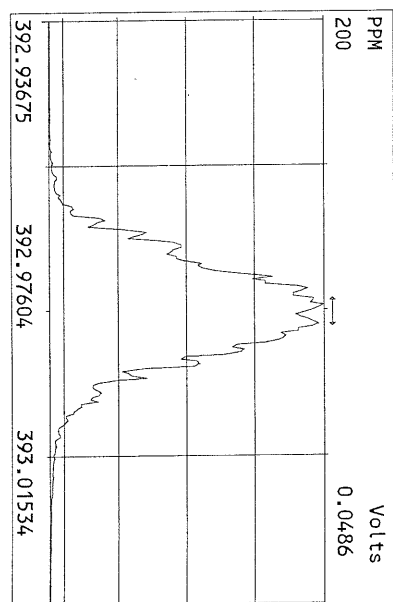
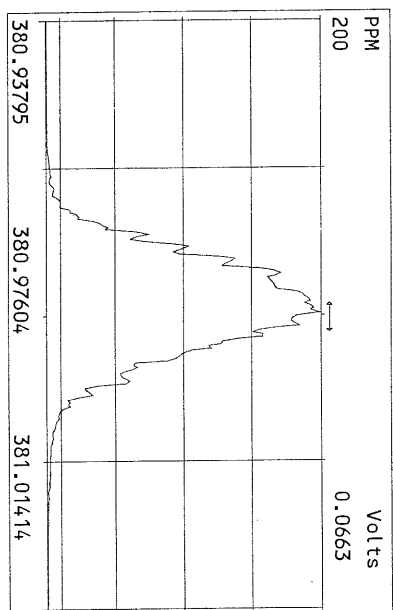
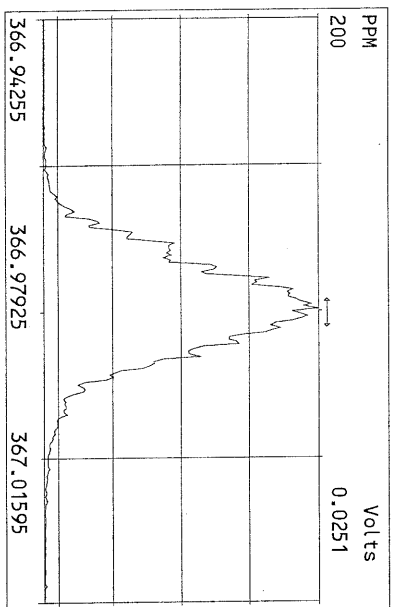
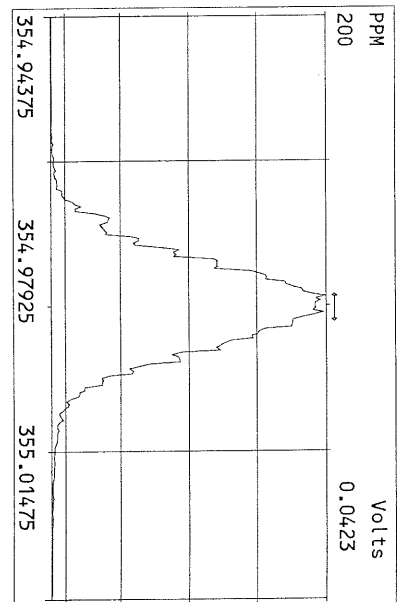
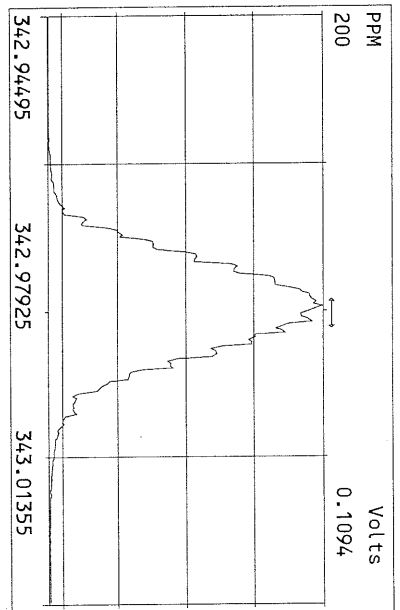
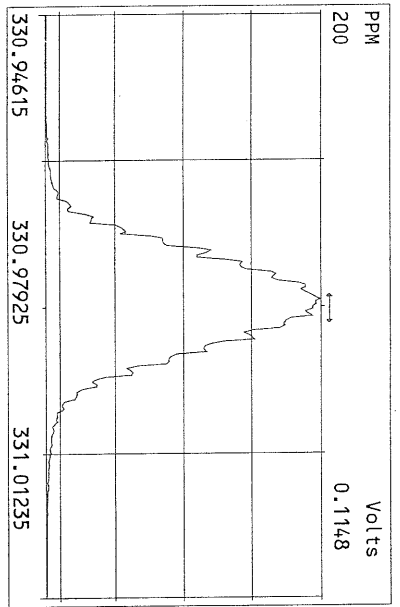
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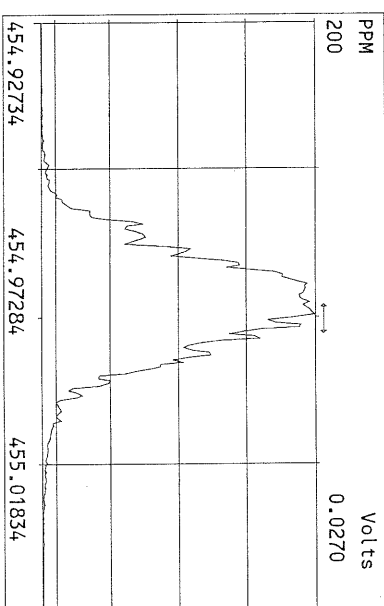
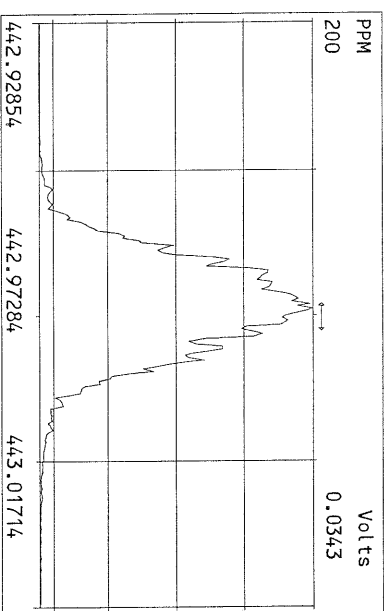
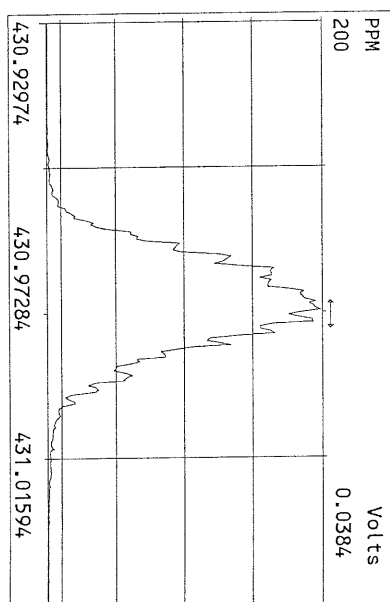
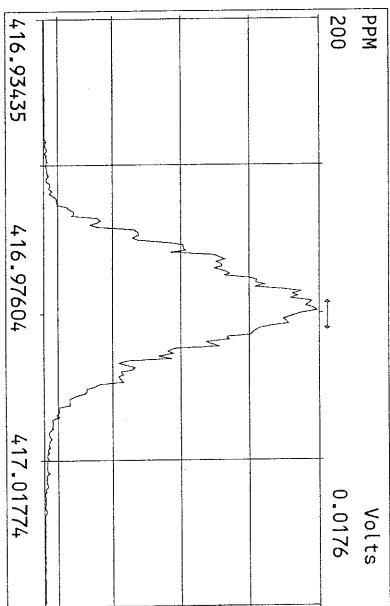
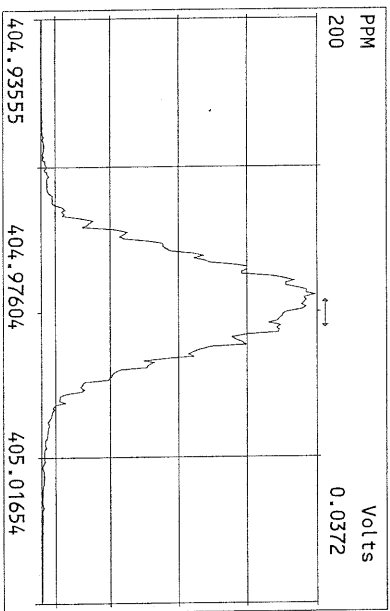
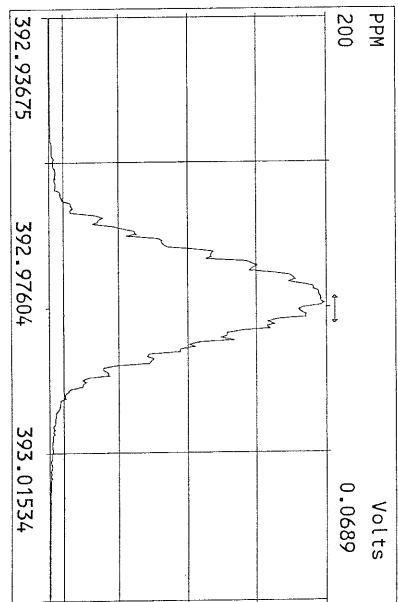
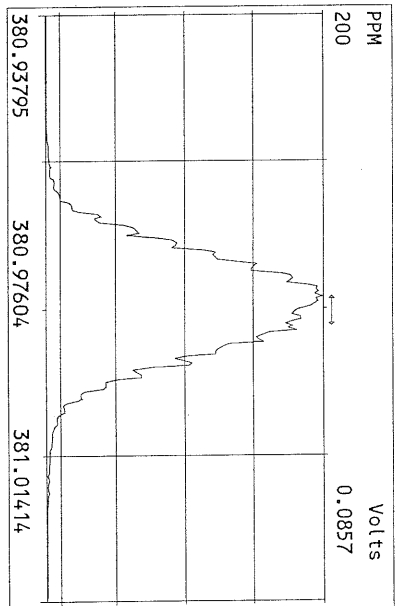
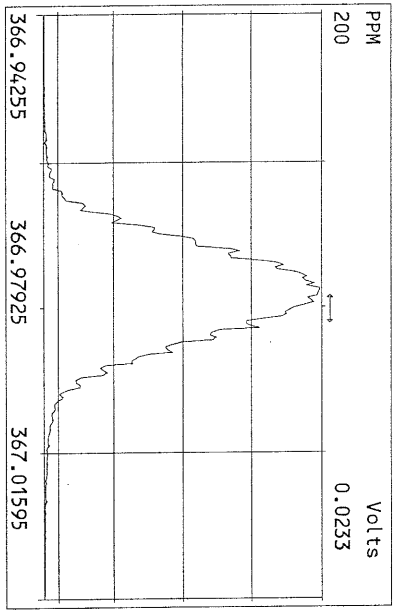
8/4/21/11

Data Backed Up: _____

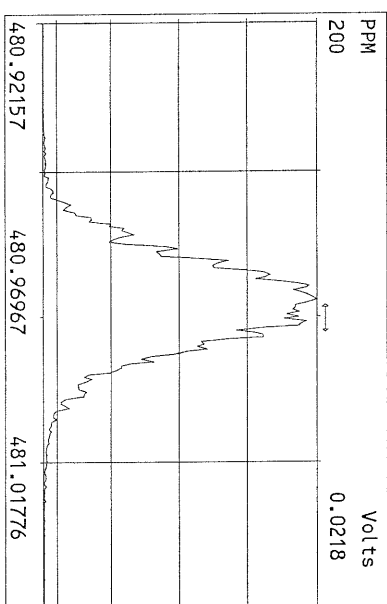
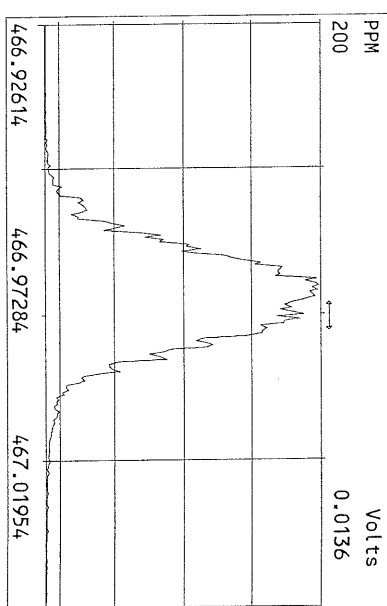
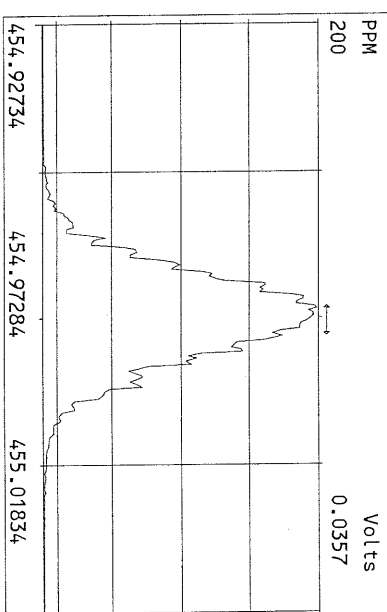
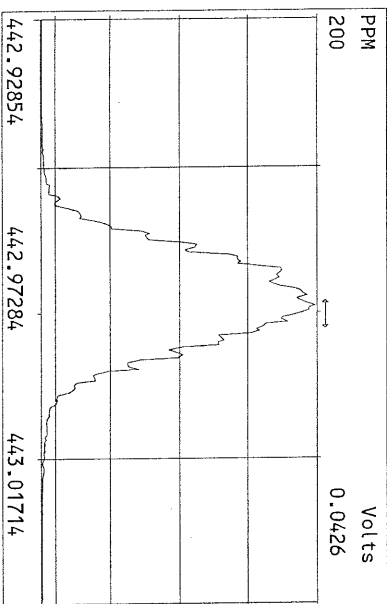
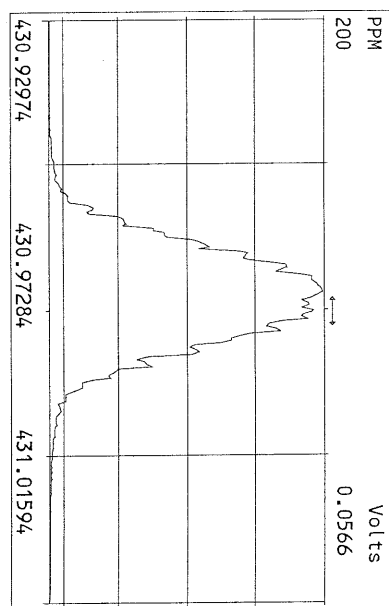
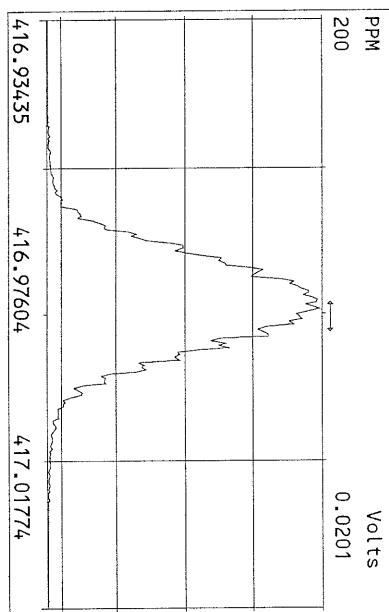
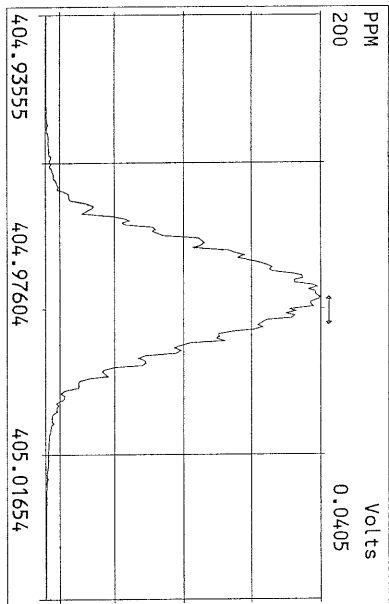
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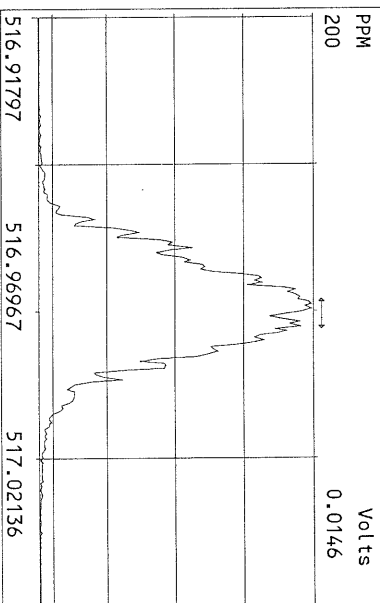
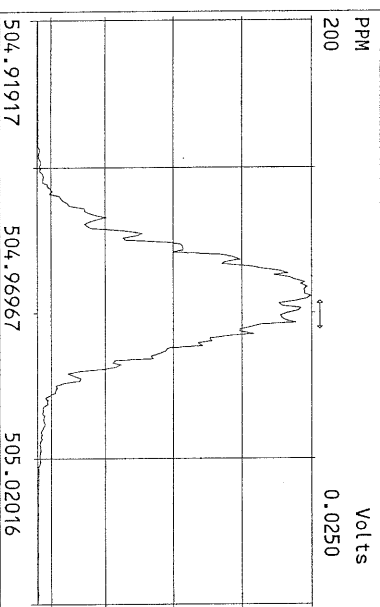
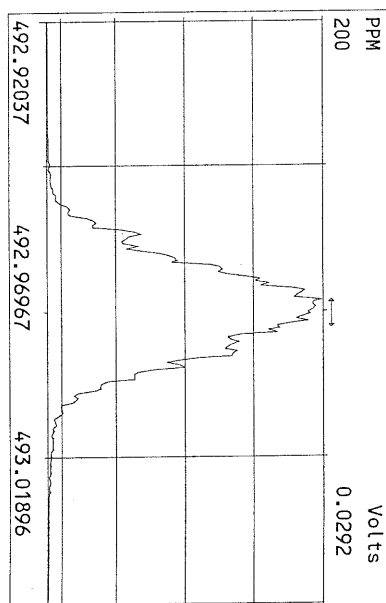
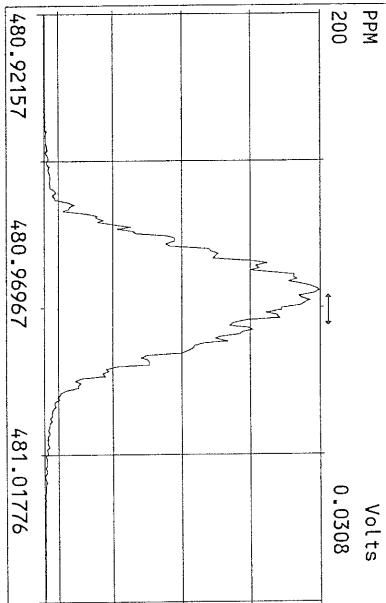
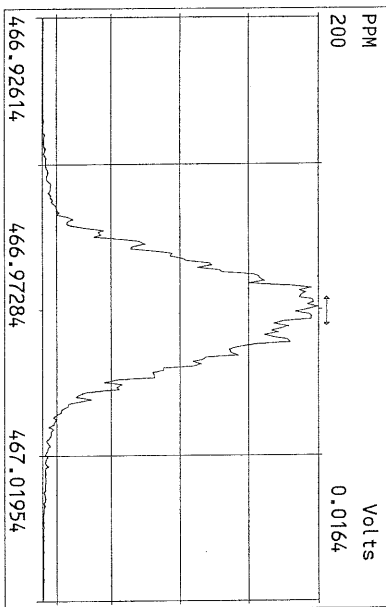
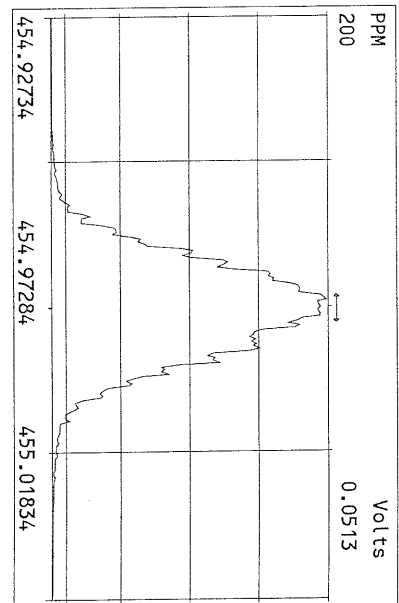
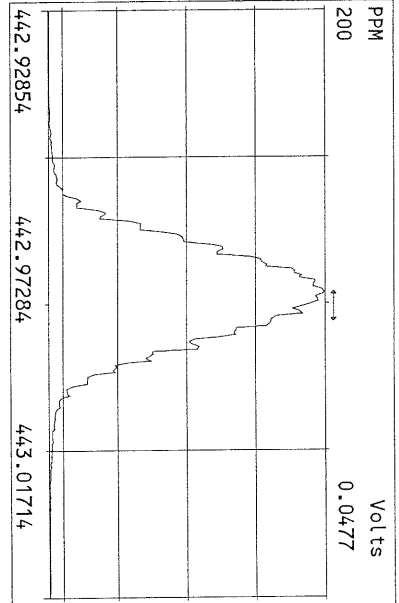
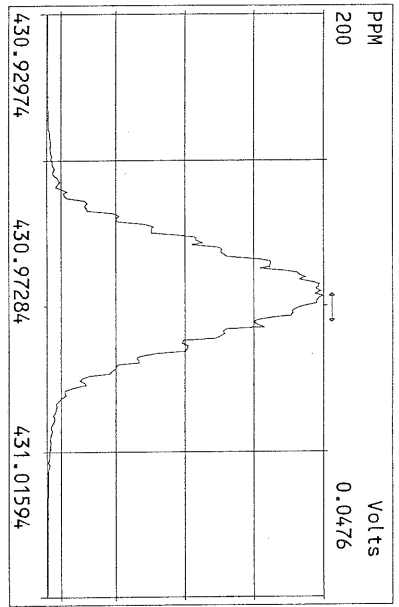




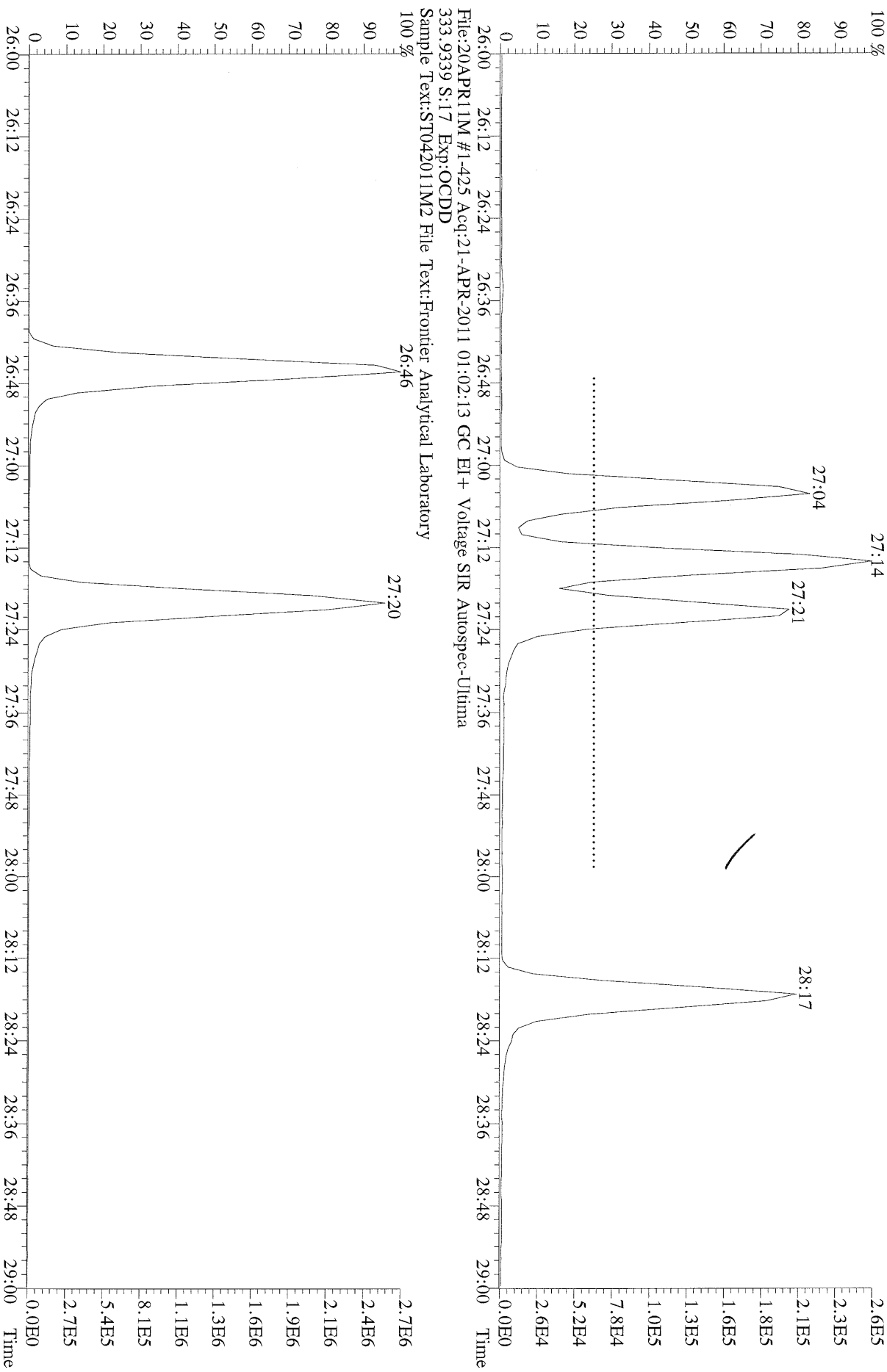


Peak Locate Examination:20-APR-2011:10:15 File:20APR11M
Experiment:OCDD Function:4 Reference:PFK

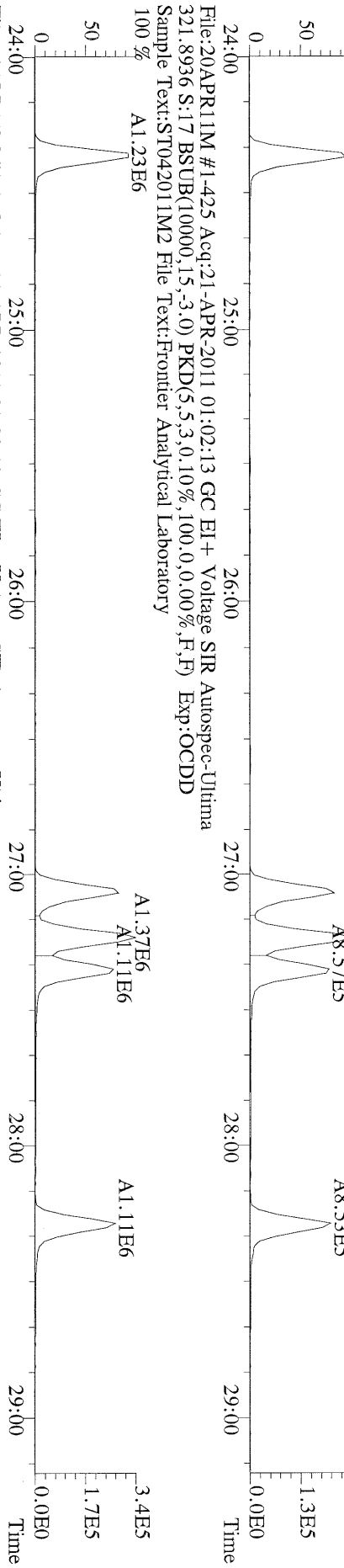




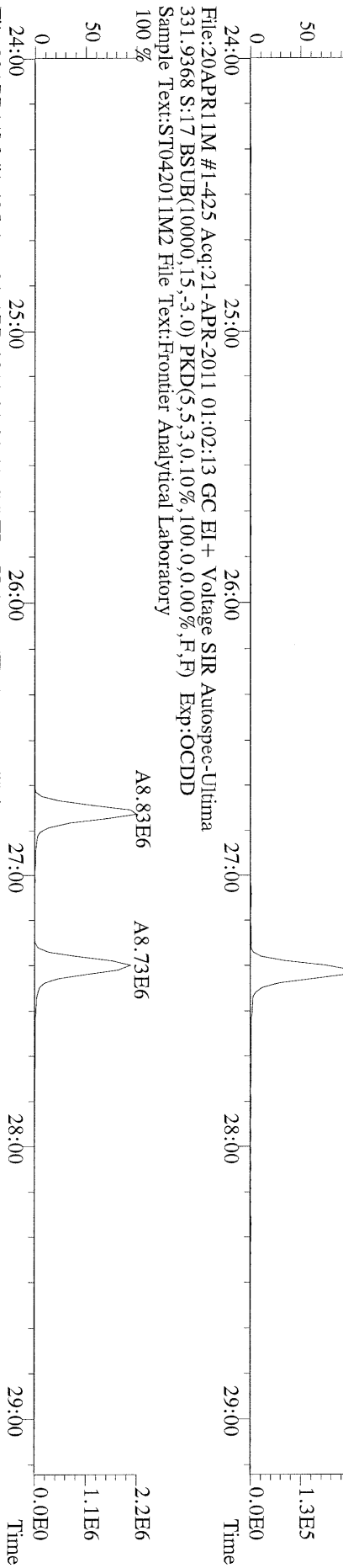
File:20APR11M #1-425 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Utima
319.8965 S:17 Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



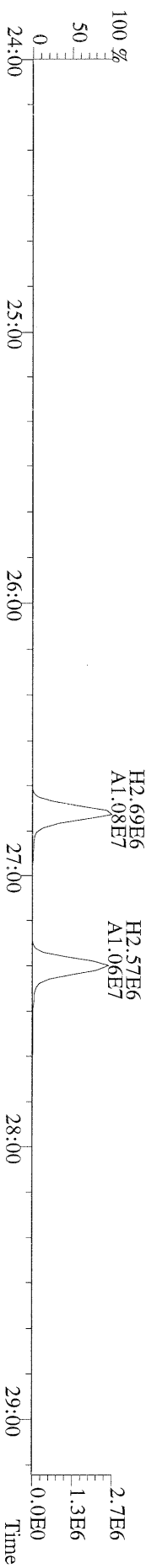
File:20APR11M #1-425 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Utima
319.8965 S:1.7 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 % A9.45E5



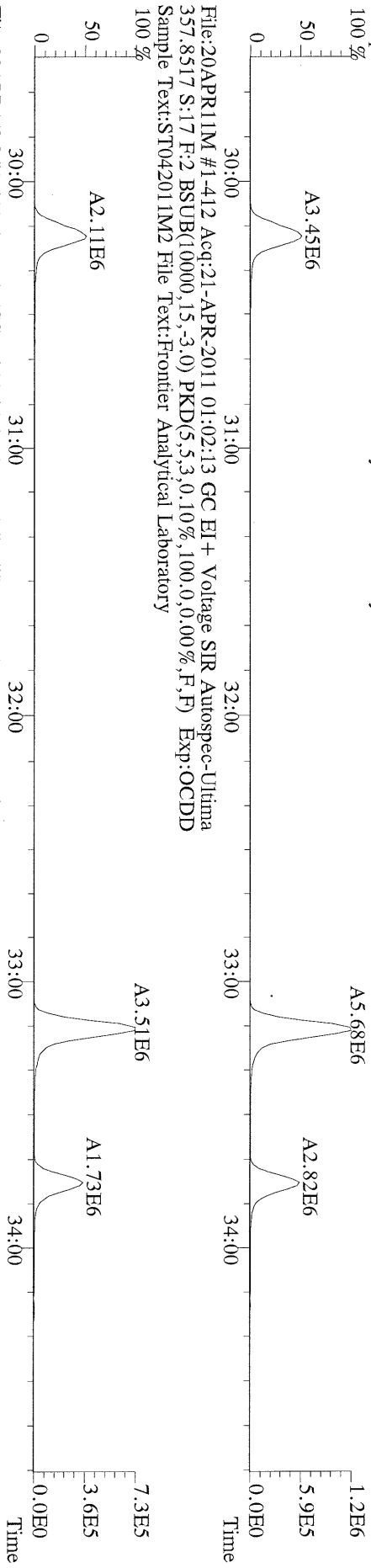
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327.8847 S:1.7 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 % A1.14E6



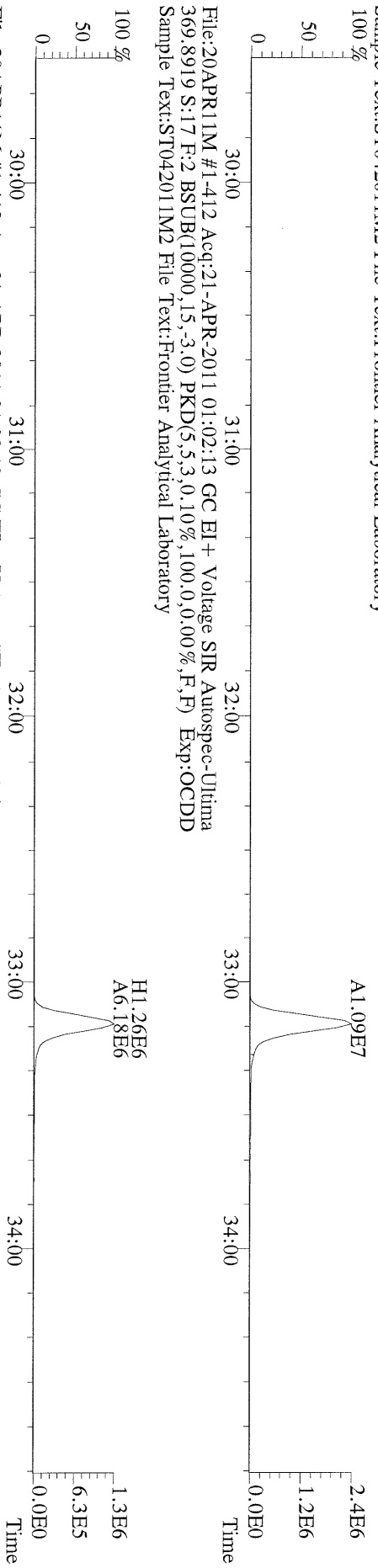
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333.9339 S:1.7 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



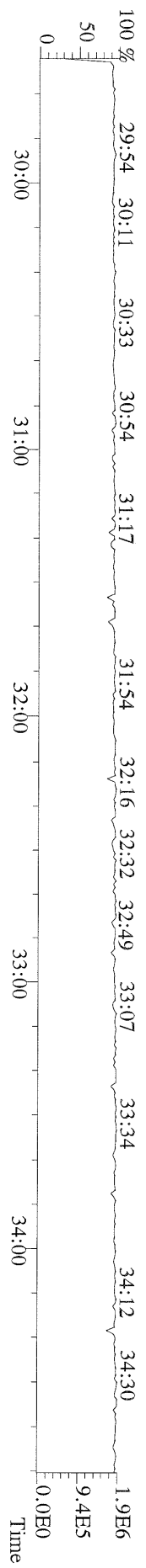
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355,8546 S:17 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



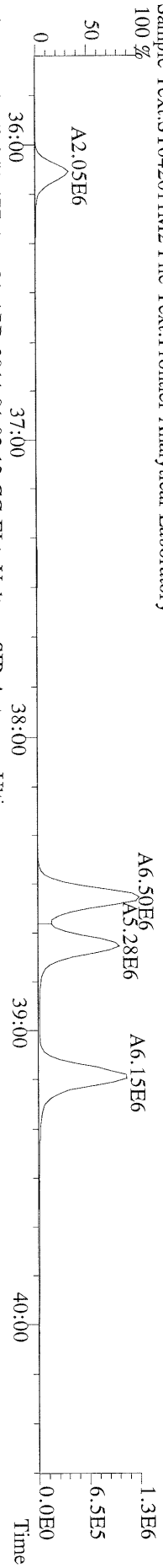
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367,8949 S:17 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



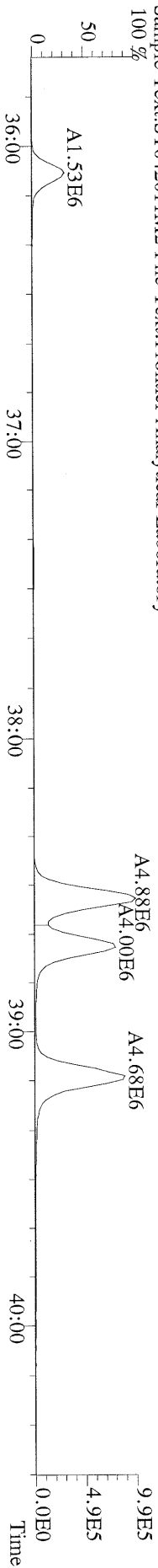
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366,9792 S:17 F:2 Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



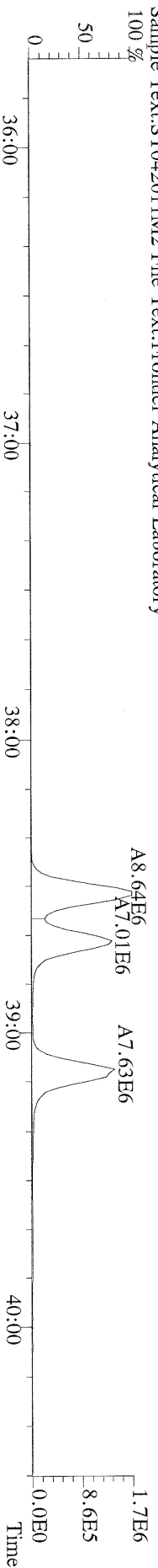
File:20APR11M #1-477 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
389.8156 S:17 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



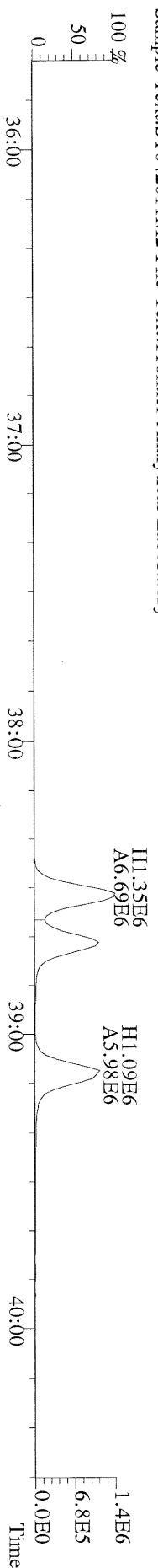
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Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



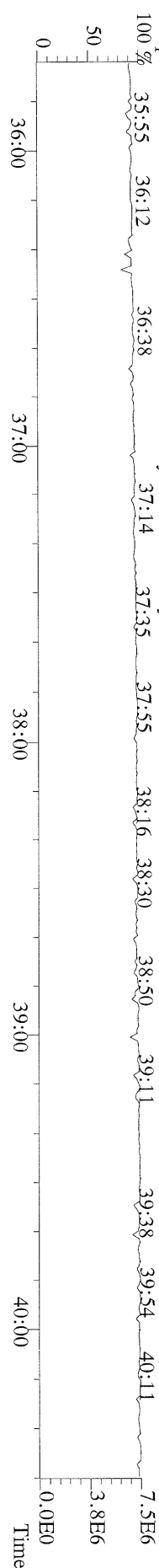
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401.8559 S:17 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



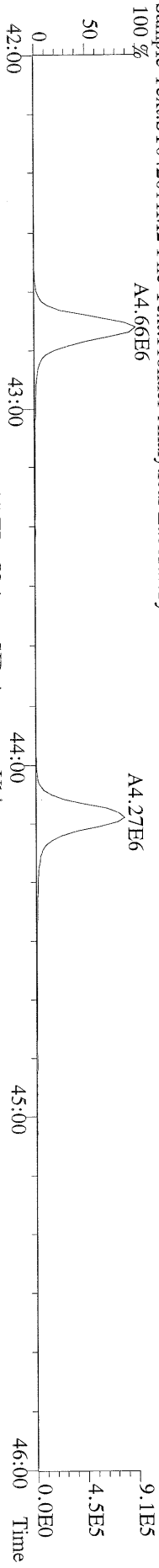
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Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



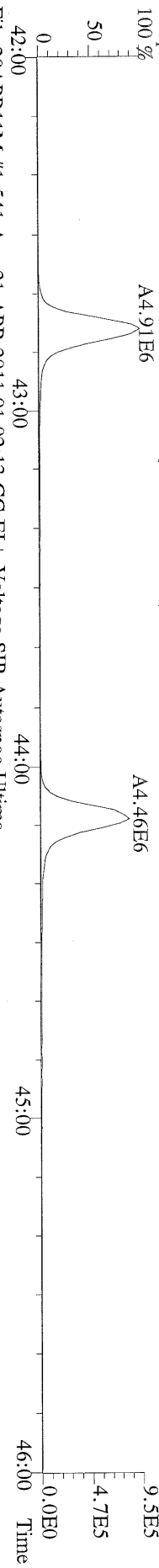
File:20APR11M #1-477 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
380.9760 S:17 F:3 Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



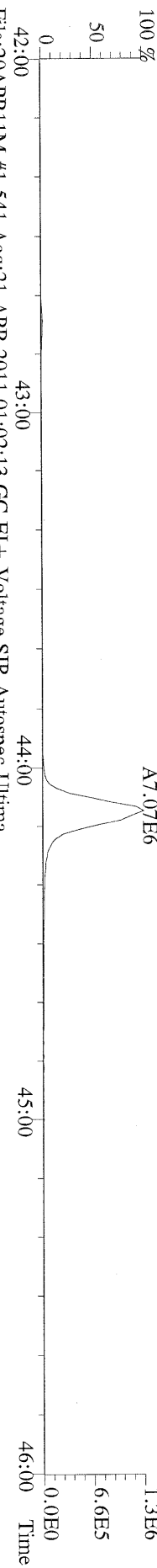
File:20APR11M #1-541 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
423.7767 S:17 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



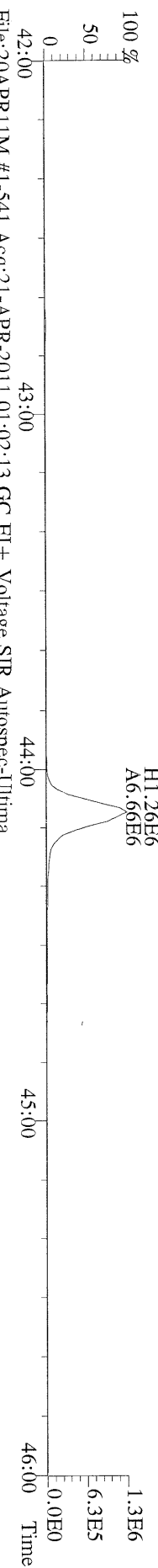
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425.7737 S:17 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



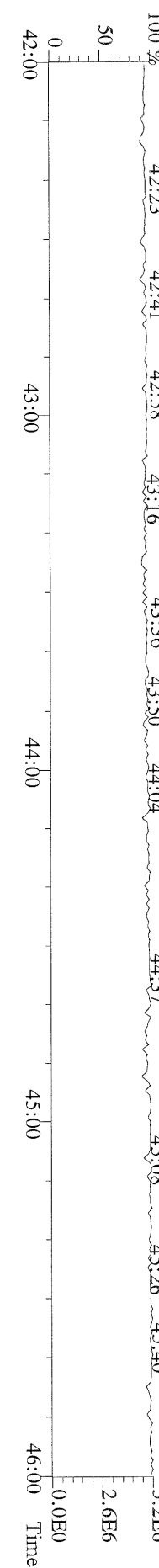
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435.8169 S:17 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



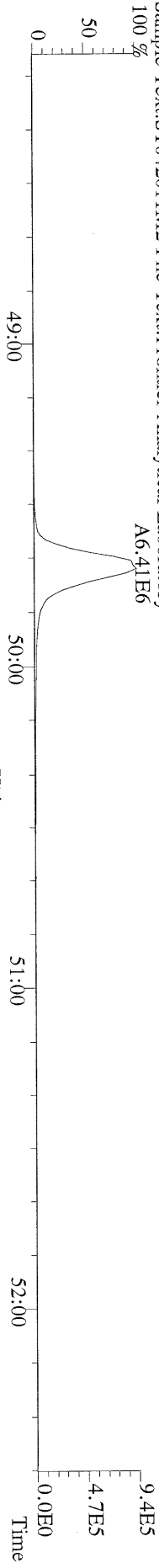
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437.8140 S:17 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



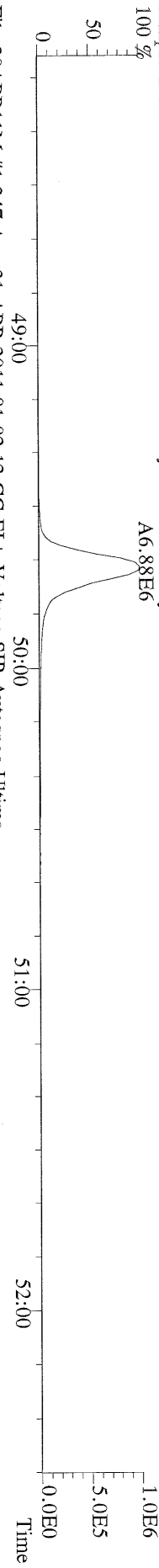
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430.9728 S:17 F:4 Exp:OCDD
Sample Text:ST04201IM2 File Text:Frontier Analytical Laboratory



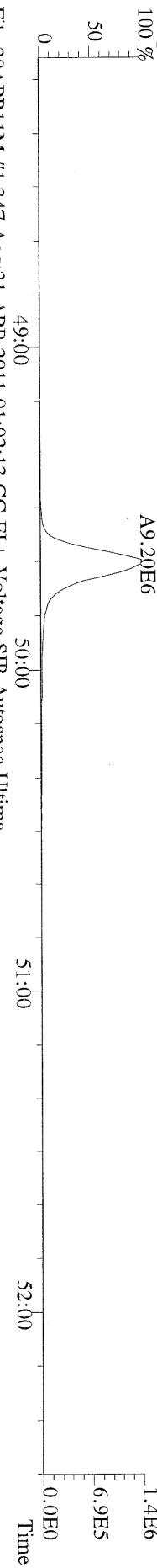
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457.7377 S:17 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Fronter Analytical Laboratory



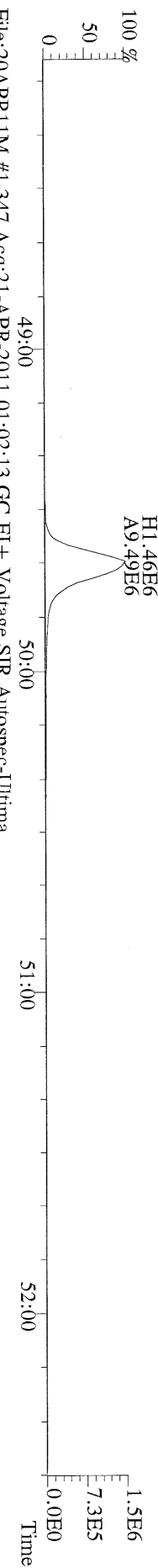
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459.7348 S:17 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Fronter Analytical Laboratory



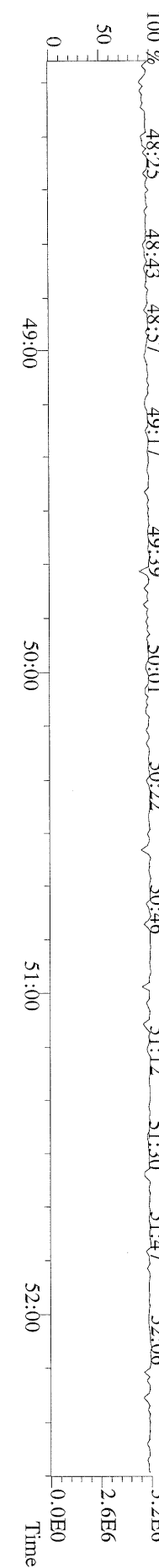
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Sample Text:ST04201IM2 File Text:Fronter Analytical Laboratory



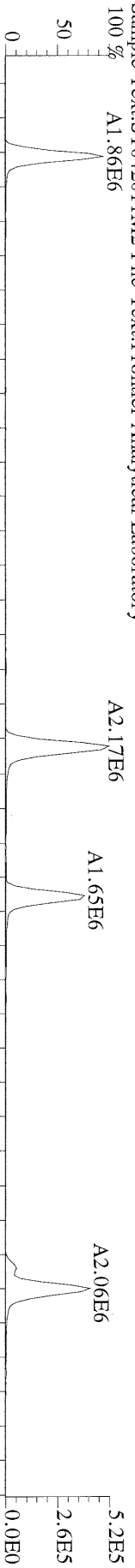
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471.7750 S:17 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM2 File Text:Fronter Analytical Laboratory



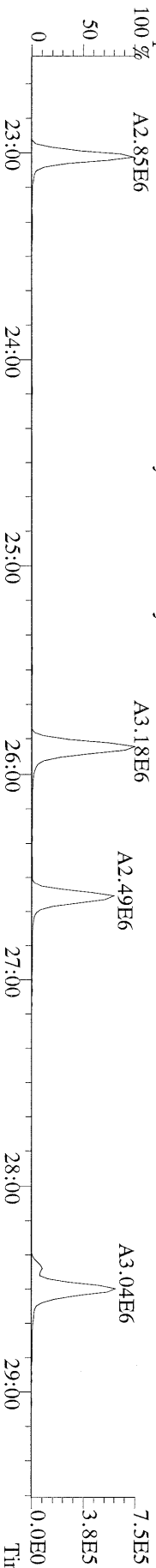
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454.9728 S:17 F:5 Exp:OCDD
Sample Text:ST04201IM2 File Text:Fronter Analytical Laboratory



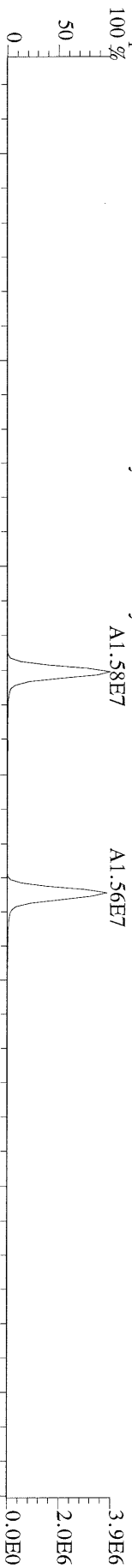
File:20APR11M #1-425 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
303.9016 S:17 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



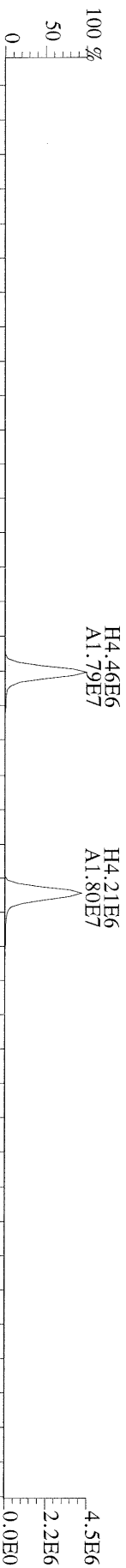
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305.8987 S:17 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



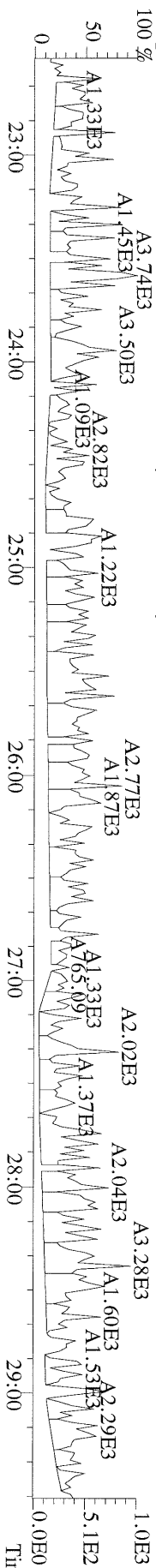
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315.9419 S:17 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



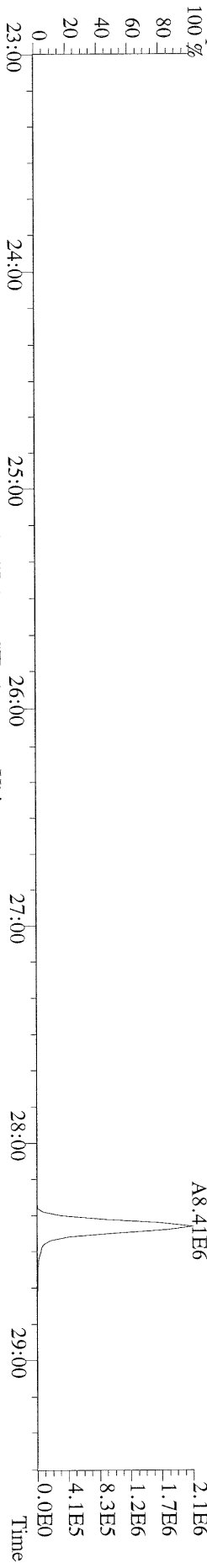
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317.9389 S:17 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



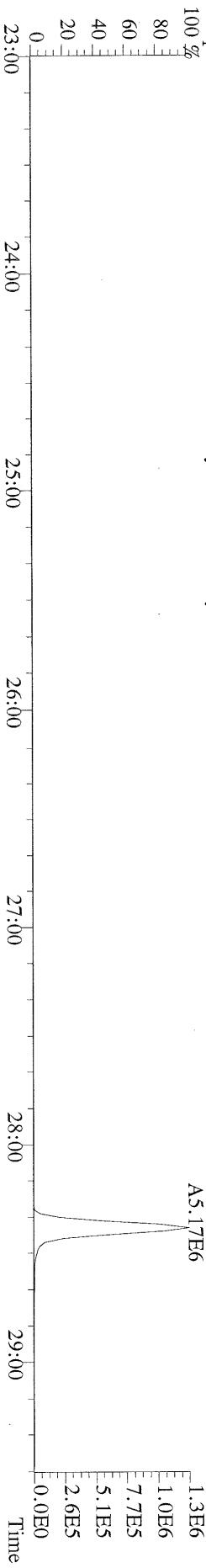
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375.8364 S:17 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



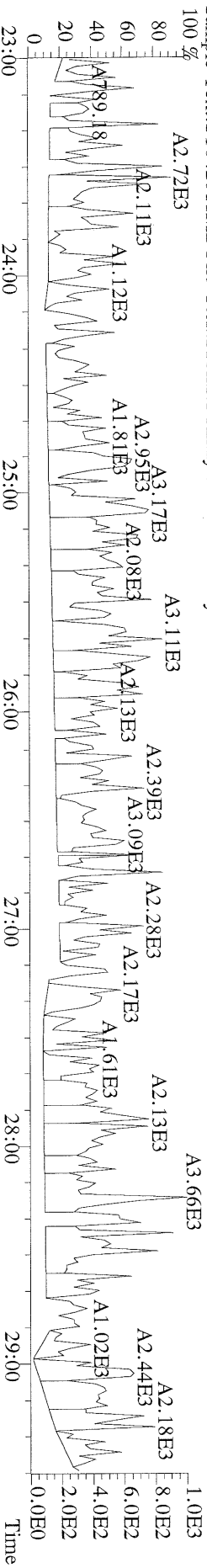
File:20APR11M #1-425 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
 339.8597 S:17 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0,0%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



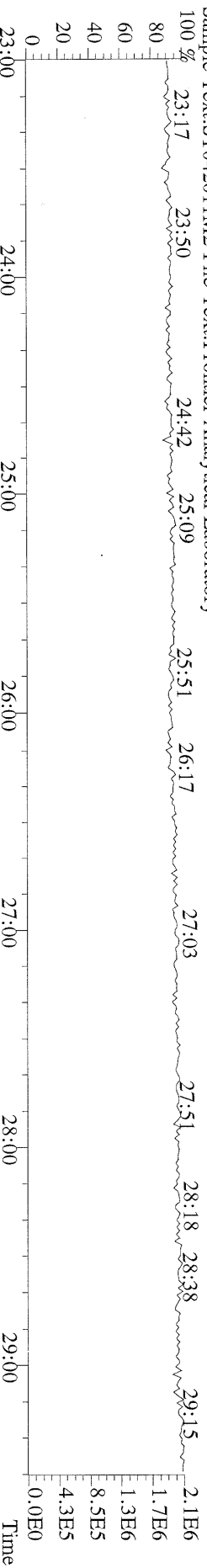
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 341.8568 S:17 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0,0%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



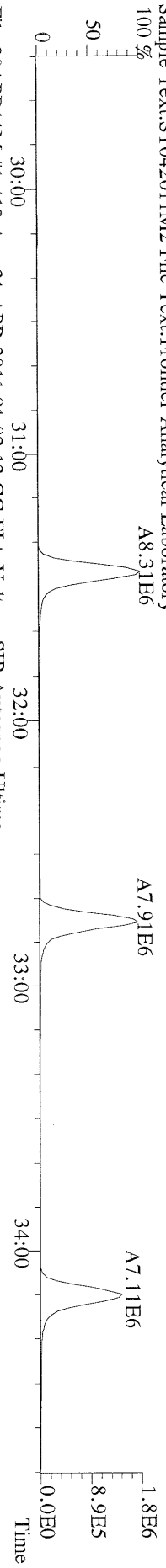
File:20APR11M #1-425 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
 409.7974 S:17 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0,0%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



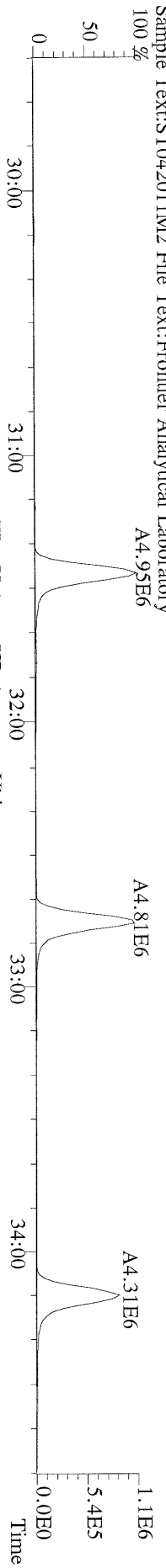
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 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



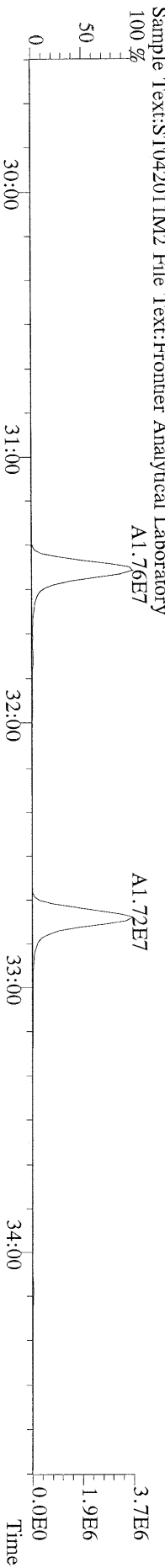
File:20APR11M #1-412 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
 339.8597 S:17 F:2 BSUB(10000,15,-3.0) PKD(5.5,3.0,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



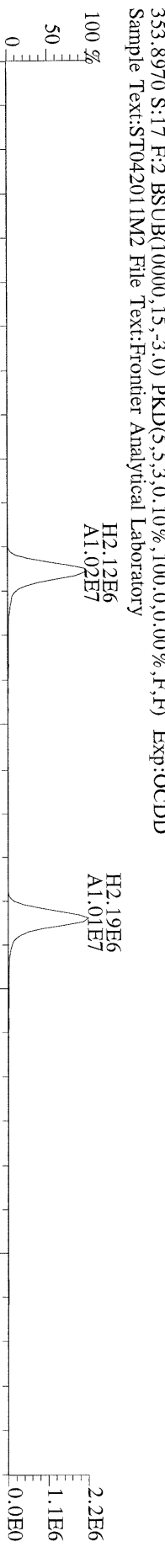
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 341.8568 S:17 F:2 BSUB(10000,15,-3.0) PKD(5.5,3.0,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



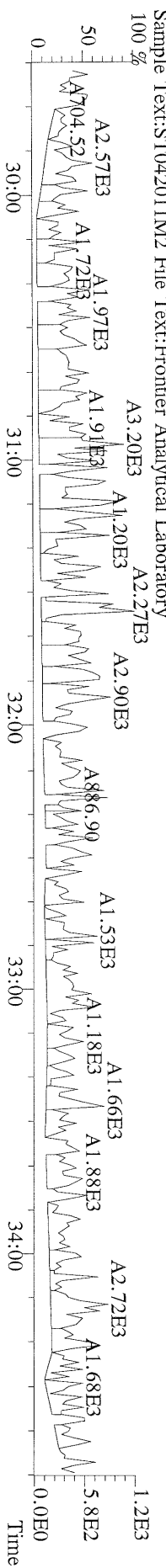
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 351.9000 S:17 F:2 BSUB(10000,15,-3.0) PKD(5.5,3.0,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



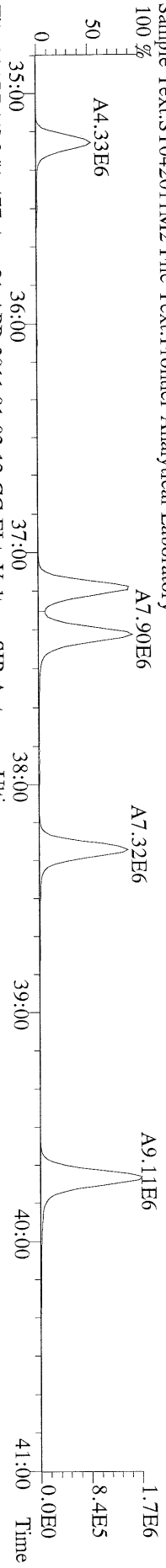
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 353.8970 S:17 F:2 BSUB(10000,15,-3.0) PKD(5.5,3.0,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



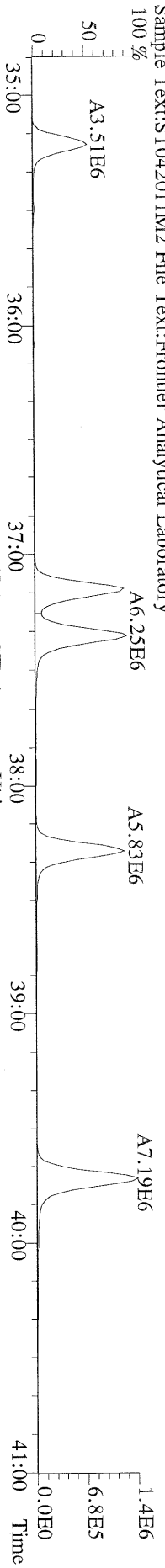
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 409.7974 S:17 F:2 BSUB(10000,15,-3.0) PKD(5.5,3.0,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



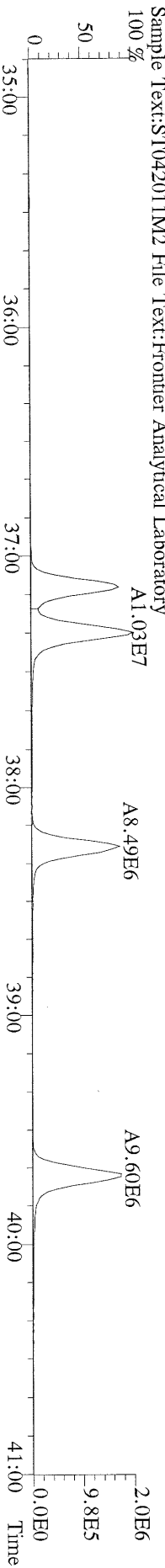
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373.8207 S:17 F:3 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



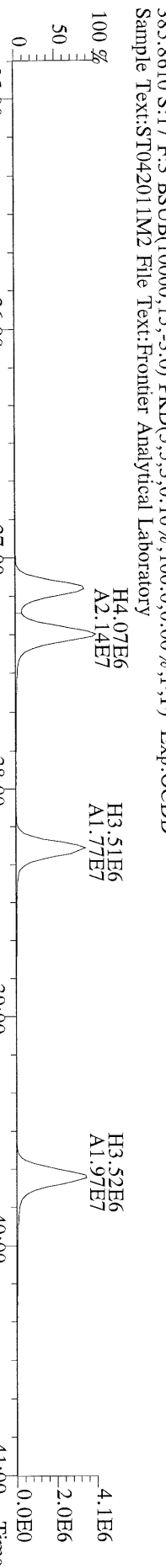
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Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



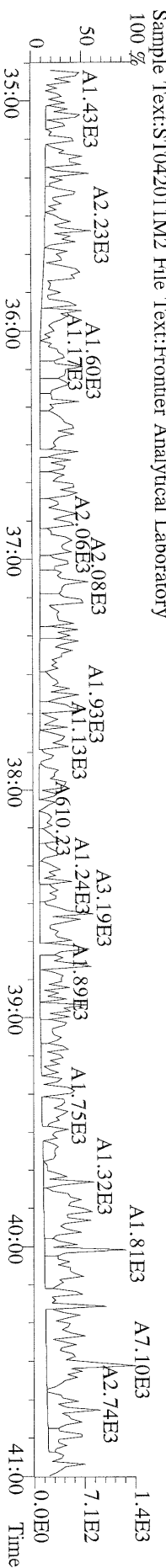
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383.8639 S:17 F:3 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



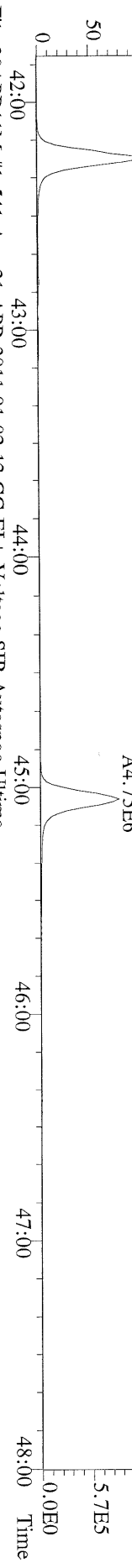
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385.8610 S:17 F:3 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



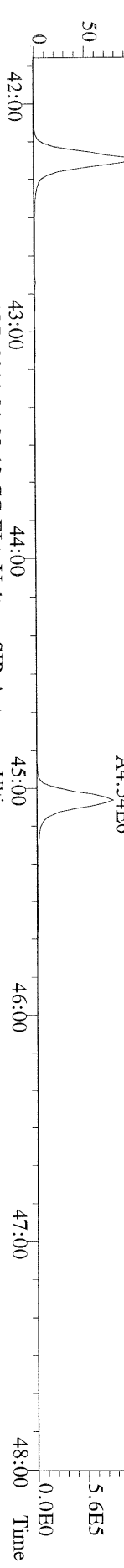
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445.7555 S:17 F:3 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



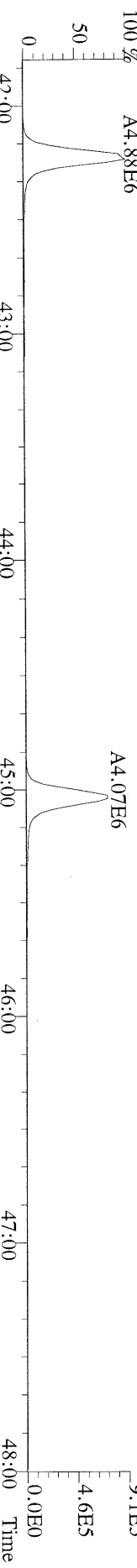
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407.7818 S:17 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 % A5.83E6



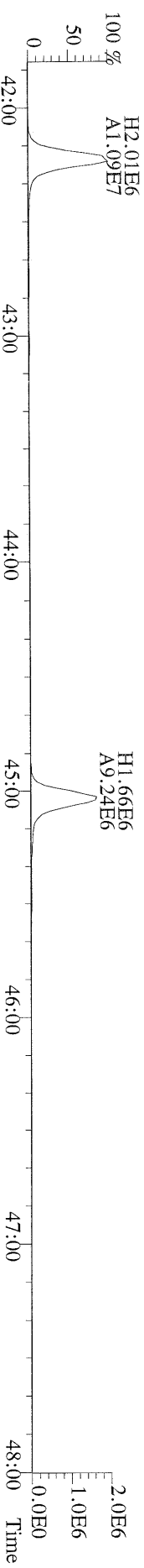
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409.7788 S:17 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 % A5.59E6



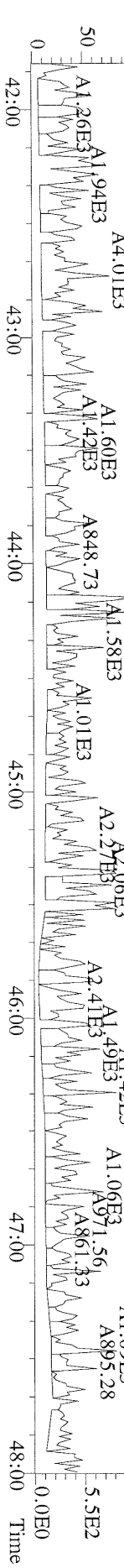
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Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 % A4.88E6



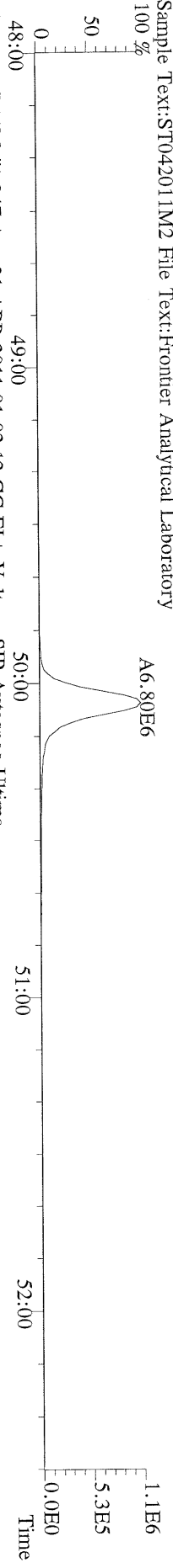
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419.8220 S:17 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



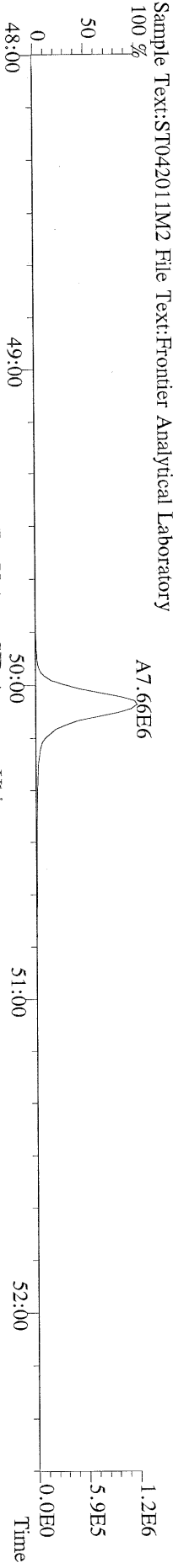
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479.7165 S:17 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,0%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory
100 %



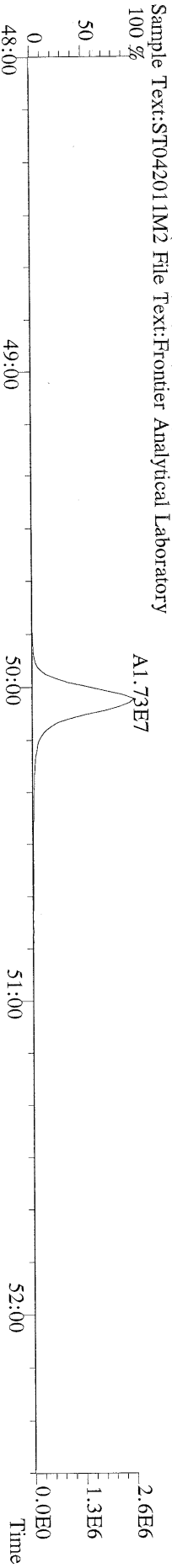
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441.7428 S:17 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



File:20APR11M #1-347 Acq:21-APR-2011 01:02:13 GC EI+ Voltage SIR Autospec-Ultima
443.7398 S:17 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory



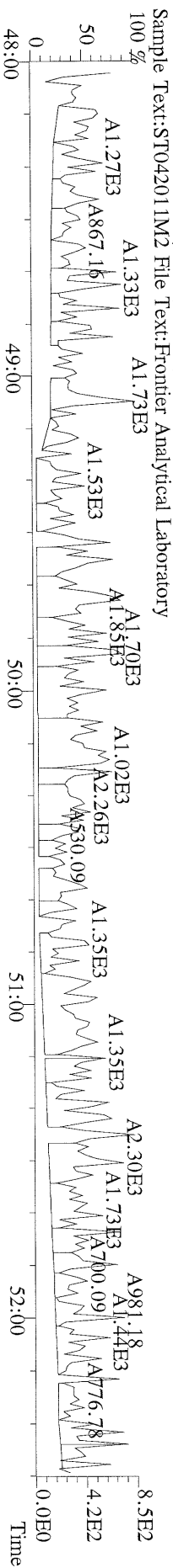
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453.7831 S:17 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory

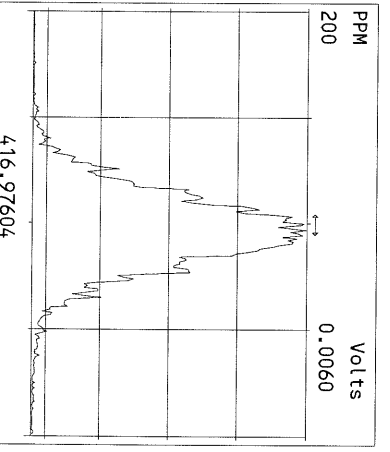
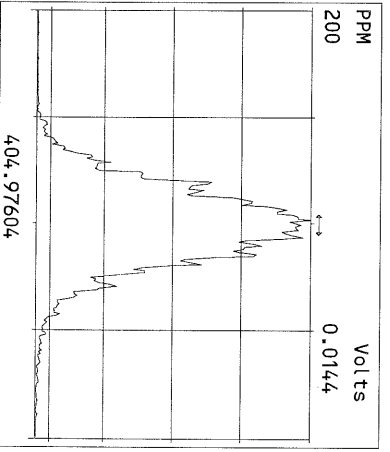
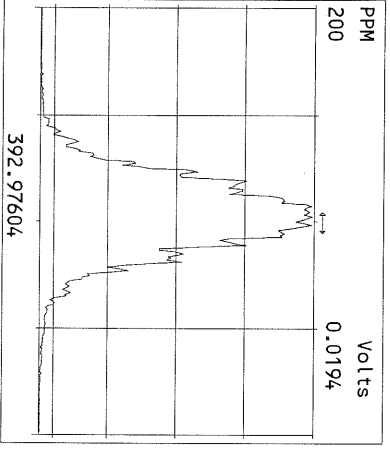
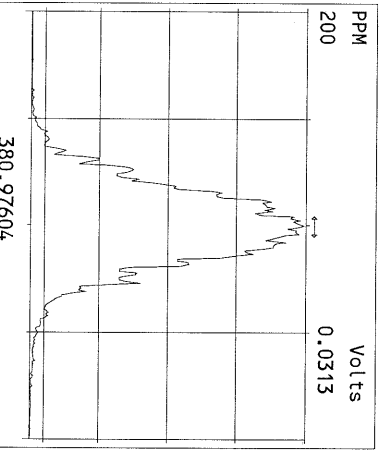
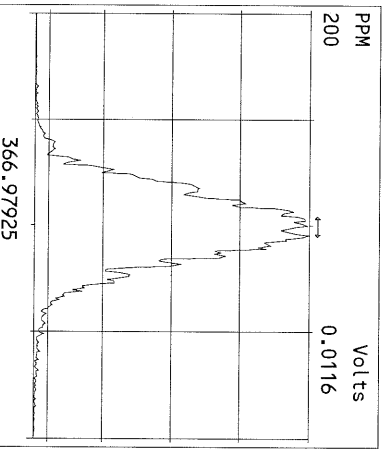
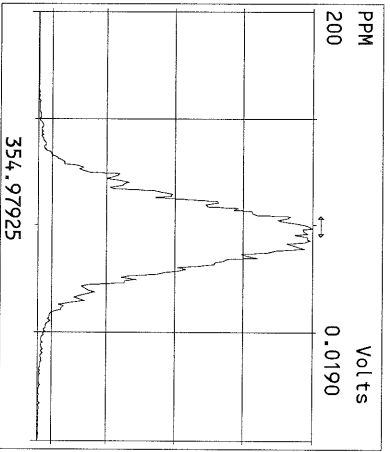
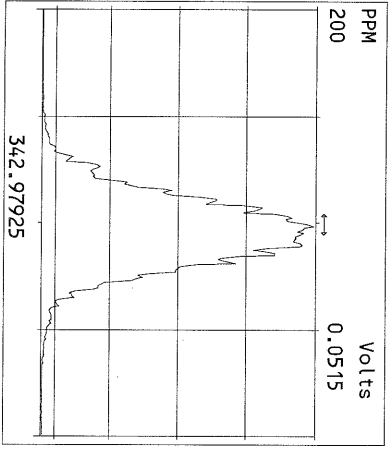
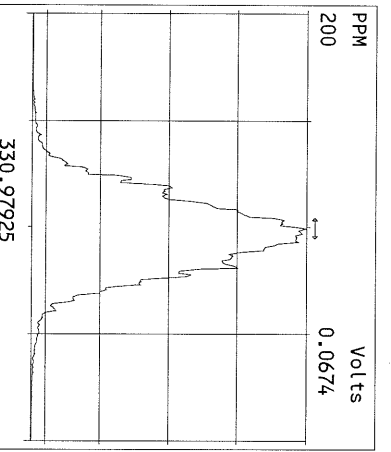
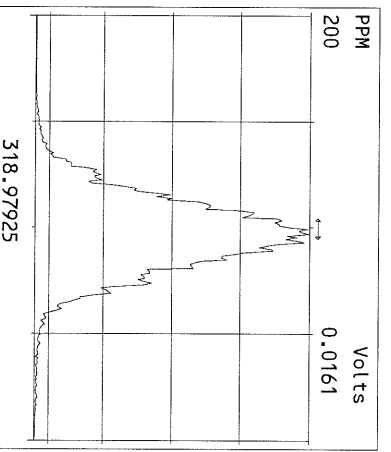
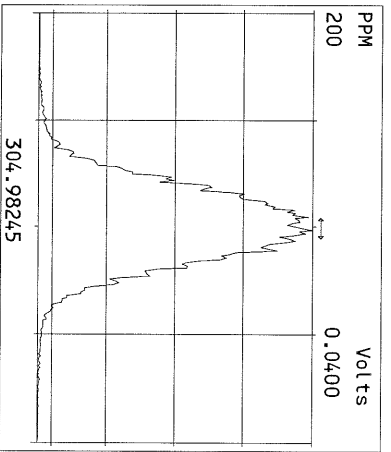
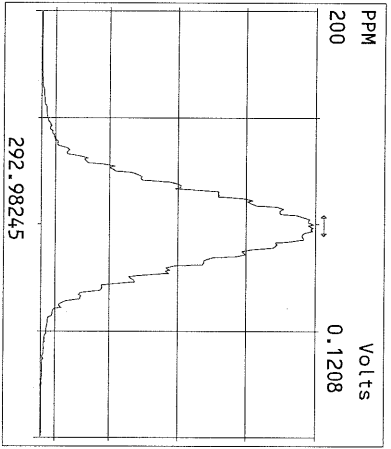


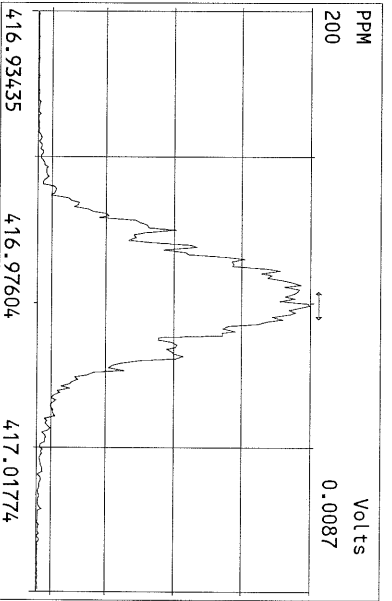
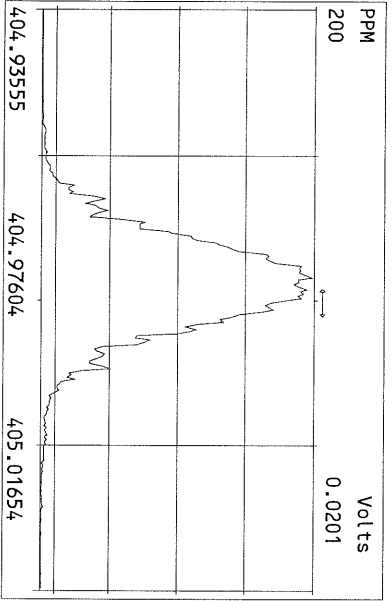
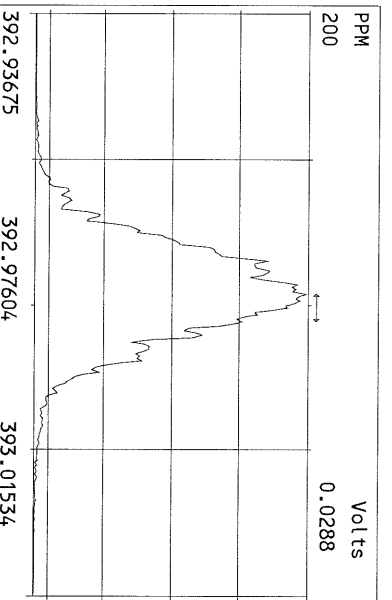
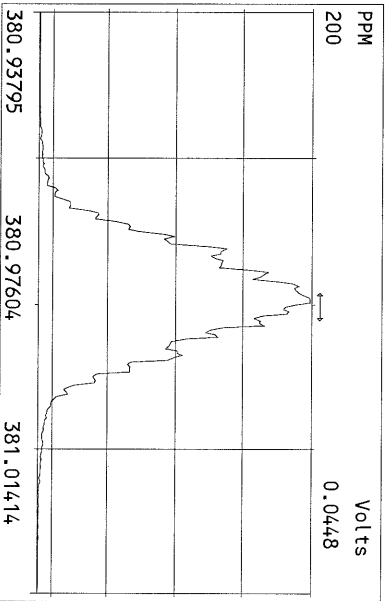
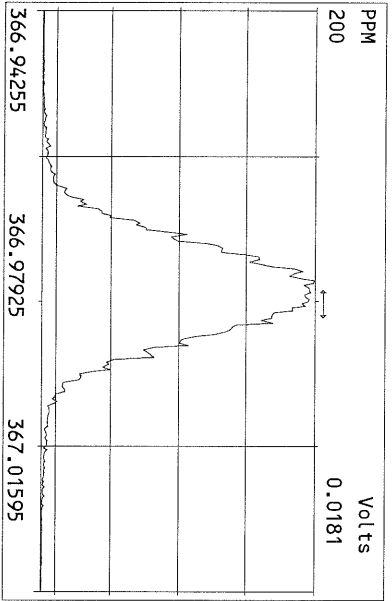
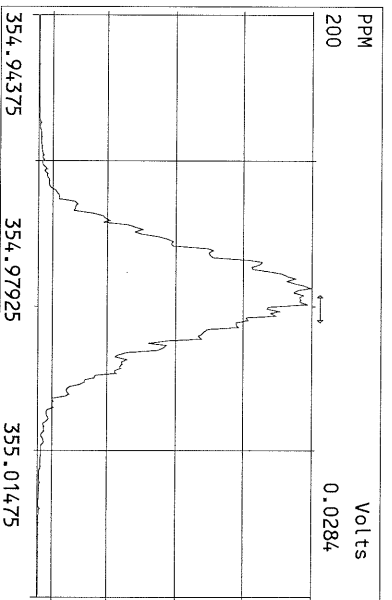
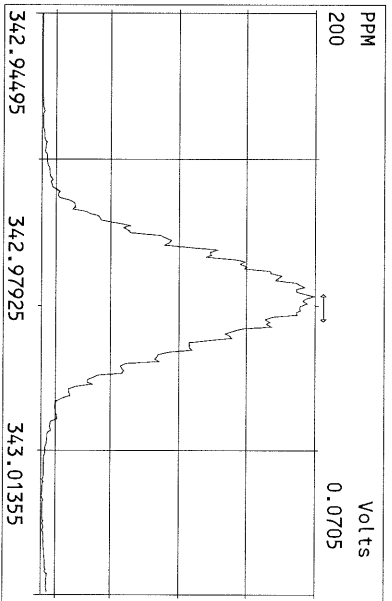
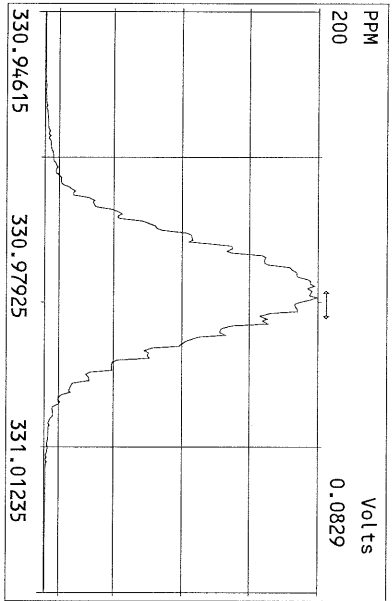
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455.7801 S:17 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory

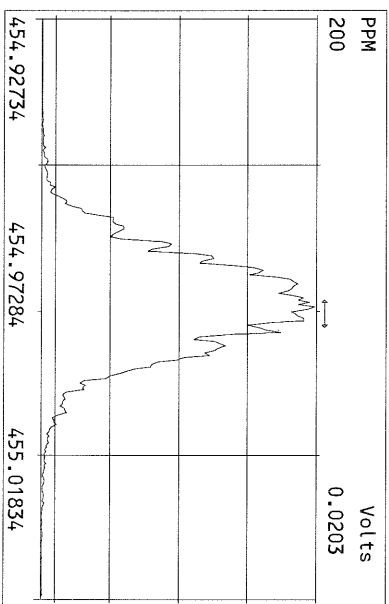
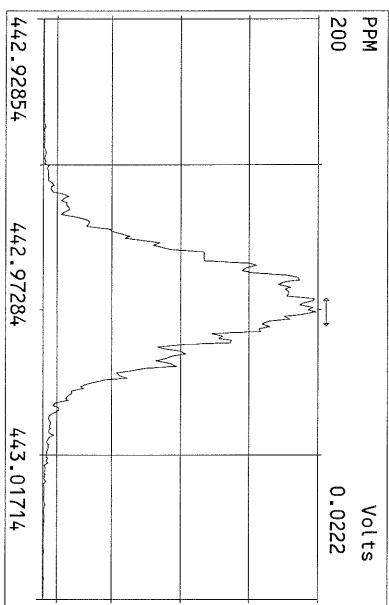
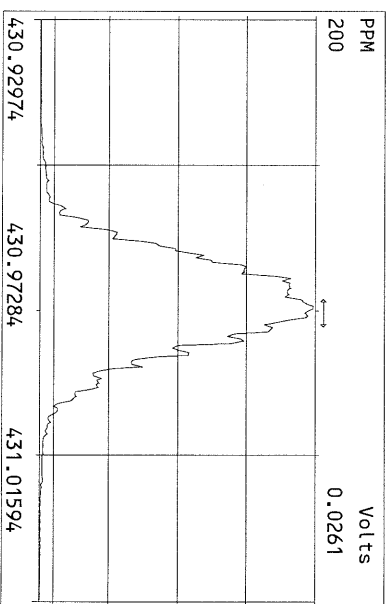
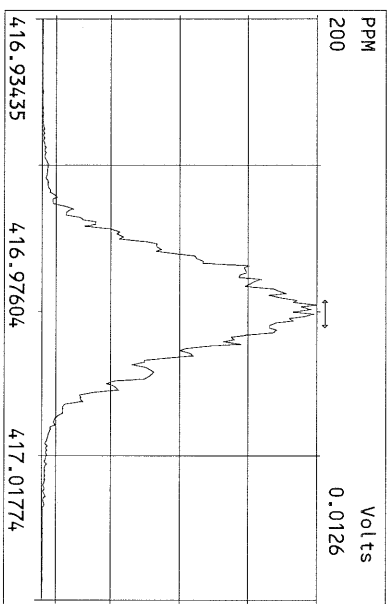
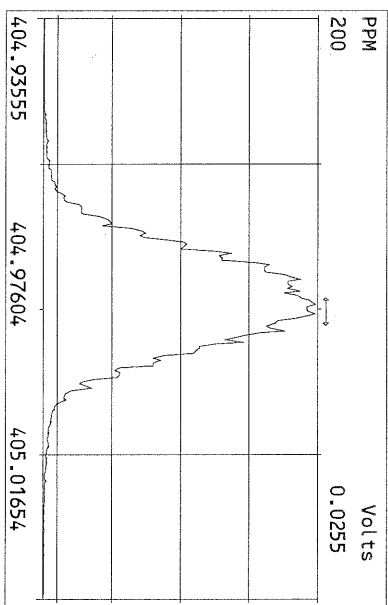
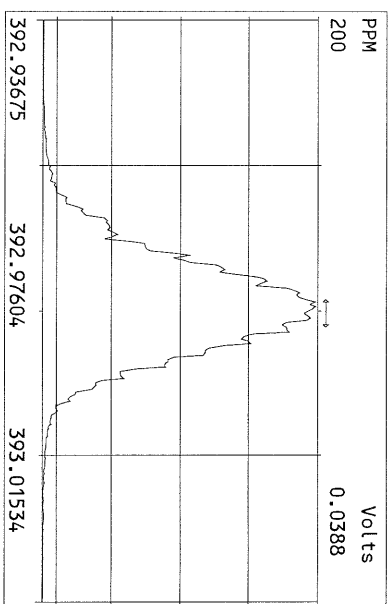
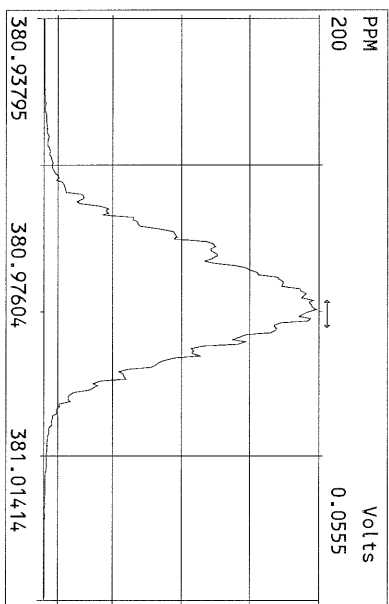
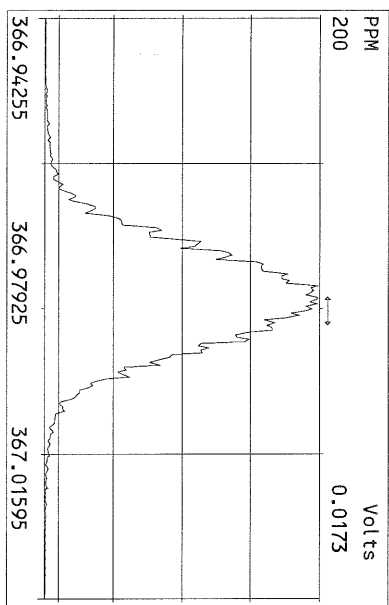


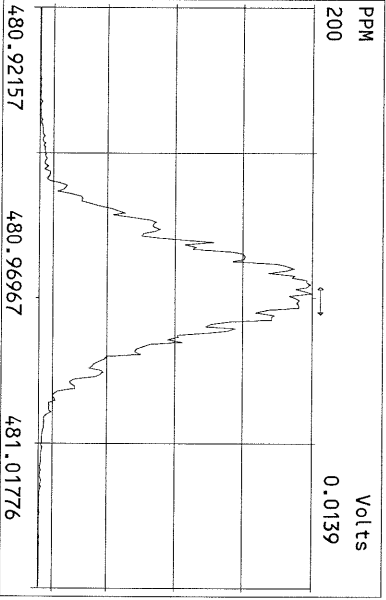
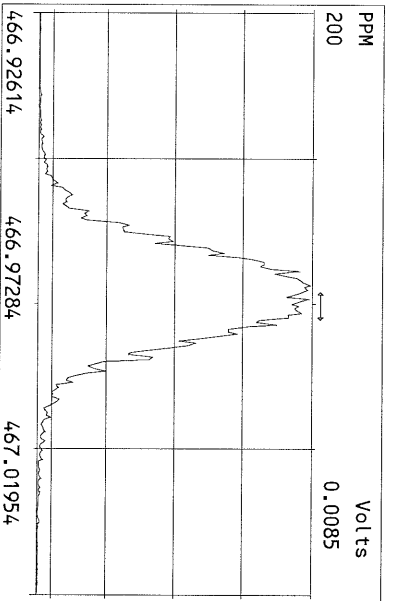
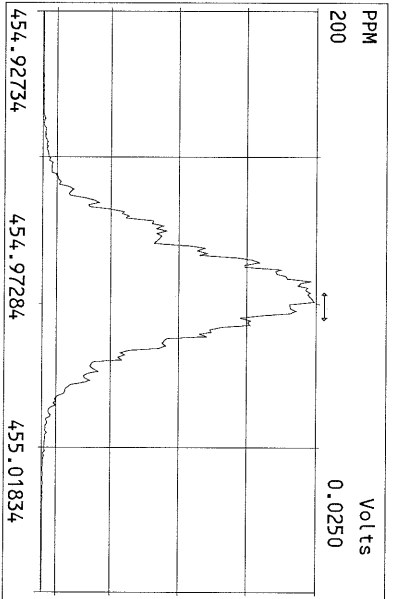
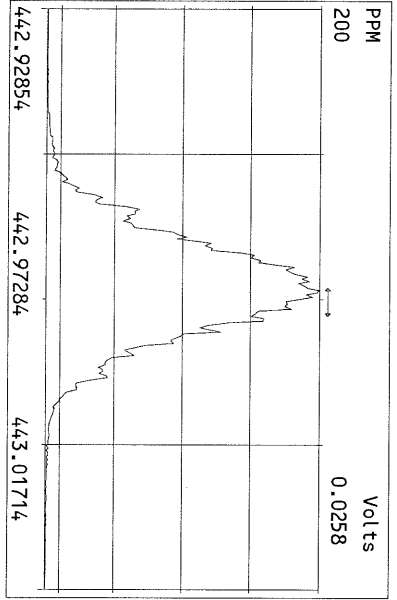
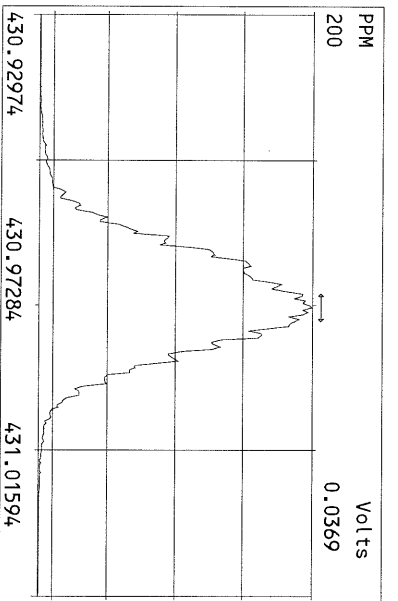
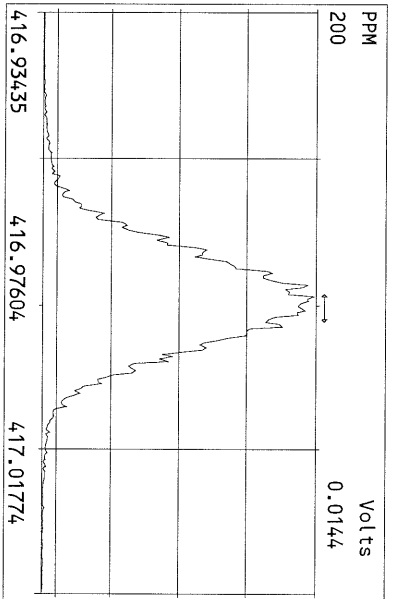
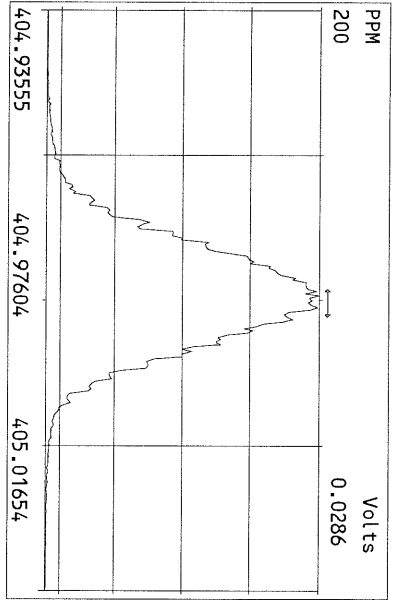
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513.6775 S:17 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M2 File Text:Frontier Analytical Laboratory

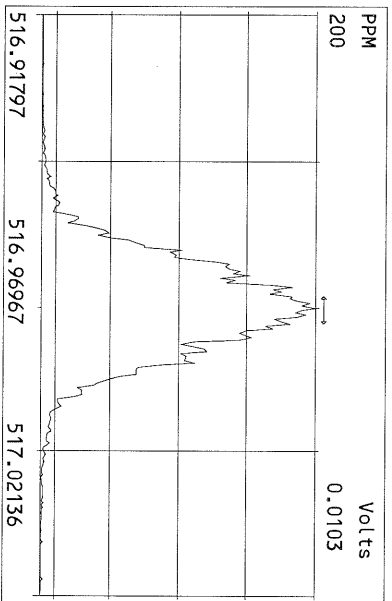
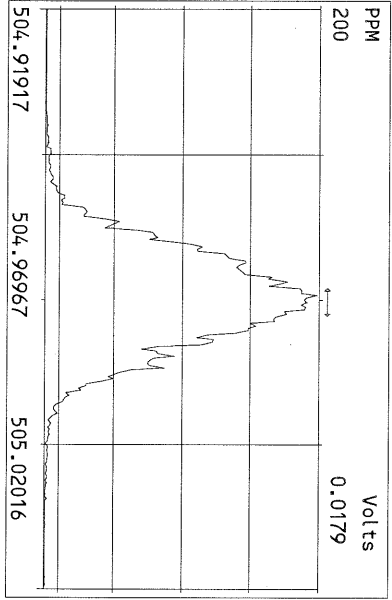
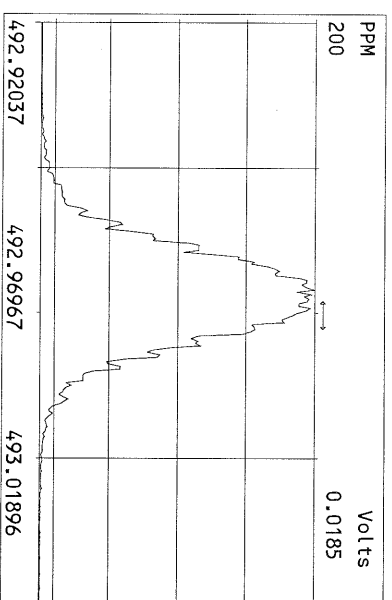
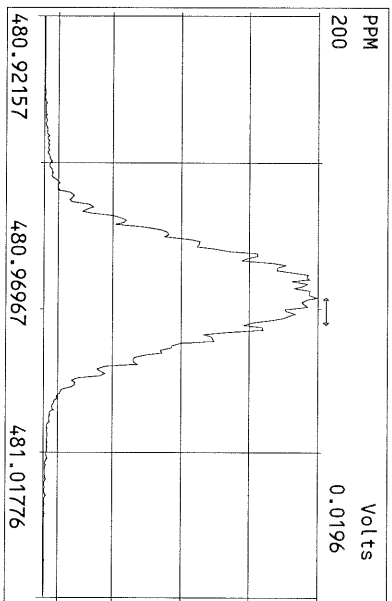
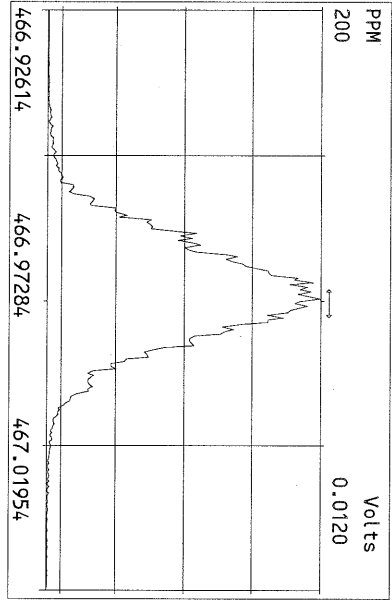
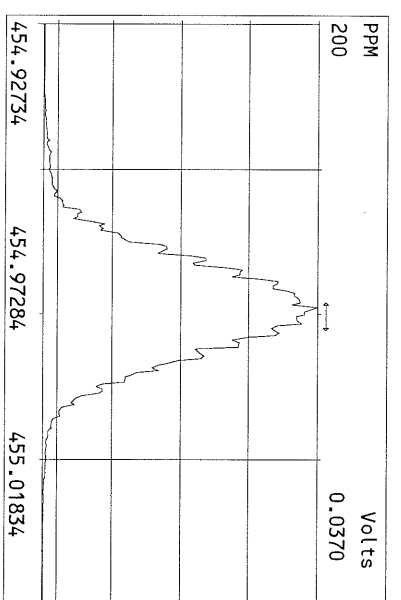
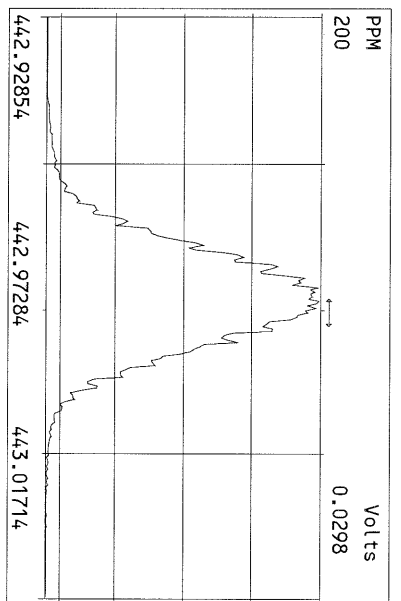
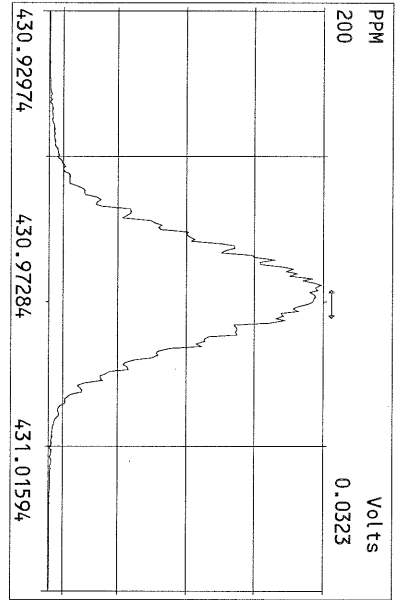












USEPA - ITD

FORM 4A

PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 20APR11M Sam:24

Analysis Date: 21-APR-11 07:29:49

NATIVE ANALYTES	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
2,3,7,8-TCDD	M/M+2	0.75	0.65-0.89	y	9.43	7.80 - 12.9 ✓
1,2,3,7,8-PeCDD	M+2/M+4	1.61	1.32-1.78	y	53.4	39.0 - 65.0 ✓
1,2,3,4,7,8-HxCDD	M+2/M+4	1.37	1.05-1.43	y	50.7	39.0 - 64.0 ✓
1,2,3,6,7,8-HxCDD	M+2/M+4	1.32	1.05-1.43	y	50.8	39.0 - 64.0 ✓
1,2,3,7,8,9-HxCDD	M+2/M+4	1.32	1.05-1.43	y	53.5	41.0 - 61.0 ✓
1,2,3,4,6,7,8-HpCDD	M+2/M+4	0.95	0.88-1.20	y	48.7	43.0 - 58.0 ✓
OCDD	M+2/M+4	0.93	0.76-1.02	y	99.5	79.0 - 126 ✓
2,3,7,8-TCDF	M/M+2	0.66	0.65-0.89	y	10.8	8.40 - 12.0 ✓
1,2,3,7,8-PeCDF	M+2/M+4	1.67	1.32-1.78	y	54.0	41.0 - 60.0 ✓
2,3,4,7,8-PeCDF	M+2/M+4	1.65	1.32-1.78	y	52.8	41.0 - 60.0 ✓
1,2,3,4,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	50.3	45.0 - 56.0 ✓
1,2,3,6,7,8-HxCDF	M+2/M+4	1.28	1.05-1.43	y	50.5	44.0 - 57.0 ✓
2,3,4,6,7,8-HxCDF	M+2/M+4	1.26	1.05-1.43	y	49.5	44.0 - 57.0 ✓
1,2,3,7,8,9-HxCDF	M+2/M+4	1.27	1.05-1.43	y	50.6	45.0 - 56.0 ✓
1,2,3,4,6,7,8-HpCDF	M+2/M+4	1.04	0.88-1.20	y	48.6	45.0 - 55.0 ✓
1,2,3,4,7,8,9-HpCDF	M+2/M+4	1.05	0.88-1.20	y	48.4	43.0 - 58.0 ✓
OCDF	M+2/M+4	0.90	0.76-1.02	y	95.6	63.0 - 159 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

Analyst: JDate: 4/21/11

USEPA - ITD

FORM 4B
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Initial Calibration Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 20APR11M Sam:24

Analysis Date: 21-APR-11 07:29:49

LABELED COMPOUNDS	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
13C-2,3,7,8-TCDD	M/M+2	0.82	0.65-0.89	y	96.6	82.0 - 121 ✓
13C-1,2,3,7,8-PeCDD	M+2/M+4	1.78	1.32-1.78	y	88.7	62.0 - 160 ✓
13C-1,2,3,4,7,8-HxCDD	M+2/M+4	1.30	1.05-1.43	y	92.6	85.0 - 117 ✓
13C-1,2,3,6,7,8-HxCDD	M+2/M+4	1.29	1.05-1.43	y	96.7	85.0 - 118 ✓
13C-1,2,3,4,6,7,8-HpCDD	M+2/M+4	1.05	0.88-1.20	y	125	72.0 - 138 ✓
13C-OCDD	M+2/M+4	0.96	0.76-1.02	y	214	96.0 - 415 ✓
13C-2,3,7,8-TCDF	M/M+2	0.88	0.65-0.89	y	101	71.0 - 140 ✓
13C-1,2,3,7,8-PeCDF	M+2/M+4	1.72	1.32-1.78	y	105	76.0 - 130 ✓
13C-2,3,4,7,8-PeCDF	M+2/M+4	1.73	1.32-1.78	y	105	77.0 - 130 ✓
13C-1,2,3,4,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	96.4	76.0 - 131 ✓
13C-1,2,3,6,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	96.2	70.0 - 143 ✓
13C-2,3,4,6,7,8-HxCDF	M/M+2	0.48	0.43-0.59	y	102	73.0 - 137 ✓
13C-1,2,3,7,8,9-HxCDF	M/M+2	0.47	0.43-0.59	y	105	74.0 - 135 ✓
13C-1,2,3,4,6,7,8-HpCDF	M/M+2	0.45	0.37-0.51	y	116	78.0 - 129 ✓
13C-1,2,3,4,7,8,9-HpCDF	M/M+2	0.43	0.37-0.51	y	128	77.0 - 129 ✓
13C-OCDF	M+2/M+4	0.94	0.76-1.02	y	211	96.0 - 415 ✓
CLEANUP STANDARD (4)						
37Cl-2,3,7,8-TCDD					8.11	7.80 - 12.8 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.

(3) Contract-required concentration range as specified in Table 6, Method 1613.

(4) No ion abundance ratio; report concentration found.

Analyst: JDate: 4/21/11

FORM 5
PCDD/PCDF RT WINDOW AND ISOMER SPECIFICITY STANDARDS

Lab Name: Frontier Analytical Laboratory Episode No.:
Contract No.: SAS No.:
Instrument ID: FAL3 Initial Calibration Date: 3/7/11
RT Window Data Filename: 20APR11M Sam:24 Analysis Date: 21-APR-11 Time: 07:29:49
DB-5 IS Data Filename: 20APR11M Sam:24 Analysis Date: 21-APR-11 Time: 07:29:49
DB-225 IS Date Filename: Analysis Date: Time:

DB-5 RT WINDOW DEFINING STANDARDS RESULTS

ISOMERS	ABSOLUTE RT	ISOMERS	ABSOLUTE RT
1,3,6,8-TCDD (F)	24:23 ✓	1,3,6,8-TCDF (F)	23:03 ✓
1,2,8,9-TCDD (L)	28:17 ✓	1,2,8,9-TCDF (L)	28:30 ✓
1,2,4,7,9-PeCDD (F)	30:12 ✓	1,3,4,6,8-PeCDF (F)	28:23 ✓
1,2,3,8,9-PeCDD (L)	33:45 ✓	1,2,3,8,9-PeCDF (L)	34:10 ✓
1,2,4,6,7,9-HxCDD (F)	36:05 ✓	1,2,3,4,6,8-HxCDF (F)	35:13 ✓
1,2,3,7,8,9-HxCDD (L)	39:09 ✓	1,2,3,7,8,9-HxCDF (L)	39:43 ✓
1,2,3,4,6,7,9-HpCDD (F)	42:46 ✓	1,2,3,4,6,7,8-HpCDF (F)	42:14 ✓
1,2,3,4,6,7,8-HpCDD (L)	44:09 ✓	1,2,3,4,7,8,9-HpCDF (L)	45:03 ✓

(F) = First eluting iosmer (DB-5); (L) = Last eluting isomer (DB-5)

ISOMER SPECIFICITY (IS) TEST STANDARD RESULTS

% VALLEY HEIGHT
BETWEEN
COMPARED PEAKS (1)

<25%

(1) To meet contract requirement, %Valley Height Between Compared Peaks shall not exceed 25% (section 15.4.2.2, Method 1613).

Analyst: 

Date: 4/21/11

USEPA - ITD

FORM 6A

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 21-APR-11 07:29:49

CS3 or VER Data Filename: 20APR11M

Sam:24

NATIVE ANALYTES	RETENTION TIME		RRT	RRT	QC LIMITS (1)
	REFERENCE				
2,3,7,8-TCDD	13C-2,3,7,8-TCDD		1.001		0.999-1.002 ✓
2,3,7,8-TCDF	13C-2,3,7,8-TCDF		1.001		0.999-1.003 ✓
1,2,3,7,8-PeCDD	13C-1,2,3,7,8-PeCDD		1.001		0.999-1.002 ✓
1,2,3,7,8-PeCDF	13C-1,2,3,7,8-PeCDF		1.001		0.999-1.002 ✓
2,3,4,7,8-PeCDF	13C-2,3,4,7,8-PeCDF		1.000		0.999-1.002 ✓
LABELED COMPOUNDS					
37Cl-2,3,7,8-TCDD	13C-1,2,3,4-TCDD		1.022		0.989-1.052 ✓
13C-2,3,7,8-TCDD			1.021		0.976-1.043 ✓
13C-2,3,7,8-TCDF			0.993		0.923-1.103 ✓
13C-1,2,3,7,8-PeCDD			1.238		1.000-1.567 ✓
13C-1,2,3,7,8-PeCDF			1.173		0.923-1.203 ✓
13C-2,3,4,7,8-PeCDF			1.223		0.923-1.303 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified in Table 2, Method 1613.

Analyst: J

Date: 4/21/11

USEPA - ITD

FORM 6B

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date: 3/7/11

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 21-APR-11 07:29:49

CS3 or VER Data Filename: 20APR11M

Sam:24

NATIVE ANALYTES	RETENTION TIME		RRT	QC LIMITS (1)
	REFERENCE	RETENTION TIME		
1,2,3,4,7,8-HxCDD	13C-1,2,3,4,7,8-HxCDD		1.000	0.999-1.001 ✓
1,2,3,6,7,8-HxCDD	13C-1,2,3,6,7,8-HxCDD		1.000	0.998-1.004 ✓
1,2,3,7,8,9-HxCDD	13C-1,2,3,6,7,8-HxCDD		1.012	1.000-1.019 ✓
1,2,3,4,7,8-HxCDF	13C-1,2,3,4,7,8-HxCDF		1.000	0.999-1.001 ✓
1,2,3,6,7,8-HxCDF	13C-1,2,3,6,7,8-HxCDF		1.001	0.997-1.005 ✓
2,3,4,6,7,8-HxCDF	13C-2,3,4,6,7,8-HxCDF		1.001	0.999-1.001 ✓
1,2,3,7,8,9-HxCDF	13C-1,2,3,7,8,9-HxCDF		1.001	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDD	13C-1,2,3,4,6,7,8-HpCDD		1.000	0.999-1.001 ✓
1,2,3,4,6,7,8-HpCDF	13C-1,2,3,4,6,7,8-HpCDF		1.000	0.999-1.001 ✓
1,2,3,4,7,8,9-HpCDF	13C-1,2,3,4,7,8,9-HpCDF		1.001	0.999-1.001 ✓
OCDD	13C-OCDD		1.000	0.999-1.001 ✓
OCDF	13C-OCDF		1.001	0.999-1.001 ✓
LABELED COMPOUNDS				
13C-1,2,3,4,7,8-HxCDD	13C-1,2,3,7,8,9-HxCDD		0.985	0.977-1.000 ✓
13C-1,2,3,6,7,8-HxCDD			0.989	0.981-1.003 ✓
13C-1,2,3,4,7,8-HxCDF			0.949	0.944-0.970 ✓
13C-1,2,3,6,7,8-HxCDF			0.954	0.949-0.975 ✓
13C-2,3,4,6,7,8-HxCDF			0.978	0.959-1.021 ✓
13C-1,2,3,7,8,9-HxCDF			1.015	0.977-1.047 ✓
13C-1,2,3,4,6,7,8-HpCDD			1.128	1.086-1.130 ✓
13C-1,2,3,4,6,7,8-HpCDF			1.079	1.043-1.085 ✓
13C-1,2,3,4,7,8,9-HpCDF			1.151	1.057-1.154 ✓
13C-OCDD			1.270	1.032-1.311 ✓
13C-OCDF			1.279	1.000-1.311 ✓

(1) Contract-required limits for Relative Retention Times (RRT) as specified
in Table 2, Method 1613.

Analyst: JDate: 4/21/11

Name	Resp	RA	RT	RRF	Conc	Qual	Fac Noise-1	Noise-2	DL	#Hom
2,3,7,8-TCDD	2.03e+06	0.75 y	27:22	1.13	9.43		2.50	-	-	*
1,2,3,7,8-PeCDD	9.28e+06	1.61 y	33:10	1.02	53.4		2.50	-	-	*
1,2,3,4,7,8-HxCDD	1.17e+07	1.37 y	38:32	1.45	50.7		2.50	-	-	*
1,2,3,6,7,8-HxCDD	9.65e+06	1.32 y	38:42	1.45	50.8		2.50	-	-	*
1,2,3,7,8,9-HxCDD	1.14e+07	1.32 y	39:09	1.47	53.5		2.50	-	-	*
1,2,3,4,6,7,8-HpCDD	9.49e+06	0.95 y	44:09	1.30	48.7		2.50	-	-	*
OCDD	1.35e+07	0.93 y	49:41	1.45	99.5		2.50	-	-	*
2,3,7,8-TCDF	4.07e+06	0.66 y	26:36	1.15	10.8		2.50	-	-	*
1,2,3,7,8-PeCDF	1.38e+07	1.67 y	31:26	0.89	54.0		2.50	-	-	*
2,3,4,7,8-PeCDF	1.32e+07	1.65 y	32:45	0.89	52.8		2.50	-	-	*
1,2,3,4,7,8-HxCDF	1.30e+07	1.26 y	37:09	1.01	50.3		2.50	-	-	*
1,2,3,6,7,8-HxCDF	1.43e+07	1.28 y	37:21	0.89	50.5		2.50	-	-	*
2,3,4,6,7,8-HxCDF	1.38e+07	1.26 y	38:17	1.02	49.5		2.50	-	-	*
1,2,3,7,8,9-HxCDF	1.67e+07	1.27 y	39:43	1.10	50.6		2.50	-	-	*
1,2,3,4,6,7,8-HpCDF	1.19e+07	1.04 y	42:14	1.48	48.6		2.50	-	-	*
1,2,3,4,7,8,9-HpCDF	9.74e+06	1.05 y	45:03	1.43	48.4		2.50	-	-	*
OCDF	1.43e+07	0.90 y	50:04	0.84	95.6		2.50	-	-	*
13C-2,3,7,8-TCDD	1.90e+07	0.82 y	27:20	1.03	96.6					Rec 96.6
13C-1,2,3,7,8-PeCDD	1.71e+07	1.78 y	33:09	1.01	88.7					88.7
13C-1,2,3,4,7,8-HxCDD	1.59e+07	1.30 y	38:31	1.19	92.6					92.6
13C-1,2,3,6,7,8-HxCDD	1.31e+07	1.29 y	38:41	0.94	96.7					96.7
13C-1,2,3,4,6,7,8-HpCDD	1.50e+07	1.05 y	44:08	0.83	125					125
13C-OCDD	1.87e+07	0.96 y	49:40	0.61	214					107
13C-2,3,7,8-TCDF	3.28e+07	0.88 y	26:35	0.98	101					101
13C-1,2,3,7,8-PeCDF	2.89e+07	1.72 y	31:25	0.83	105					105
13C-2,3,4,7,8-PeCDF	2.79e+07	1.73 y	32:44	0.80	105					105
13C-1,2,3,4,7,8-HxCDF	2.56e+07	0.48 y	37:08	1.84	96.4					96.4
13C-1,2,3,6,7,8-HxCDF	3.18e+07	0.48 y	37:20	2.29	96.2					96.2
13C-2,3,4,6,7,8-HxCDF	2.73e+07	0.48 y	38:15	1.86	102					102
13C-1,2,3,7,8,9-HxCDF	3.00e+07	0.47 y	39:41	1.98	105					105
13C-1,2,3,4,6,7,8-HpCDF	1.65e+07	0.45 y	42:13	0.99	116					116
13C-1,2,3,4,7,8,9-HpCDF	1.41e+07	0.43 y	45:01	0.77	128					128
13C-OCDF	3.55e+07	0.94 y	50:02	1.17	211					105
37Cl-2,3,7,8-TCDD	1.13e+06		27:22	0.73	8.11					81.1
13C-1,2,3,4-TCDD	1.91e+07	0.83 y	26:46	-	50.2					
13C-1,2,3,4-TCDF	3.31e+07	0.88 y	25:31	-	46.0					
13C-1,2,3,7,8,9-HxCDD	1.44e+07	1.29 y	39:07	-	58.2					
Total Tetra-Dioxins	1.07e+07		22:38	1.13	49.9		2.50	-	-	* 25
Total Penta-Dioxins	1.96e+07		30:12	1.02	113		2.50	-	-	* 4
Total Hexa-Dioxins	3.65e+07		36:05	1.46	173		2.50	-	-	* 7
Total Hepta-Dioxins	1.94e+07		42:46	1.30	99.4		2.50	-	-	* 11
Total Tetra-Furans	1.88e+07		23:03	1.15	49.8		2.50	-	-	* 18
1st Fn. Tot Penta-Furans	1.36e+07		28:23	0.89	54.0		2.50	-	-	* PeCDF 1
Total Penta-Furans	3.89e+07		30:07	0.89	154		2.50	-	-	* 208 8
Total Hexa-Furans	6.61e+07		35:13	1.00	230		2.50	-	-	* 15
Total Hepta-Furans	2.16e+07		42:14	1.46	97.1		2.50	-	-	* 3

Analyst: J

Date: 4/21/11

Frontier Analytical Laboratory - Acquisition Log

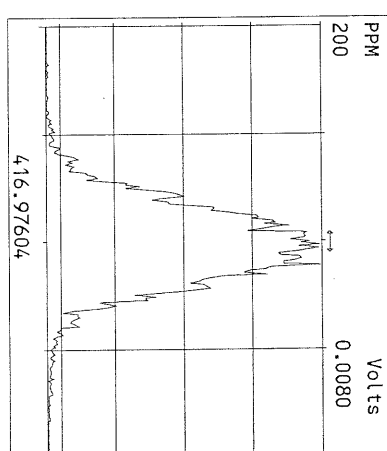
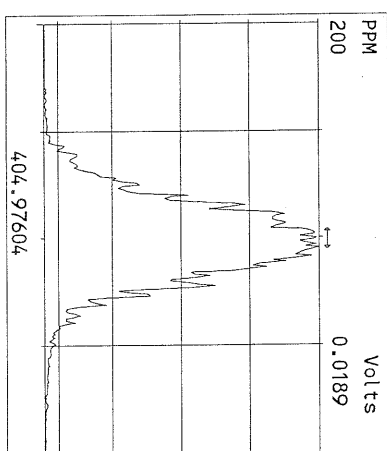
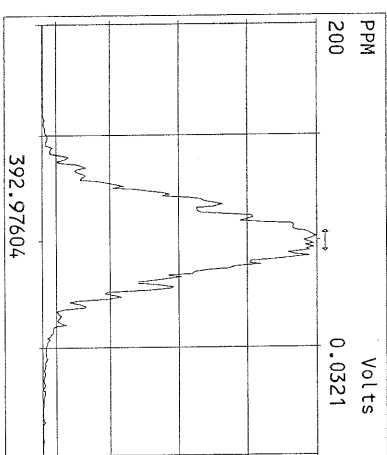
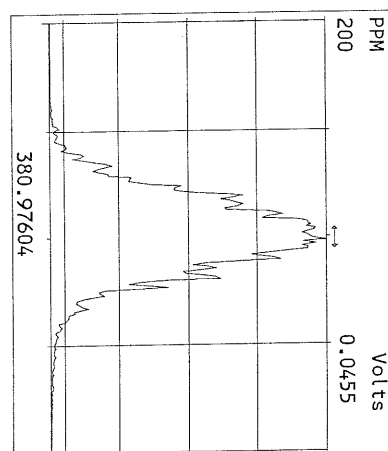
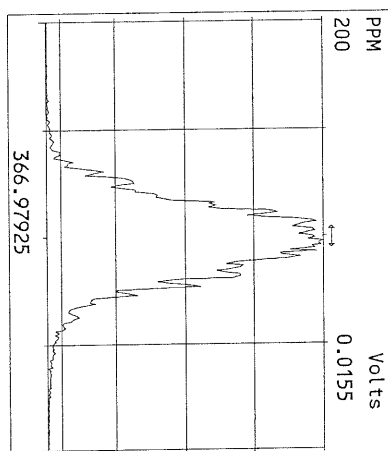
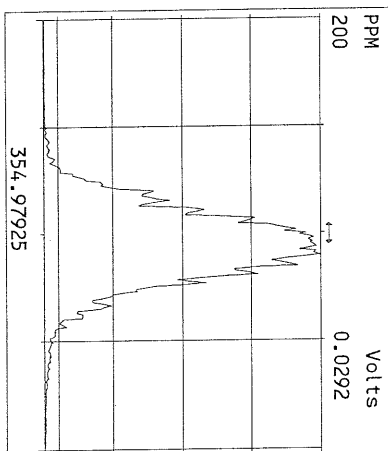
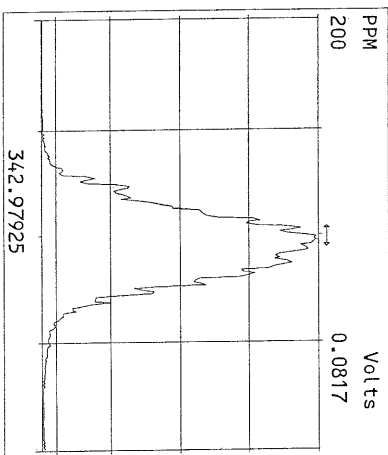
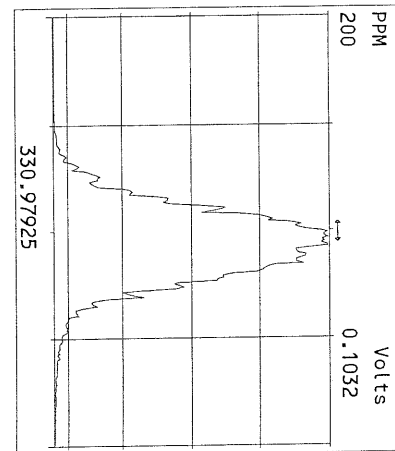
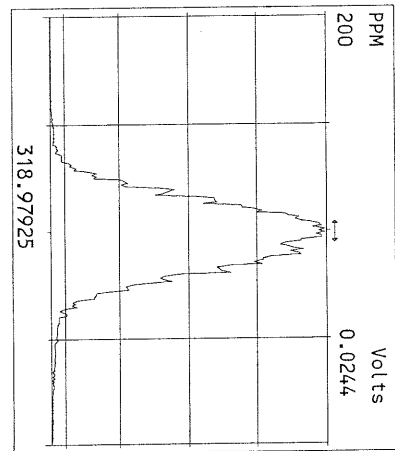
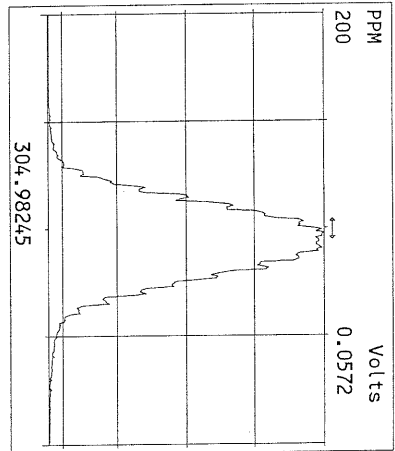
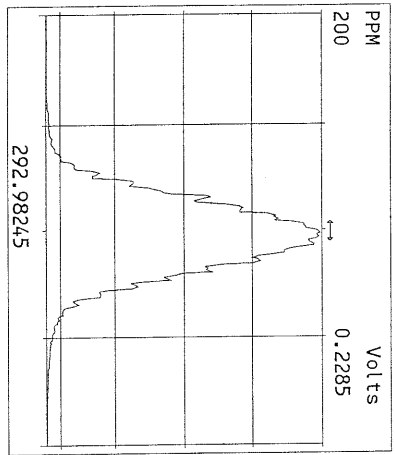
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Data File	S	FAL ID	Client ID	Acquired	ConCal	EndCal	Analyst
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20APR11M	2	2272-001-0001-OPR	OPR	20-APR-11 11:11:39	ST042011M1	ST042011M2	TC
20APR11M	3	2272-001-0001-MB	Method Blank	20-APR-11 12:07:02	ST042011M1	ST042011M2	TC
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20APR11M	6	6701-009-0001-SA	MC-SED3-0-10-032911	20-APR-11 14:53:07	ST042011M1	ST042011M2	TC
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20APR11M	14	6701-002-0001-MSD	LL-SED2-0-15-032911	20-APR-11 22:15:58	ST042011M1	ST042011M2	TC
20APR11M	15	SB042011M1	Solvent Blank	20-APR-11 23:11:21	ST042011M1	ST042011M2	TC
20APR11M	16	SB042011M2	Solvent Blank	21-APR-11 00:06:50	ST042011M1	ST042011M2	TC
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20APR11M	21	6701-001-0001-SA	LL-SED1-0-15-032911	21-APR-11 04:43:45	ST042011M2	ST042011M3	TC
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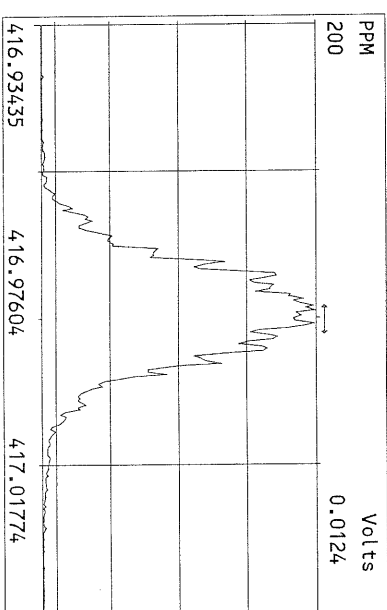
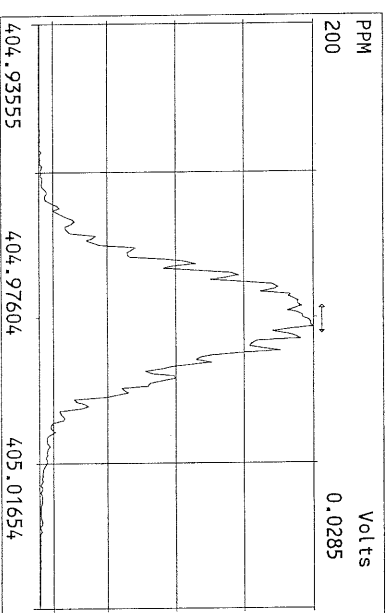
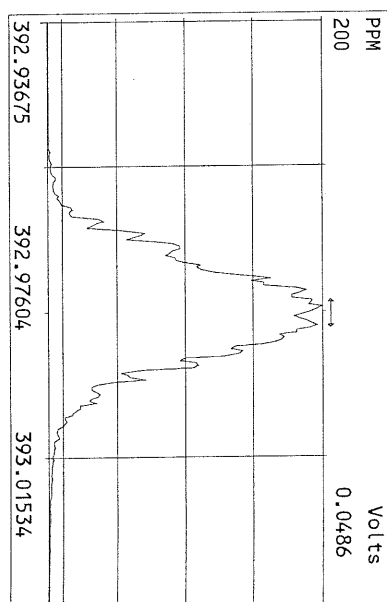
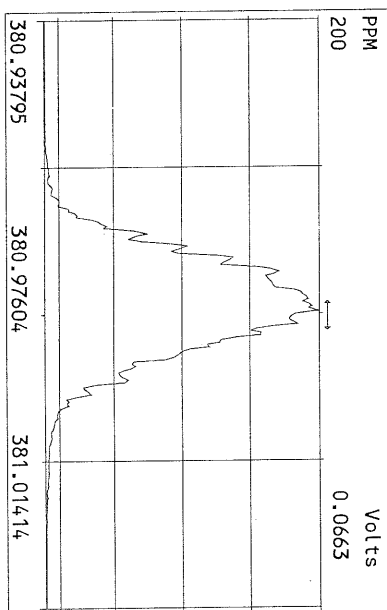
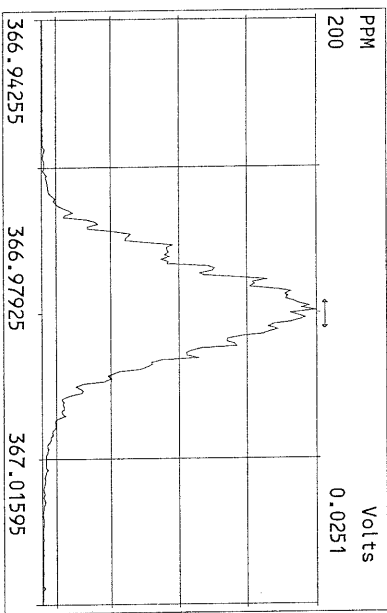
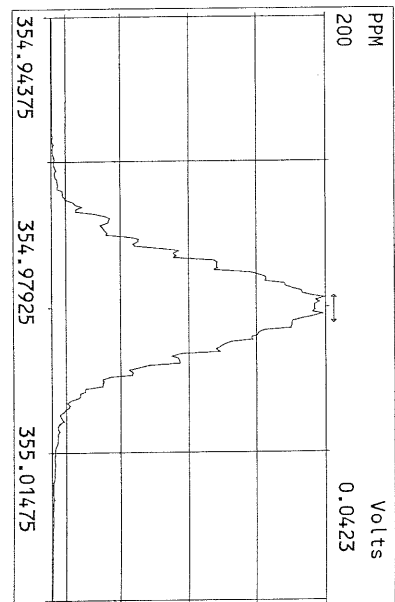
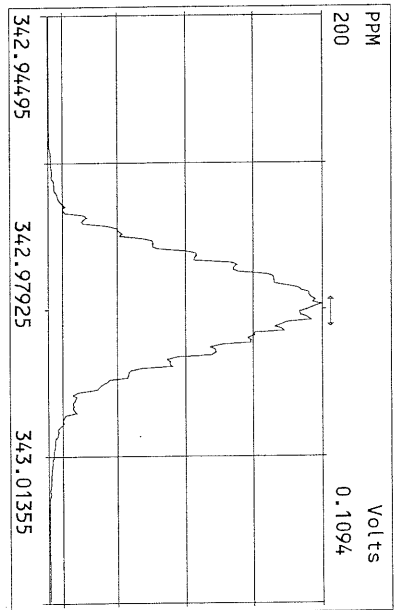
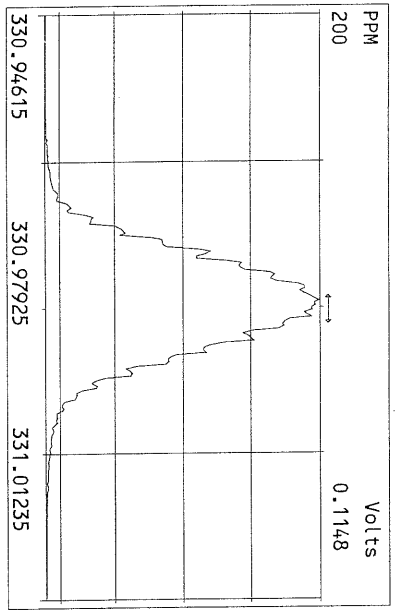
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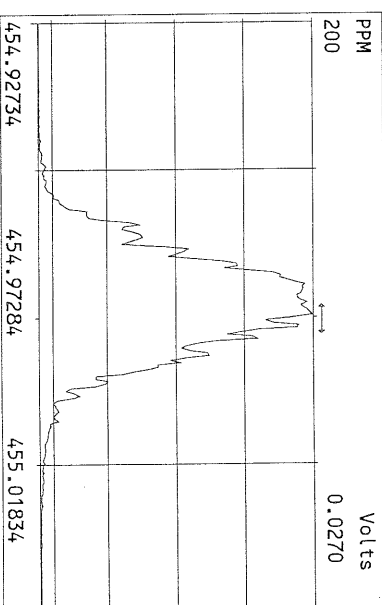
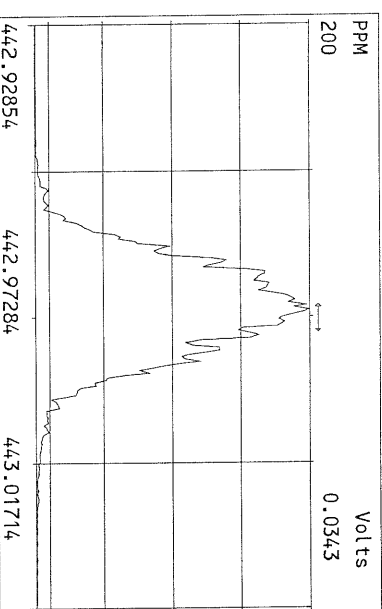
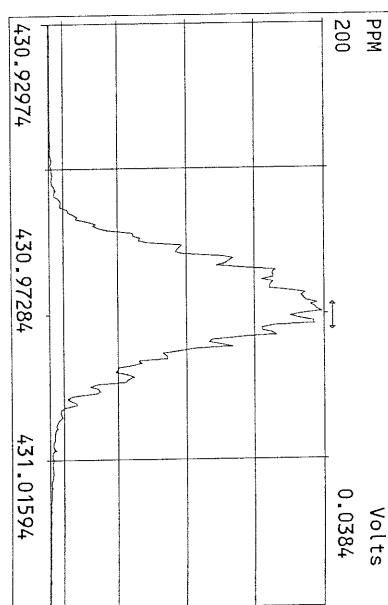
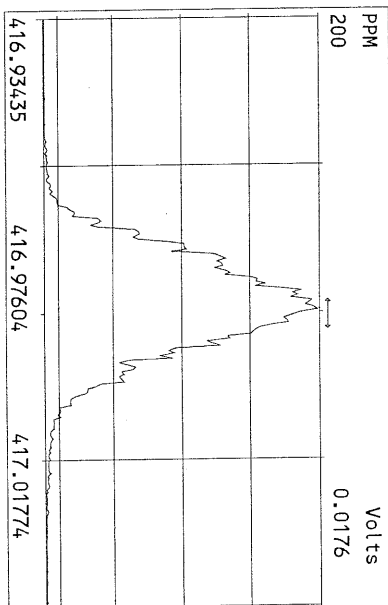
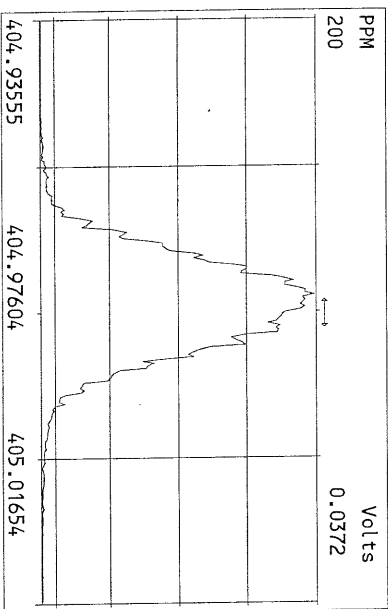
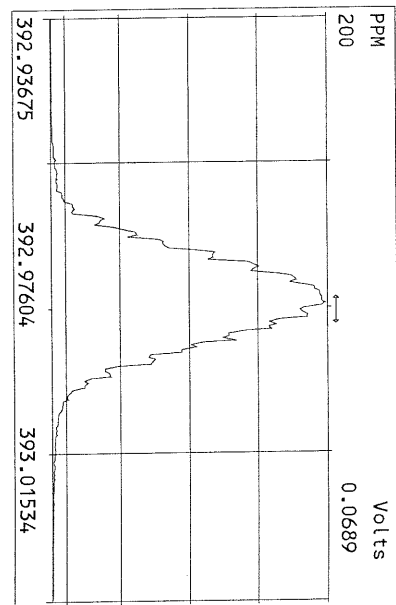
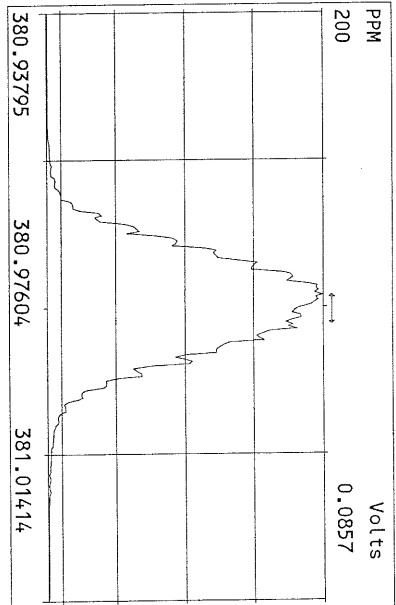
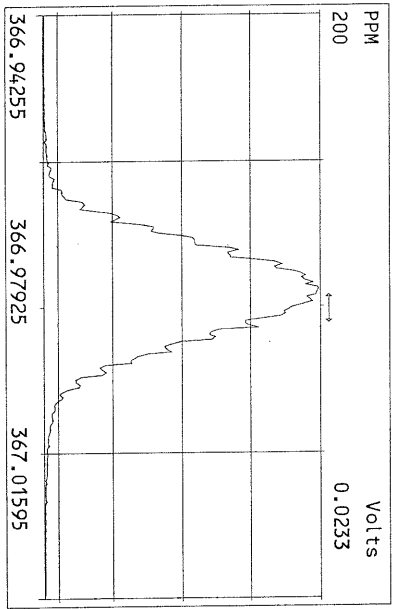
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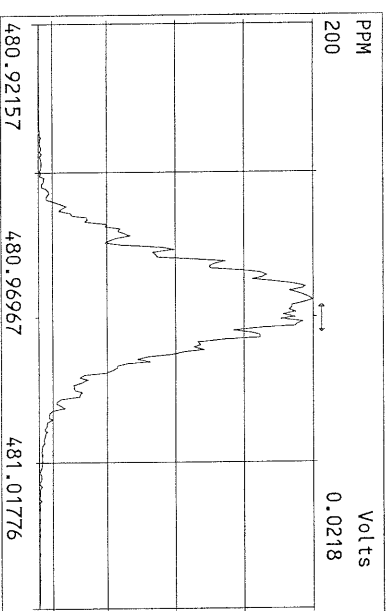
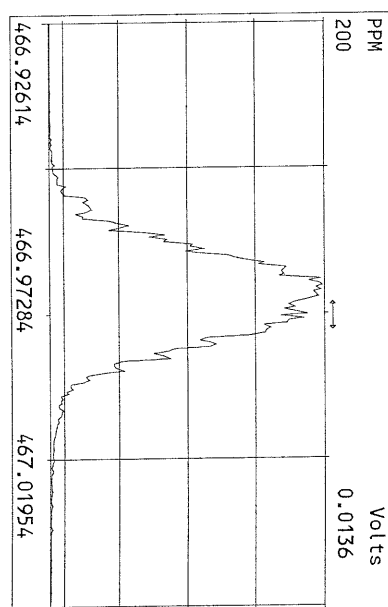
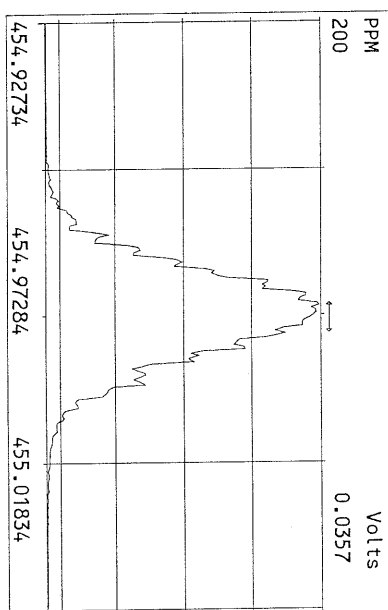
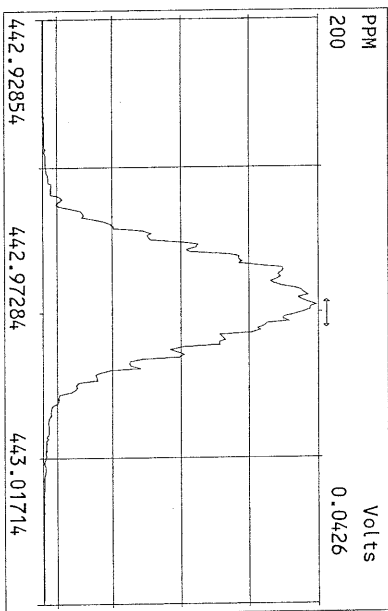
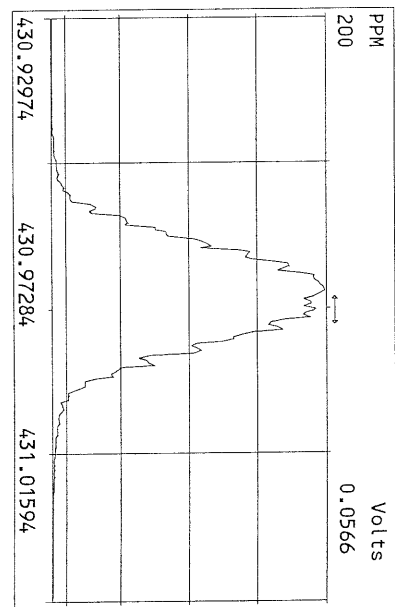
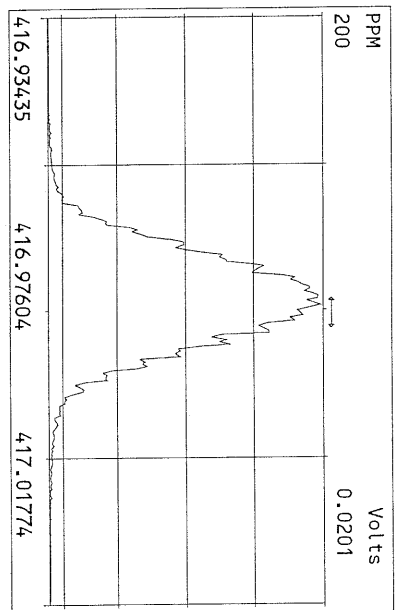
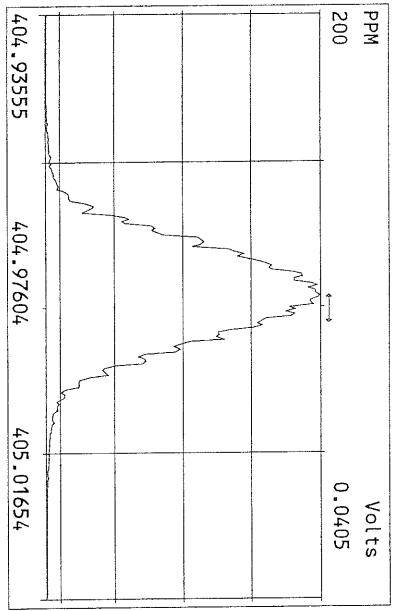
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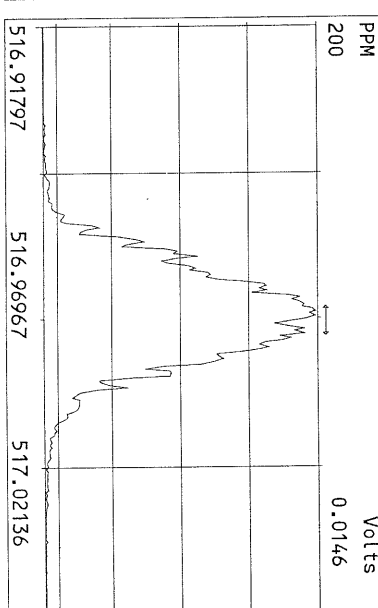
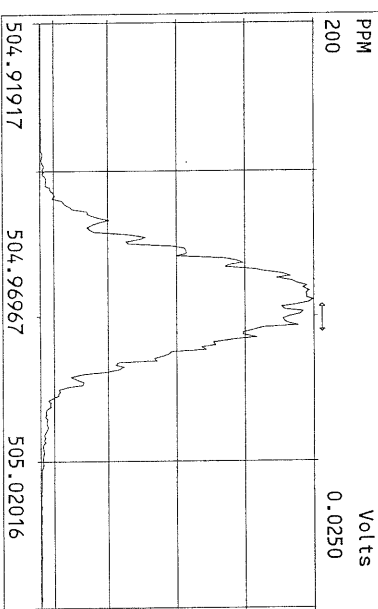
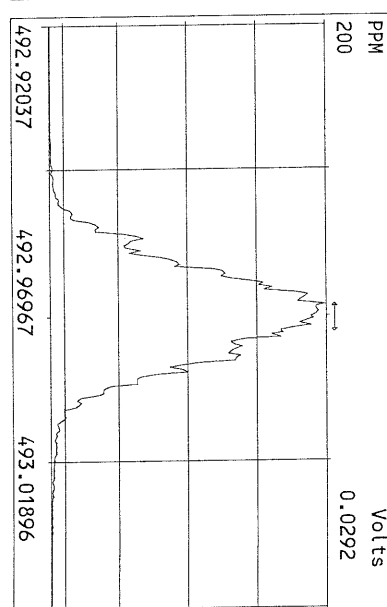
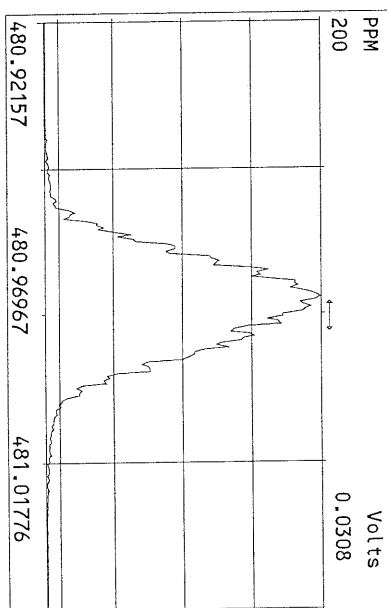
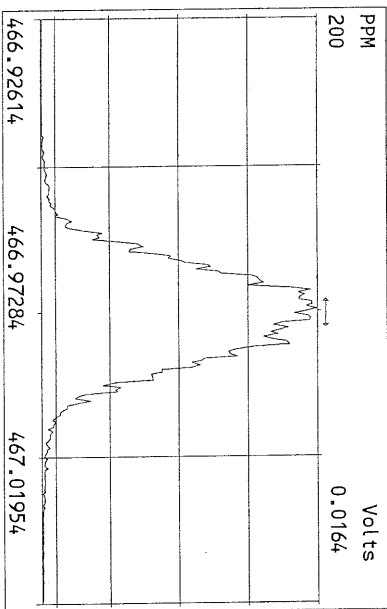
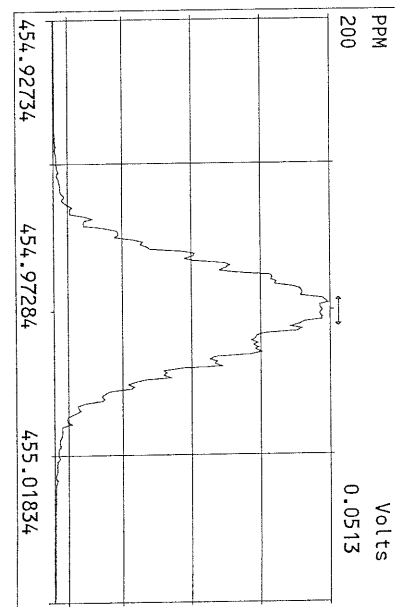
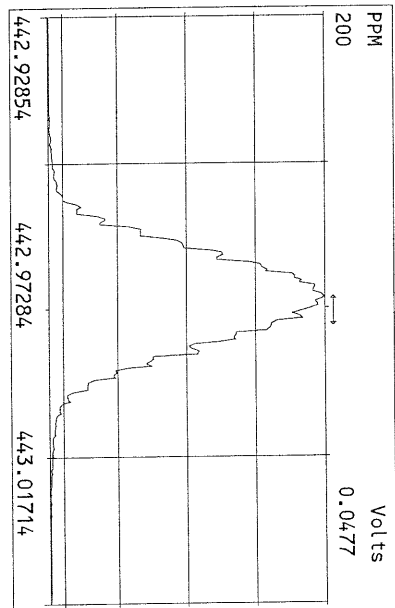
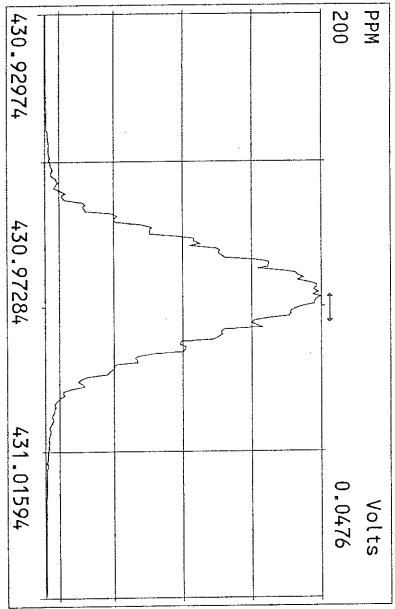


Peak Locate Examination:20-APR-2011:10:15 File:20APR11M
Experiment:OCDD Function:2 Reference:PFK

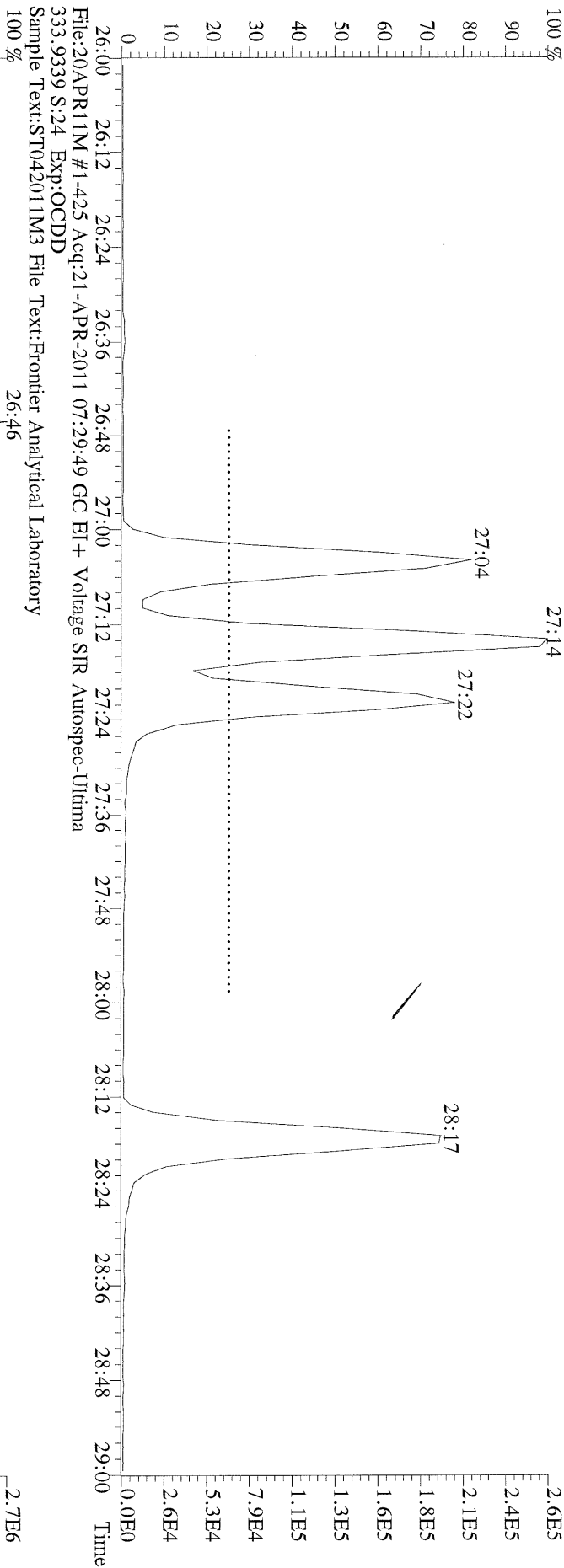




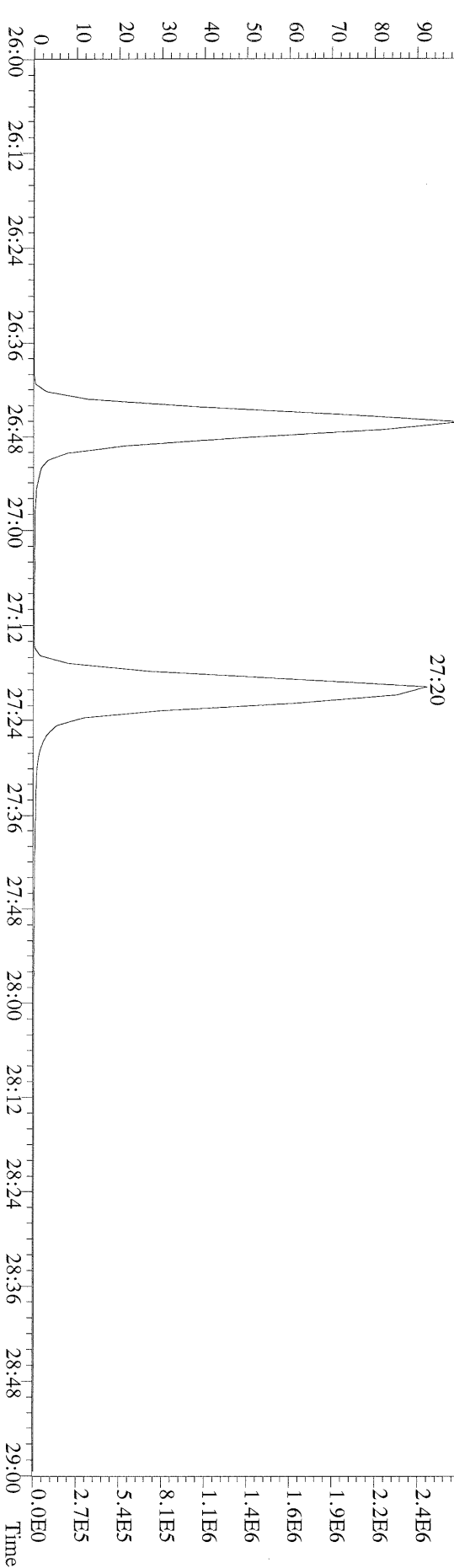




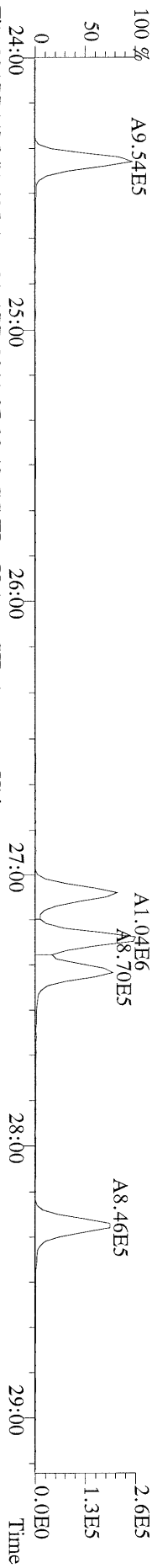
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
319.8965 S:24 Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



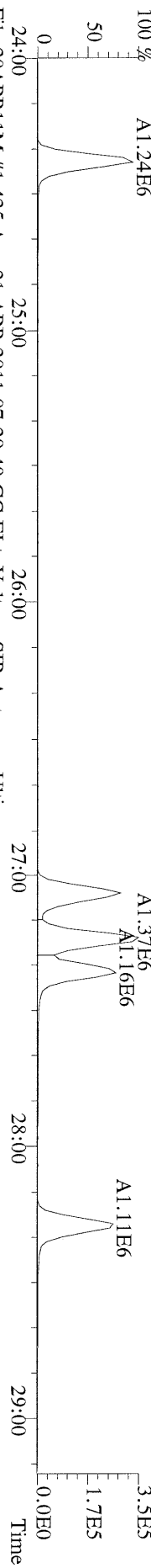
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333.9339 S:24 Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



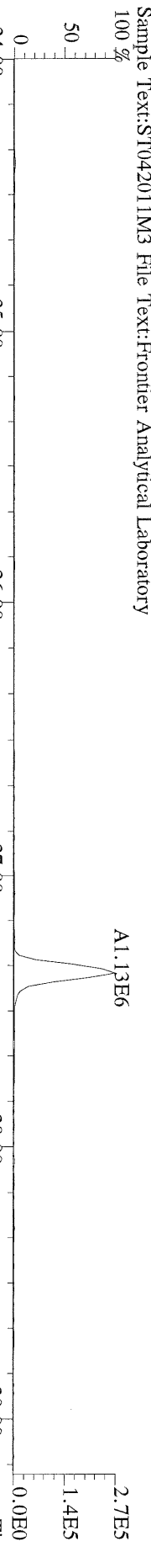
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319,8965 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



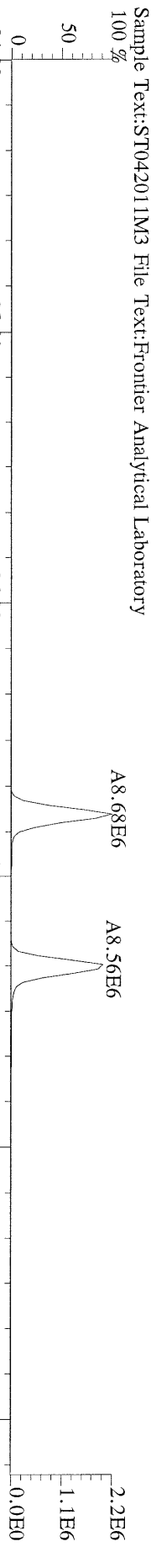
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321,8936 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



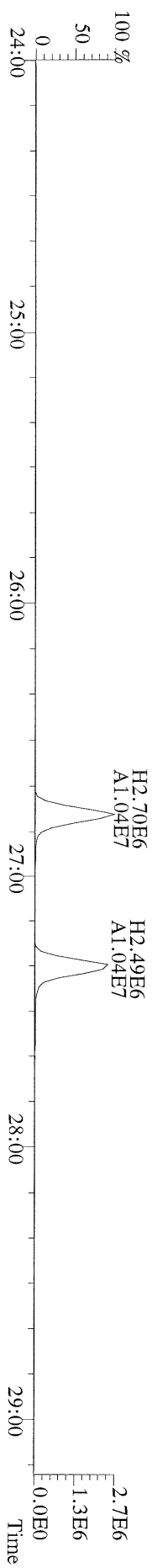
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327,8847 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



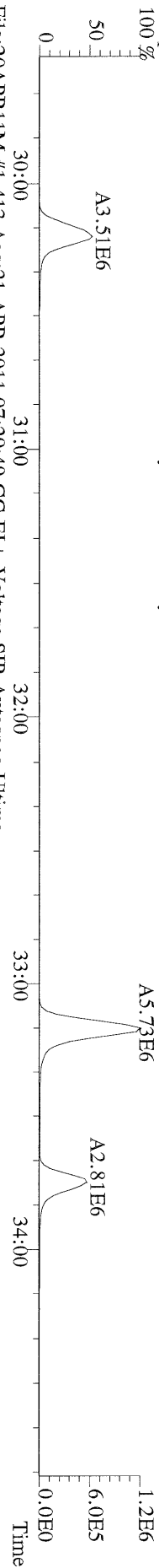
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331,9368 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



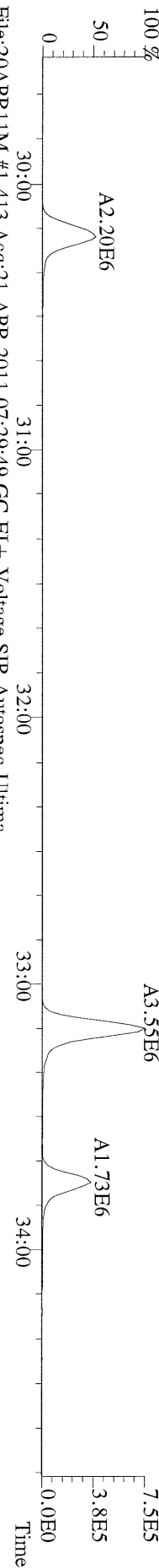
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333,9339 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



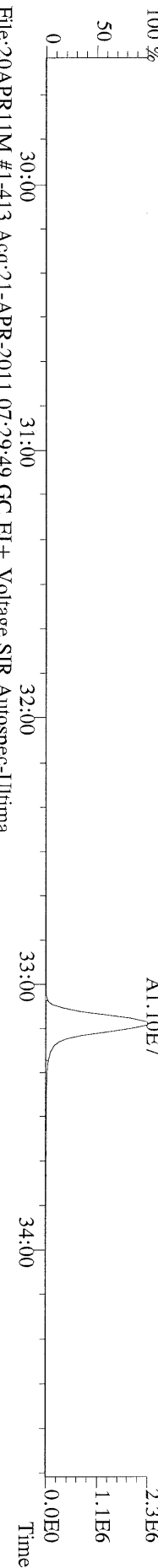
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355.8546 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 %



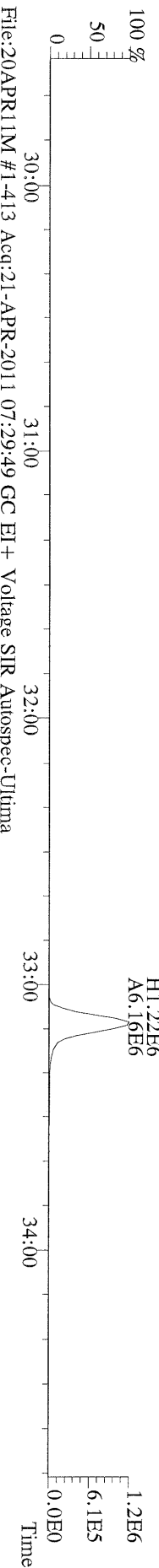
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357.8517 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 %



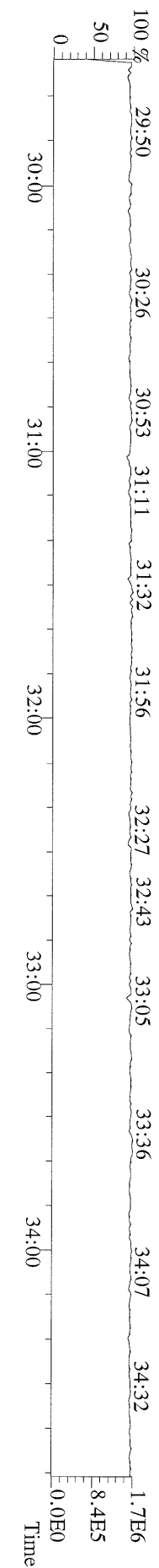
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367.8949 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 %



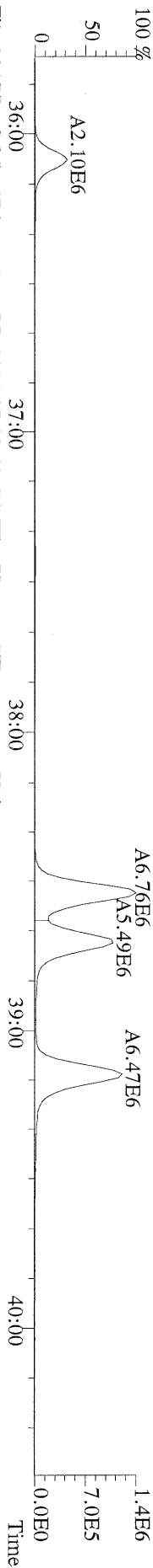
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369.8919 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



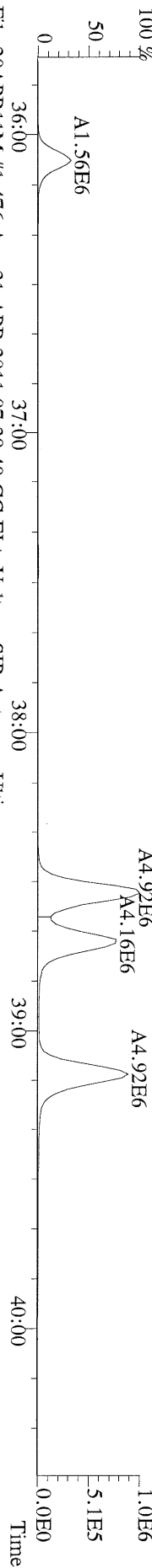
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366.9792 S:24 F:2 Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



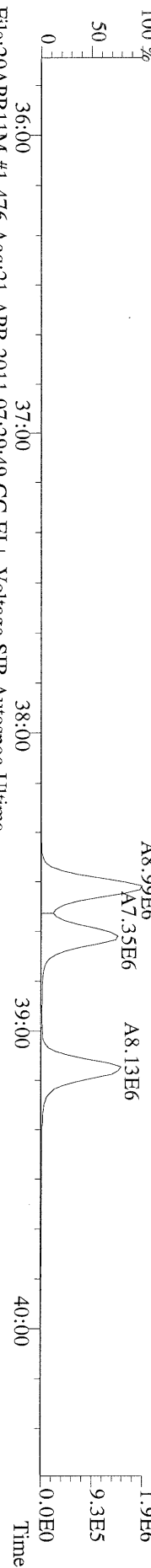
File:20APR11M #1-476 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
389.8156 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



File:20APR11M #1-476 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
391.8127 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



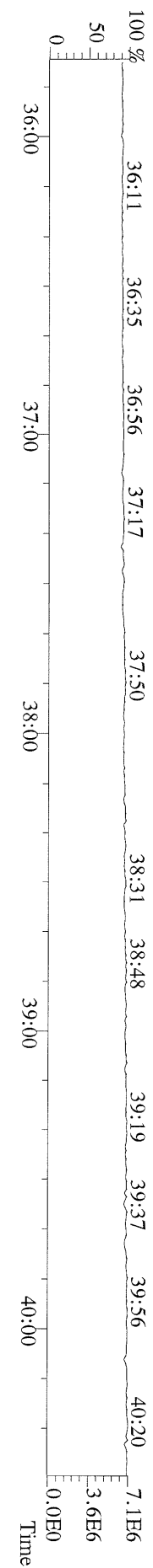
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401.8559 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



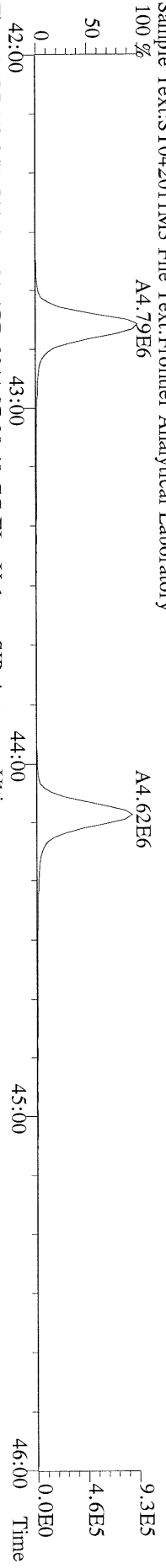
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403.8530 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



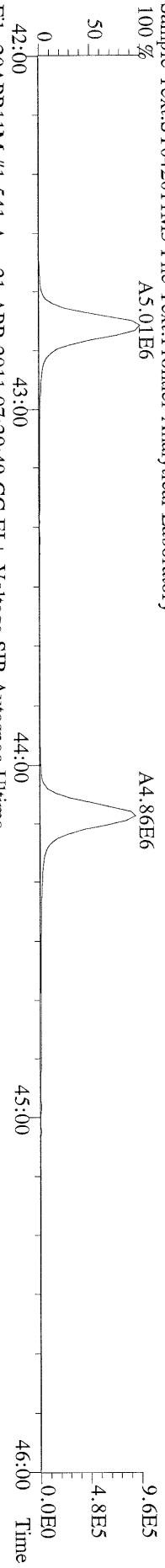
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380.9760 S:24 F:3 Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



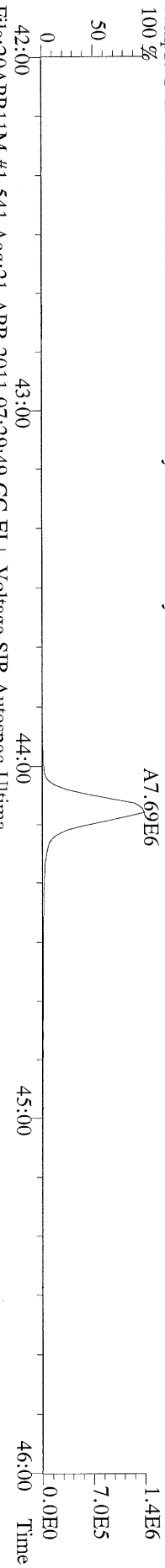
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423.7767 S:24 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



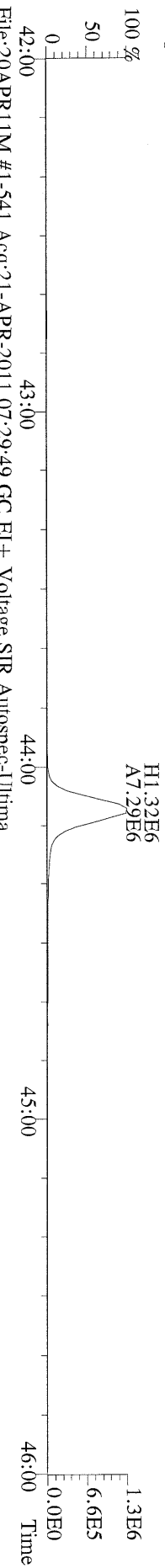
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425.7737 S:24 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



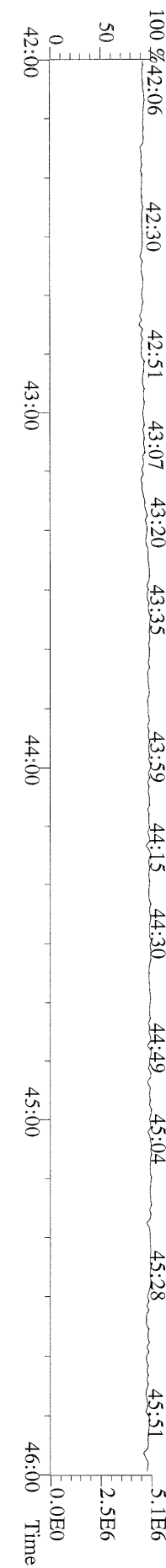
File:20APR11M #1-541 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
435.8169 S:24 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



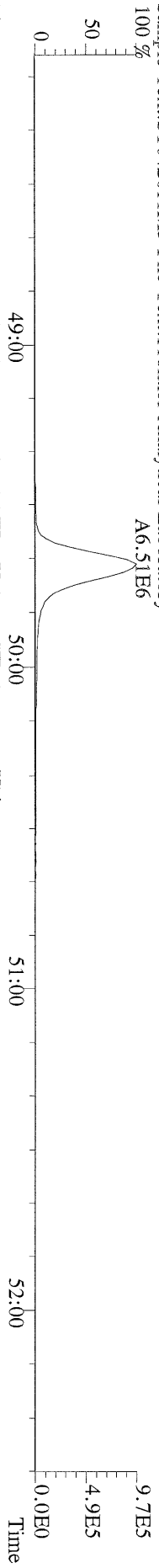
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437.8140 S:24 F:4 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



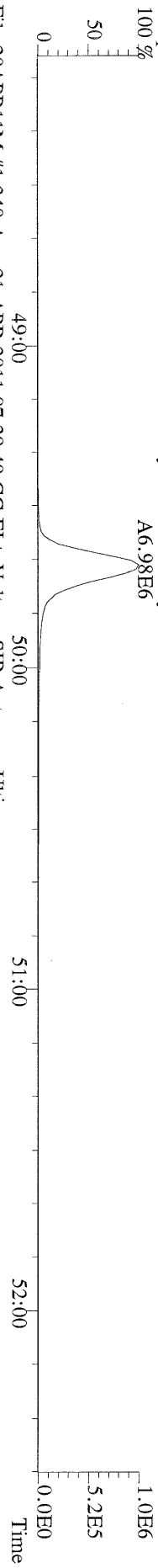
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430.9728 S:24 F:4 Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



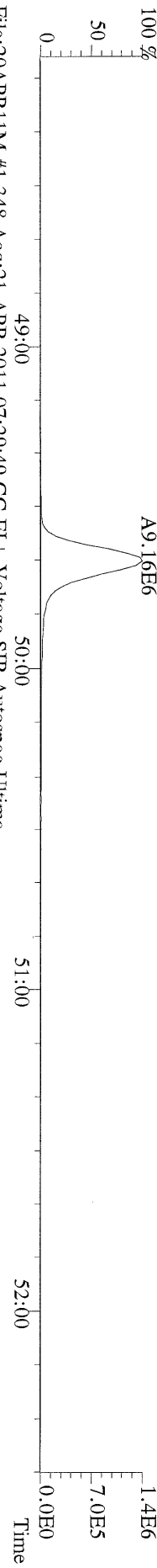
File:20APR11M #1-348 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
457.7377 S:24 F:5 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100%



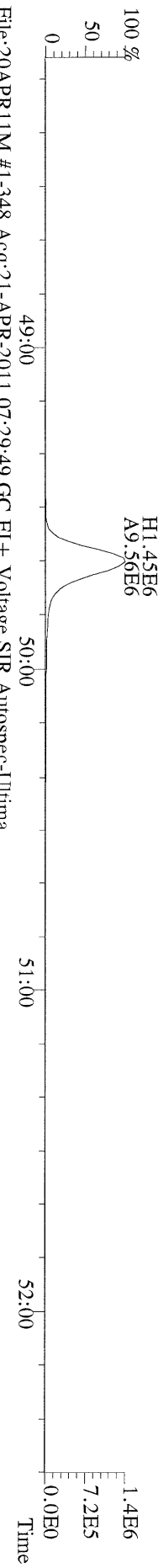
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459.7348 S:24 F:5 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100%



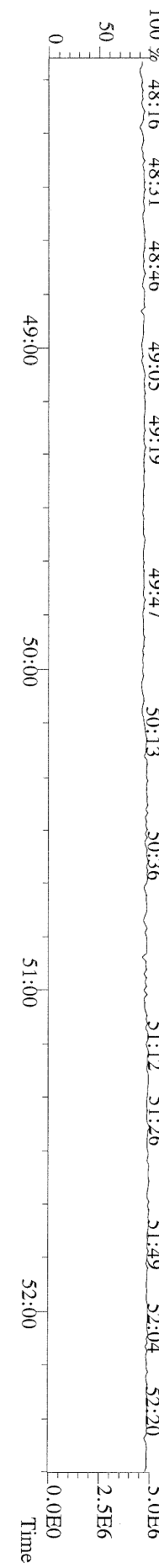
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469.7780 S:24 F:5 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
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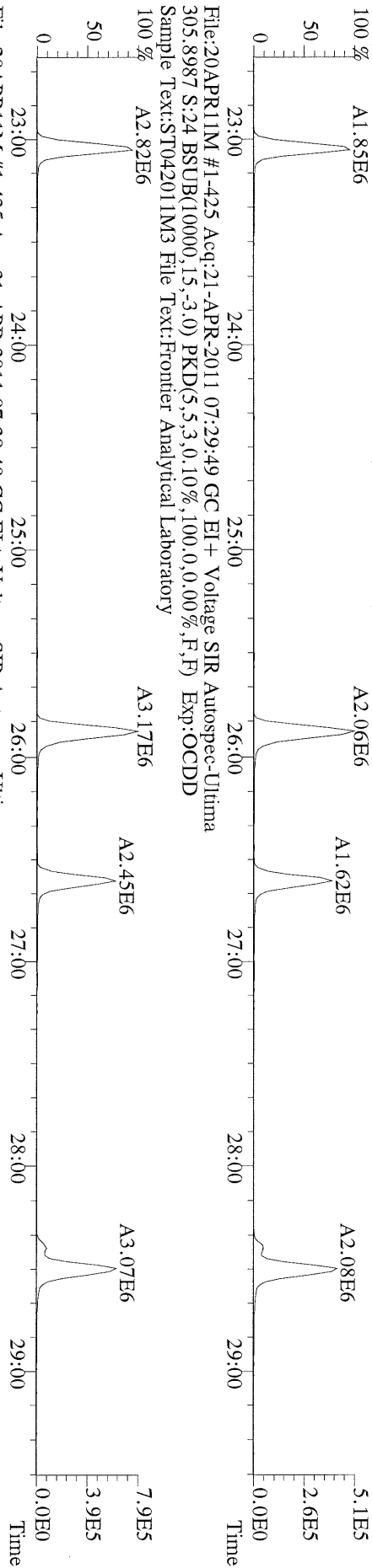
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471.7750 S:24 F:5 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100%



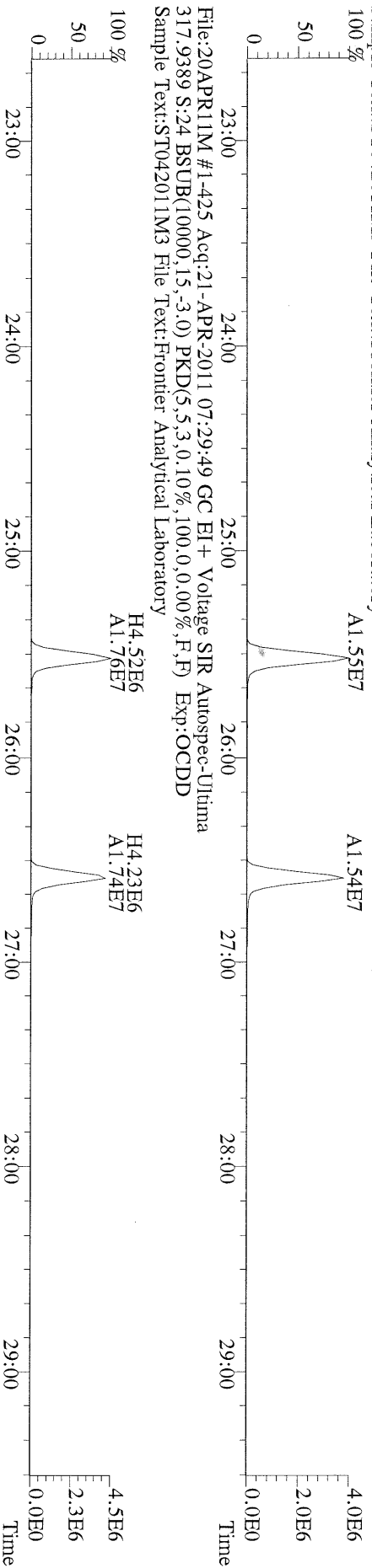
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454.9728 S:24 F:5 Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100%



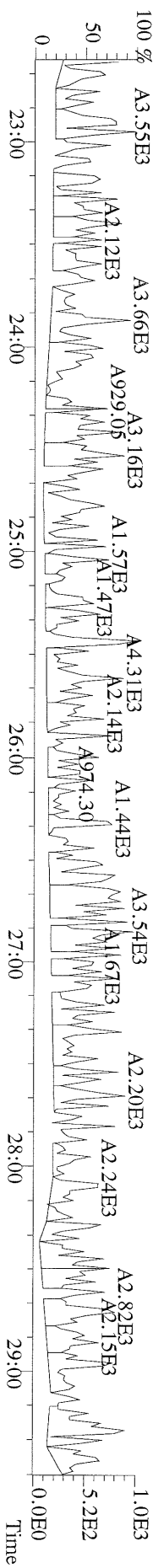
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
303.9016 S:24 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



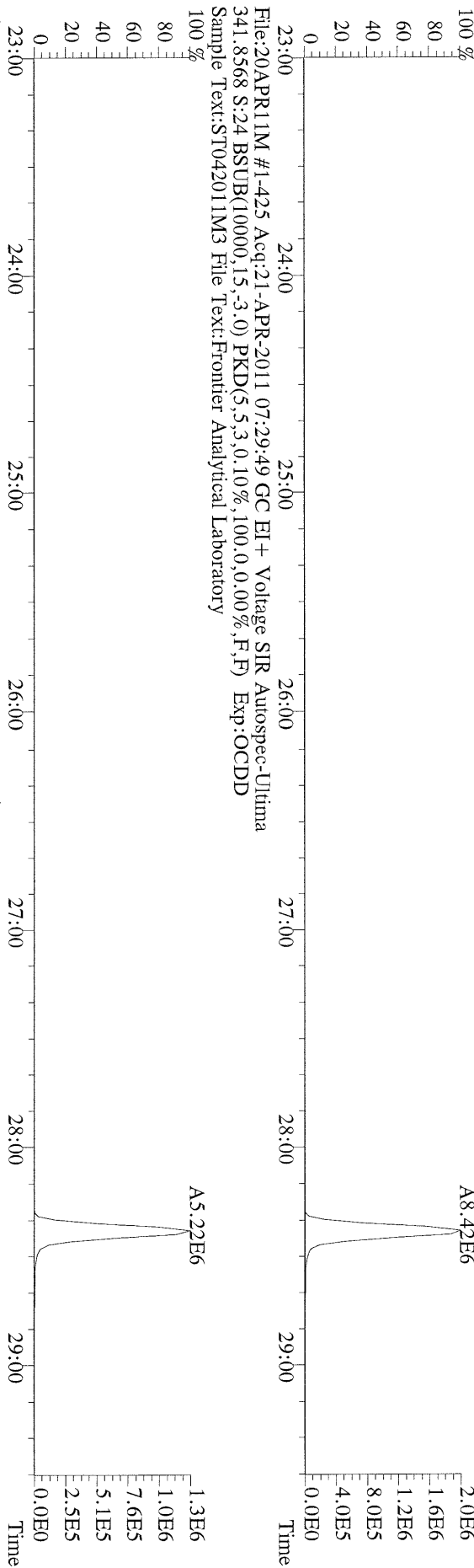
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
315.9419 S:24 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



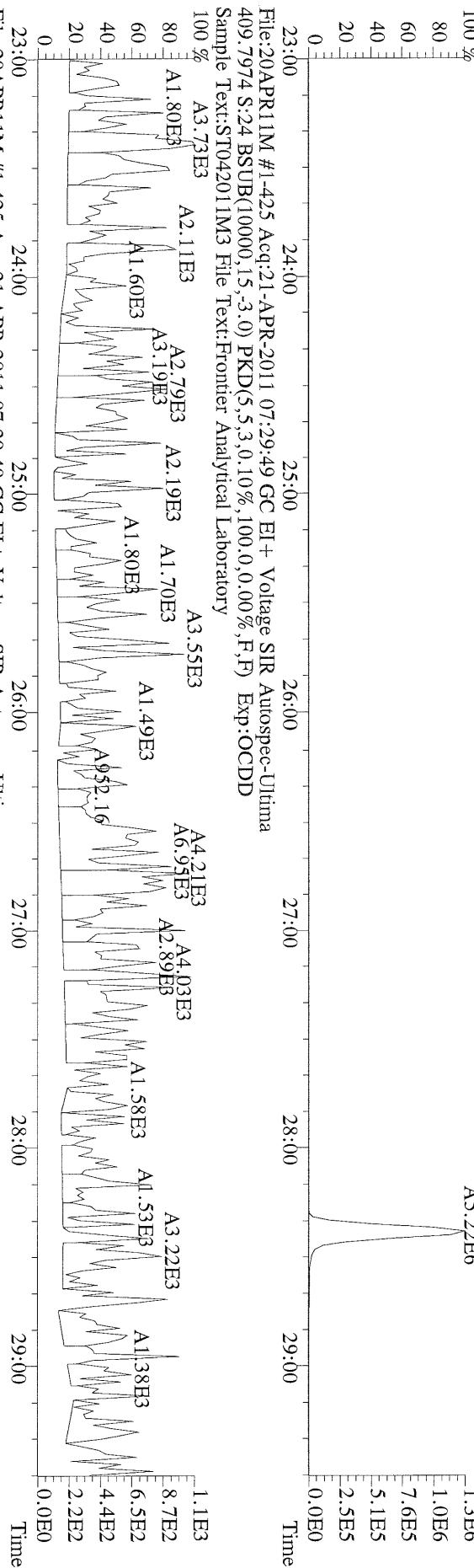
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
375.8364 S:24 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0,0.00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



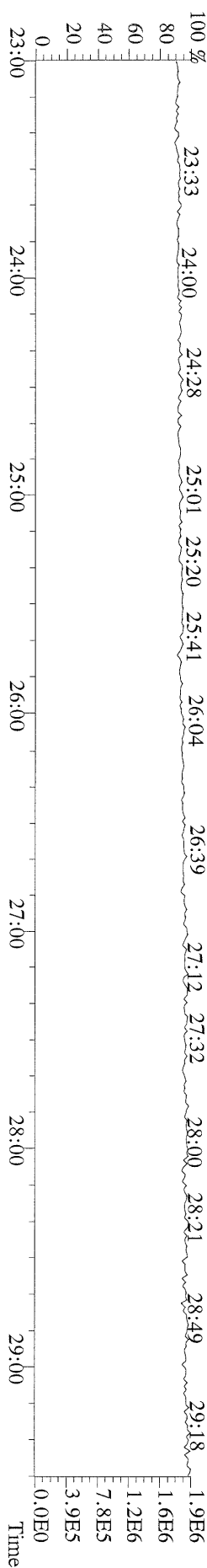
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Utima
 339.8597 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



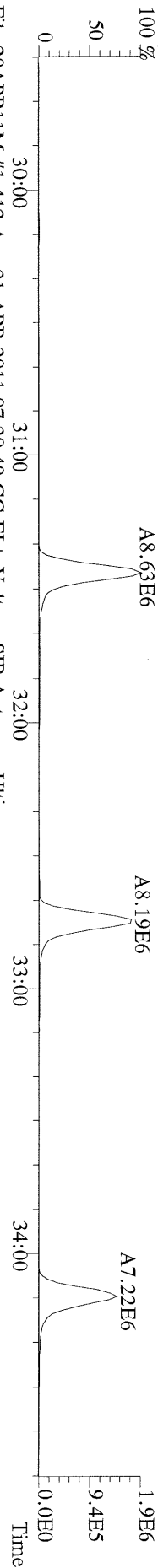
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 341.8568 S:24 BSUB(10000,15,-3.0) PKD(5,5,3.0,10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



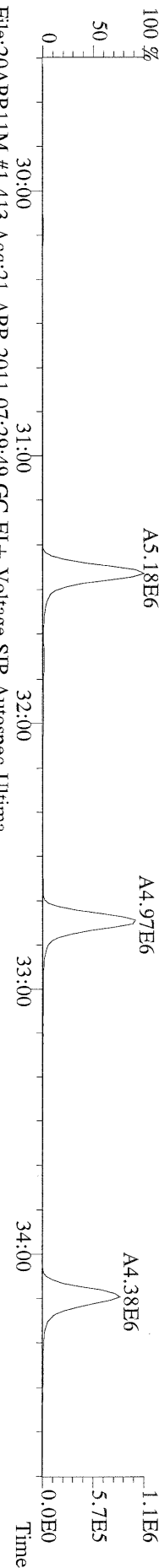
File:20APR11M #1-425 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Utima
 316.9824 S:24 Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



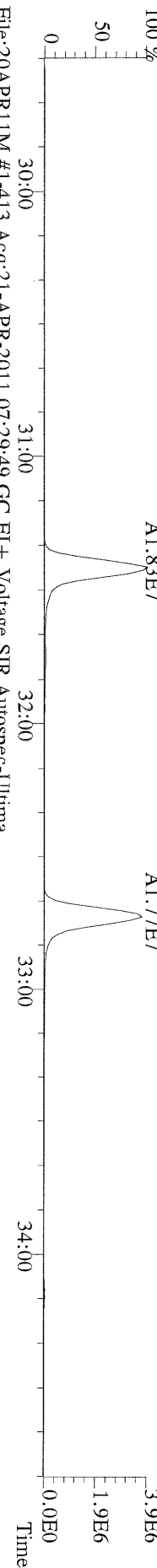
File:20APR11M #1-413 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
339.8597 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



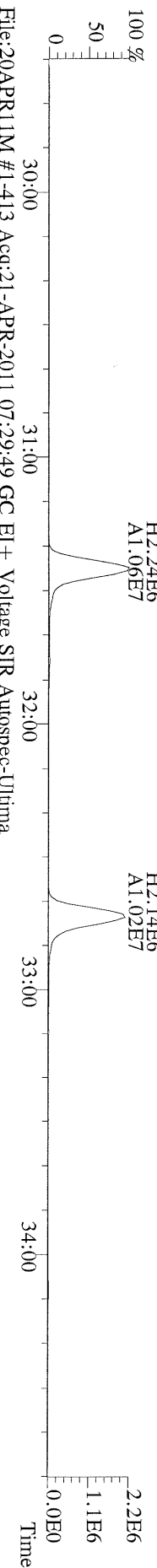
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341.8568 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



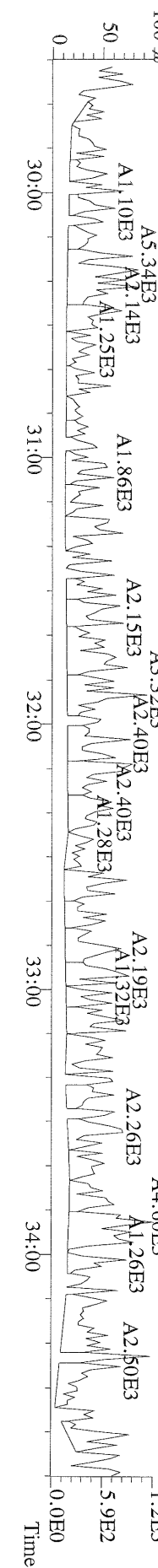
File:20APR11M #1-413 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
351.9000 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



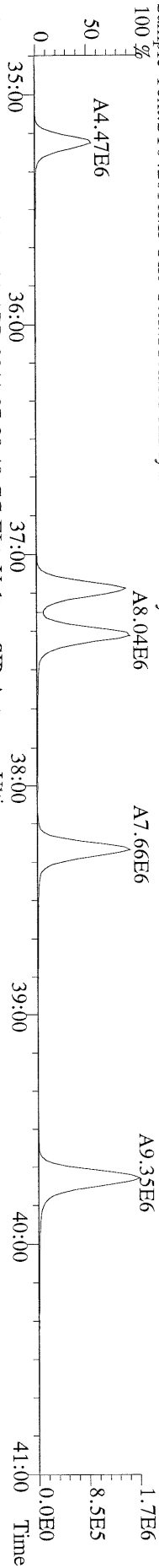
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353.8970 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



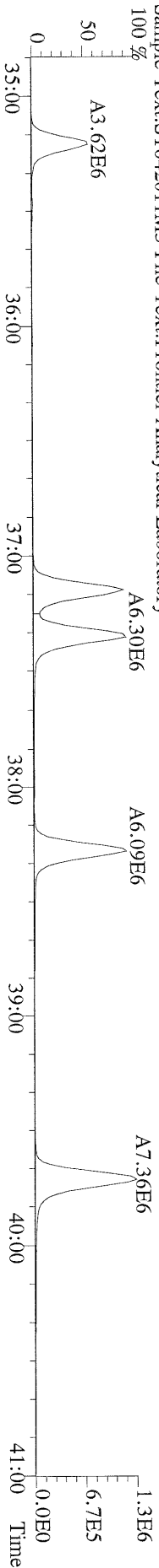
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409.7974 S:24 F:2 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory



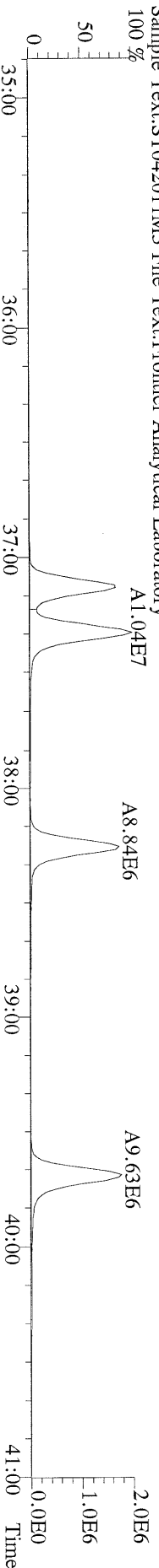
File:20APR11M #1-476 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Utima
 373.8207 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



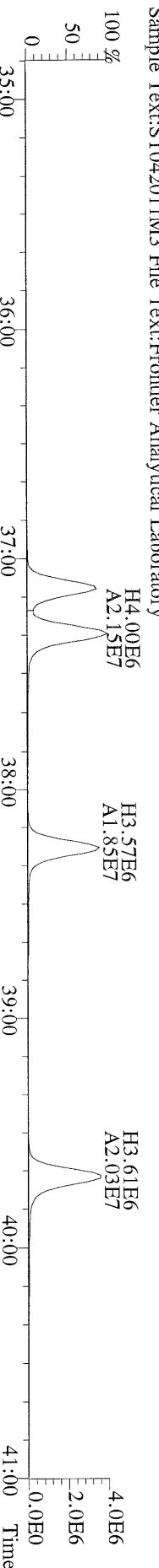
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 375.8178 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



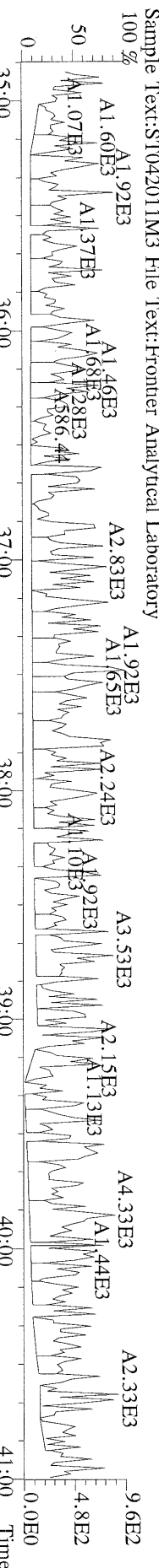
File:20APR11M #1-476 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Utima
 383.8639 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



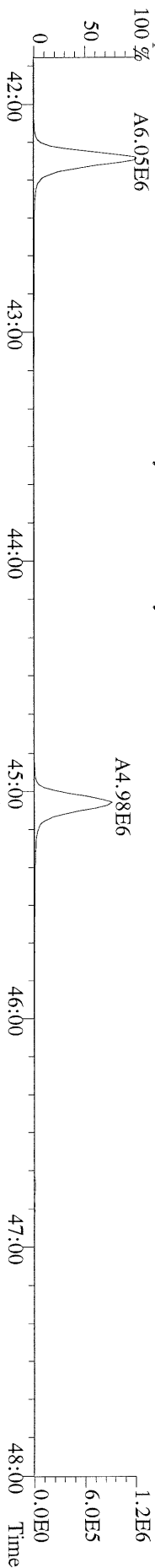
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 385.8610 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



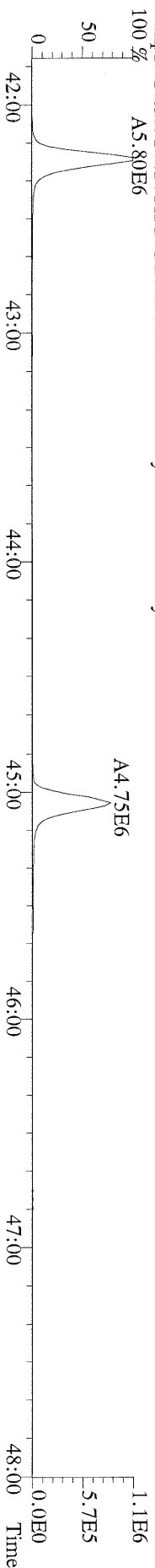
File:20APR11M #1-476 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Utima
 445.7555 S:24 F:3 BSUB(10000,15,-3.0) PKD(5.5,3,0.10%,100.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



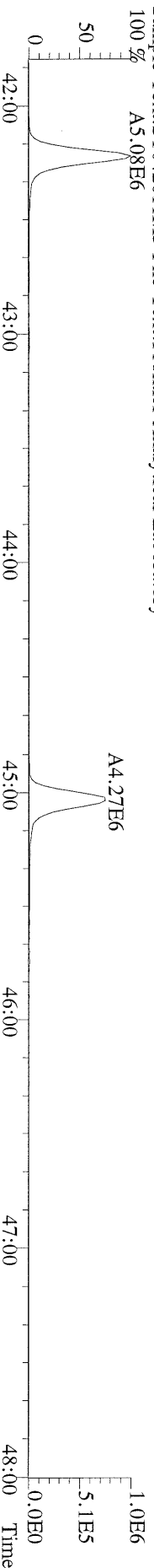
File:20APR11M #1-541 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
407.7818 S:24 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 % A6.05E6



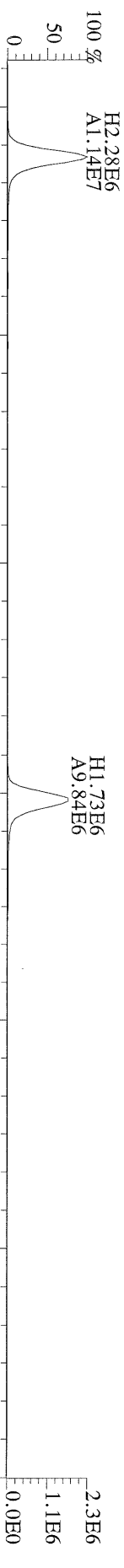
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409.7788 S:24 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 % A5.80E6



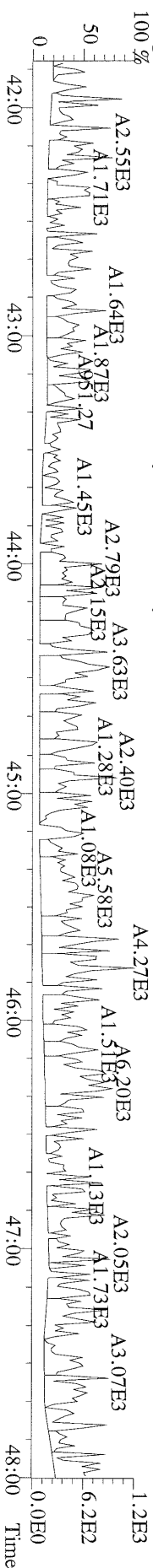
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417.8253 S:24 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory
100 % A5.08E6



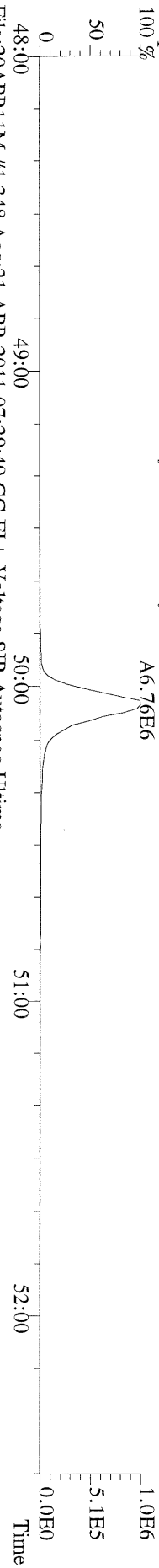
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419.8220 S:24 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



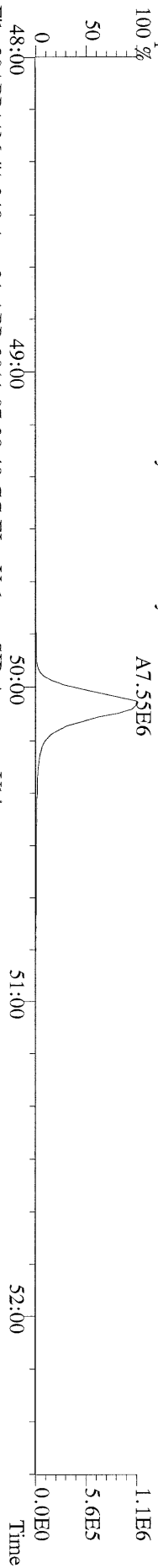
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479.7165 S:24 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST042011M3 File Text:Frontier Analytical Laboratory



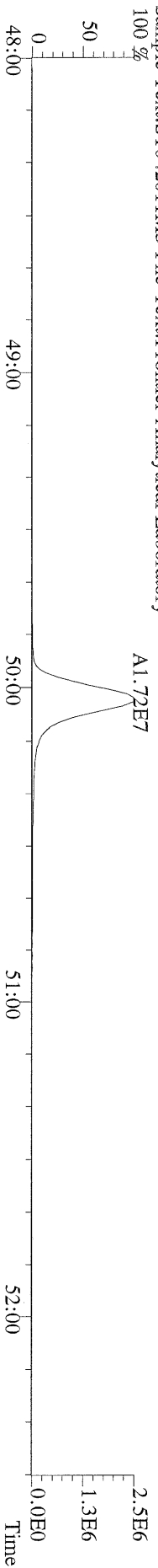
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441.7428 S:24 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100 %



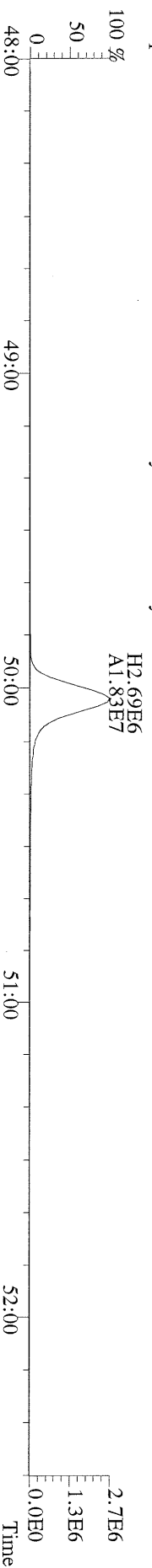
File:20APR11M #1-348 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
443.7398 S:24 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100 %



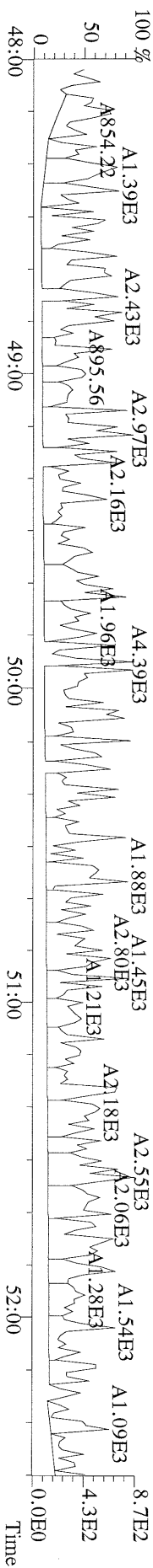
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453.7831 S:24 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100 %

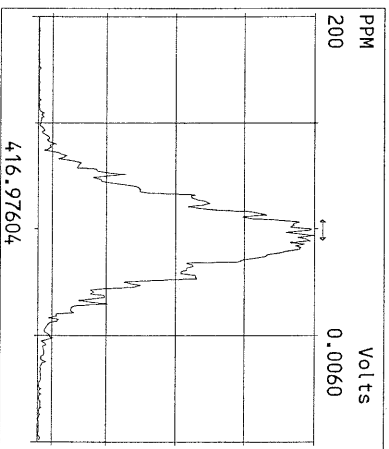
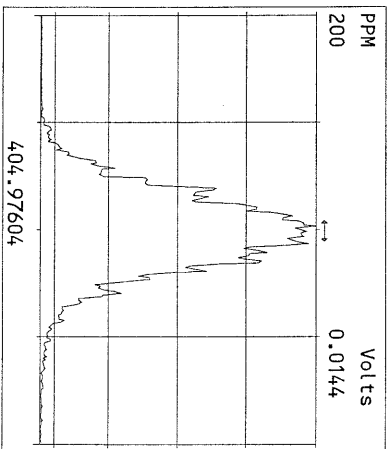
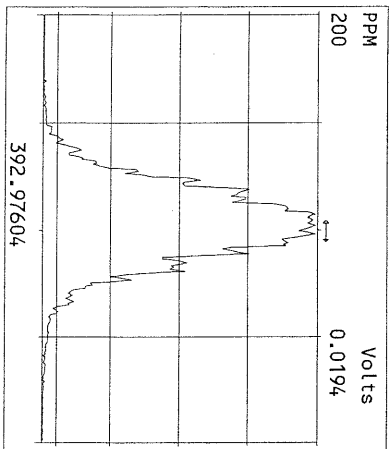
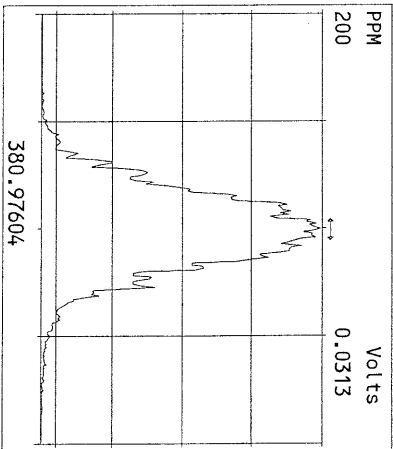
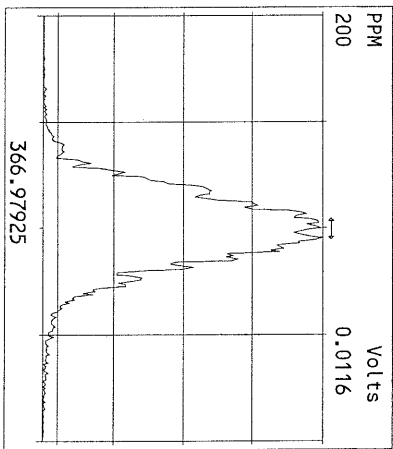
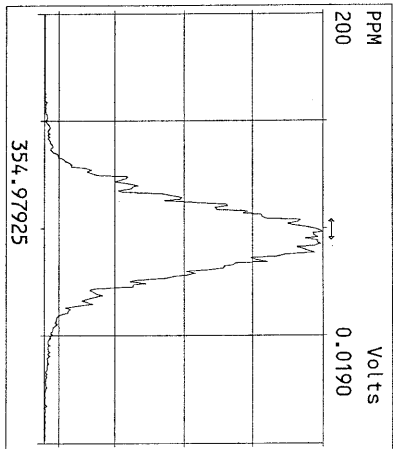
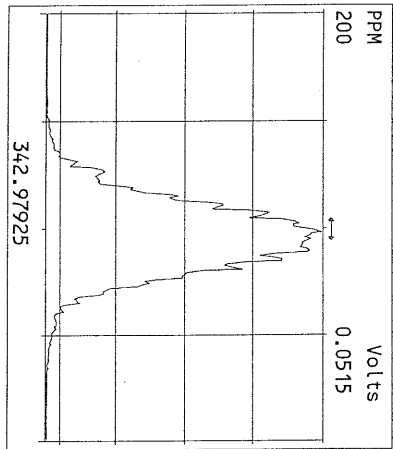
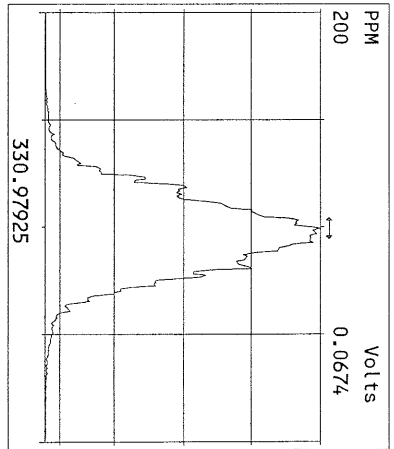
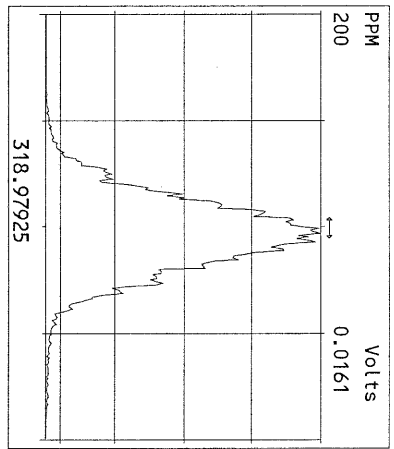
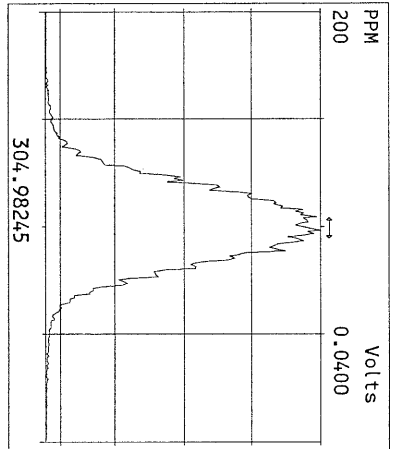
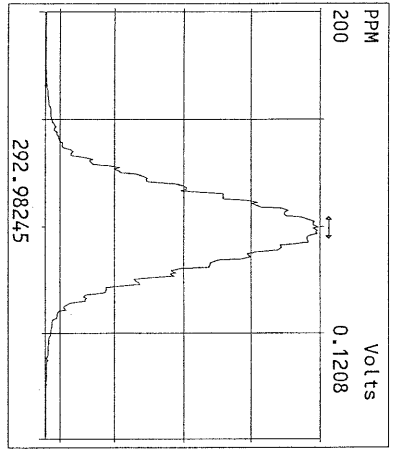


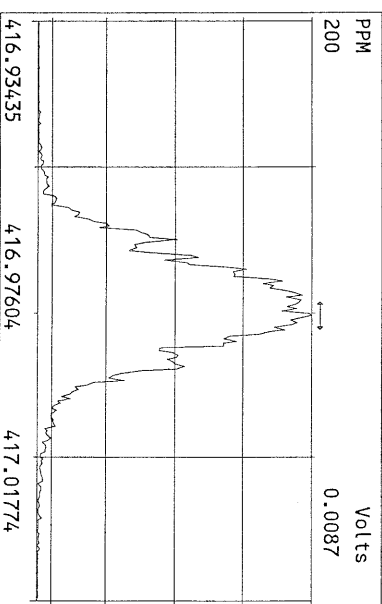
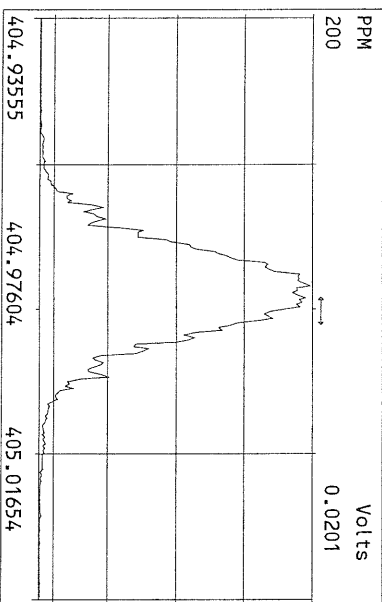
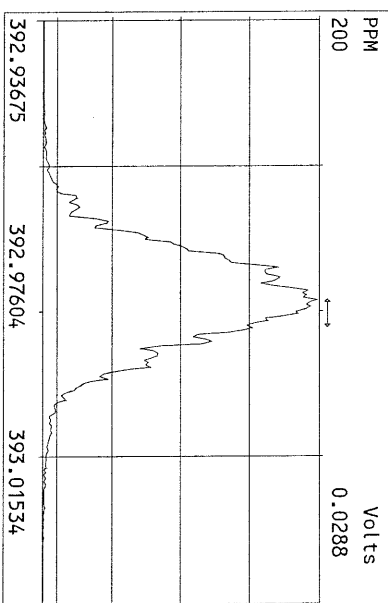
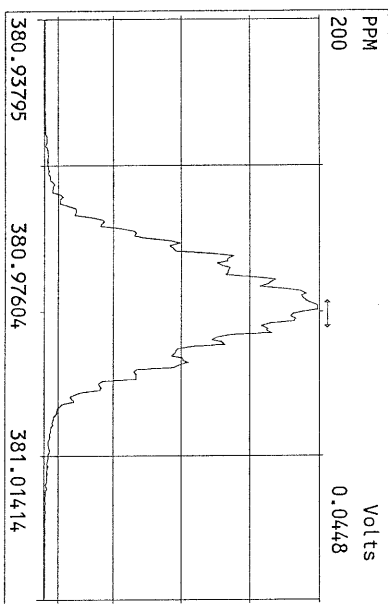
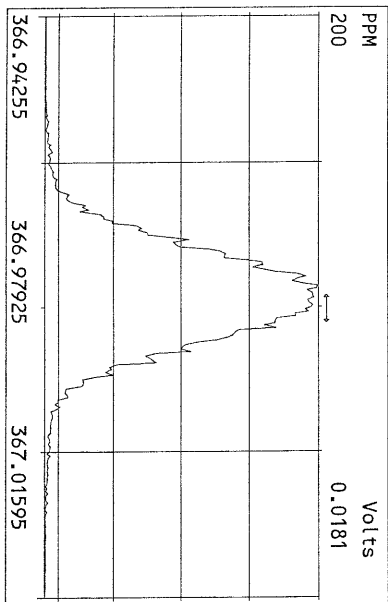
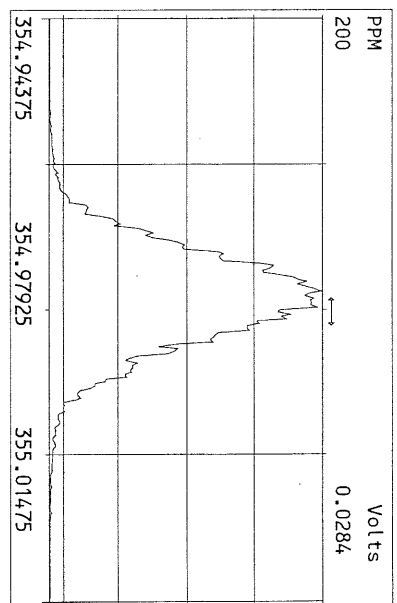
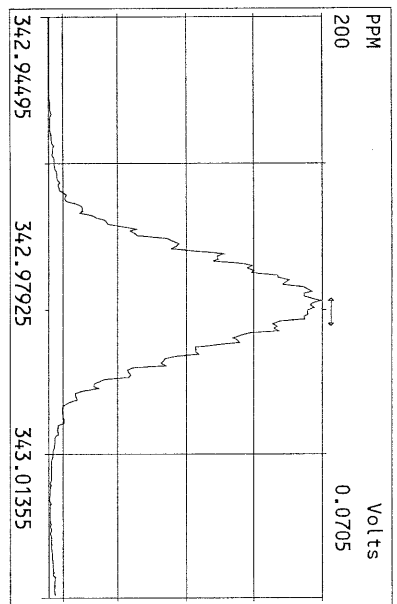
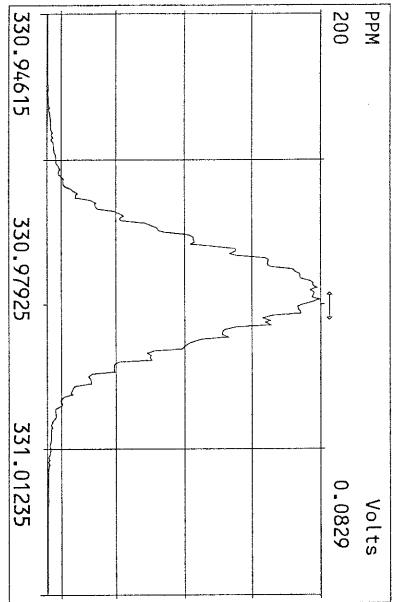
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455.7801 S:24 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory

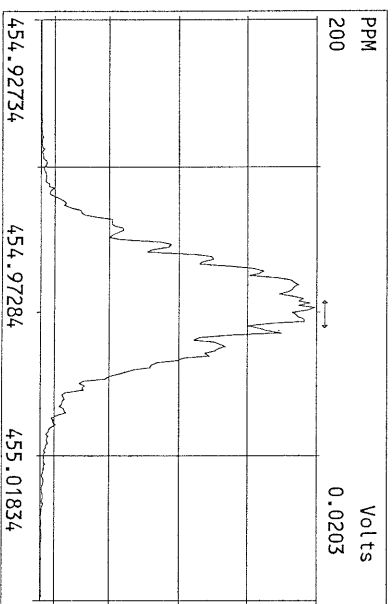
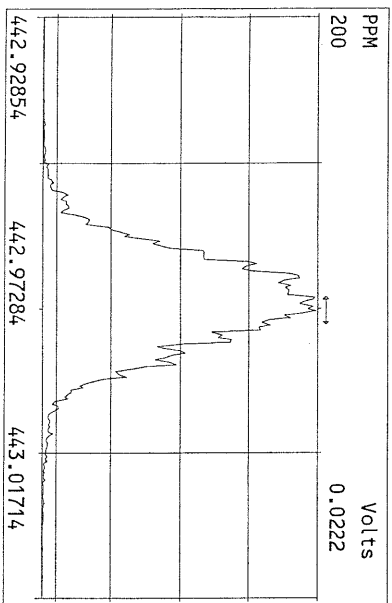
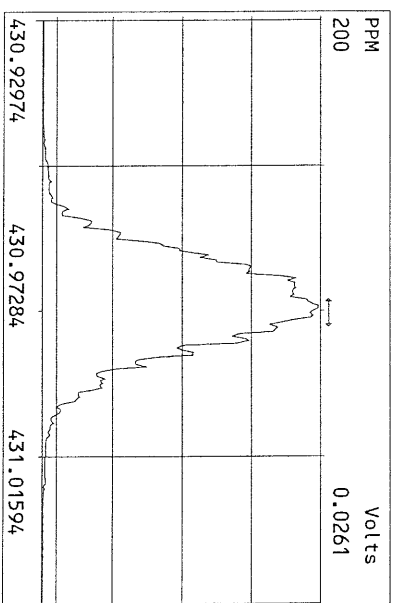
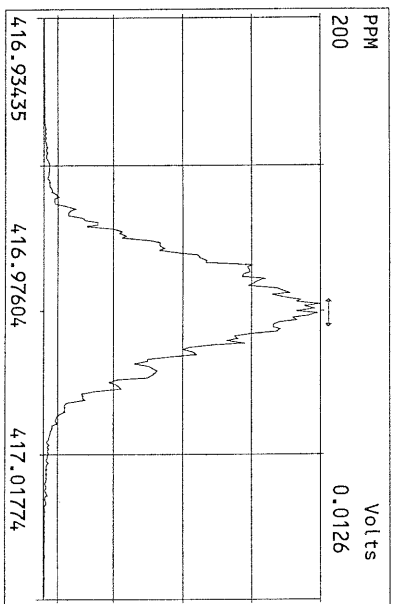
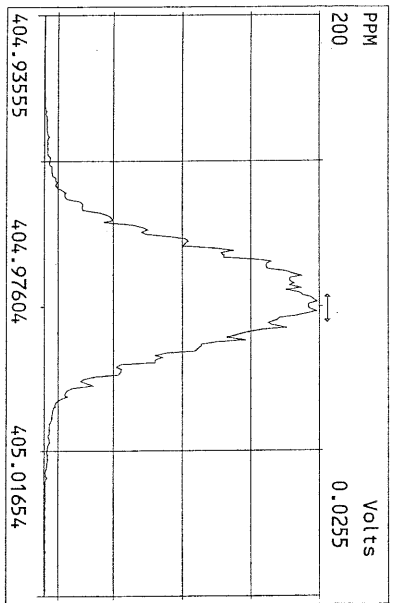
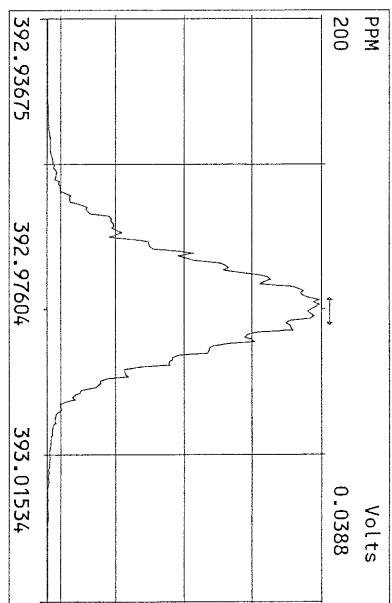
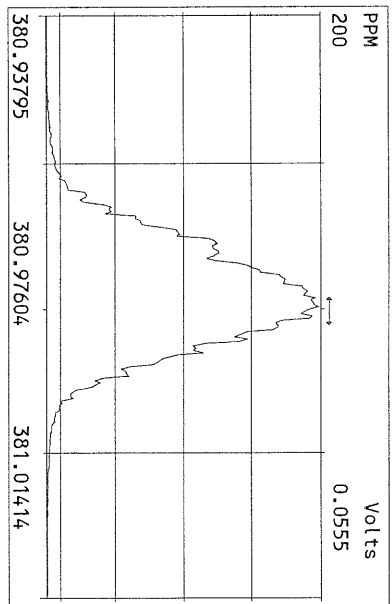
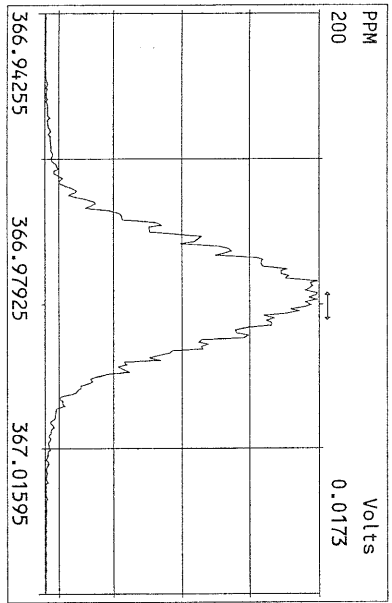


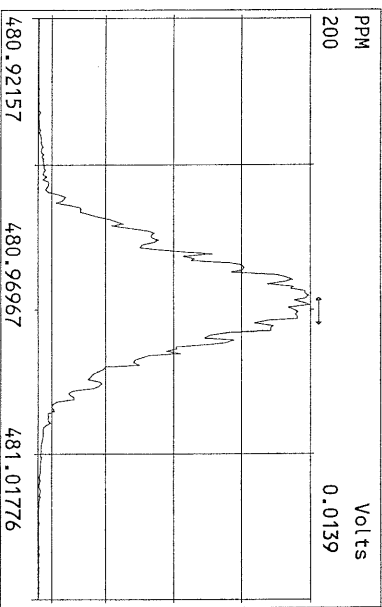
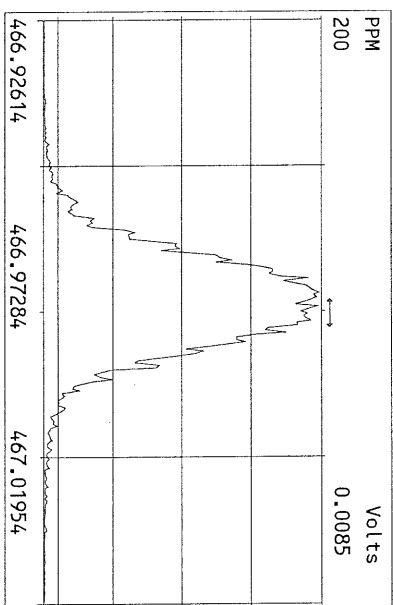
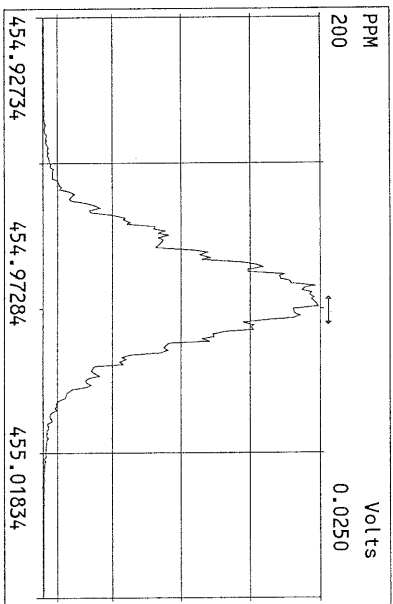
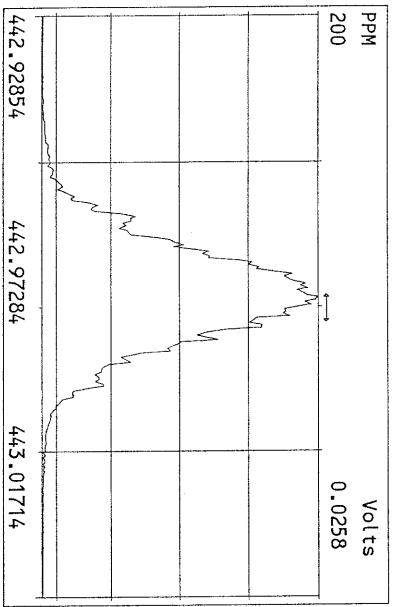
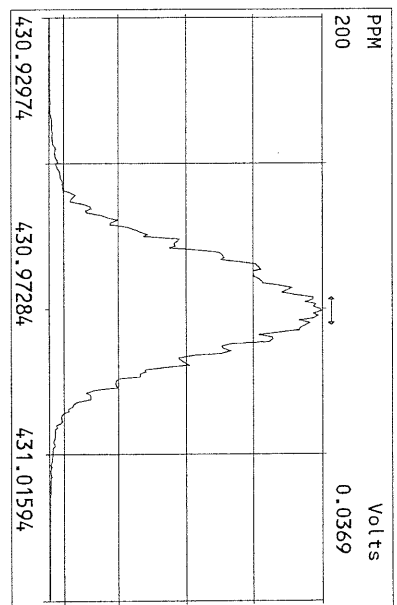
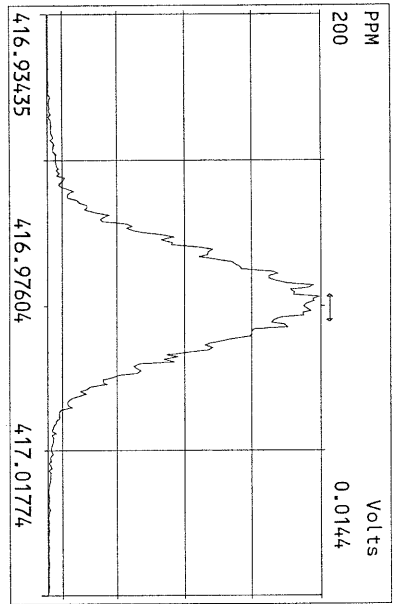
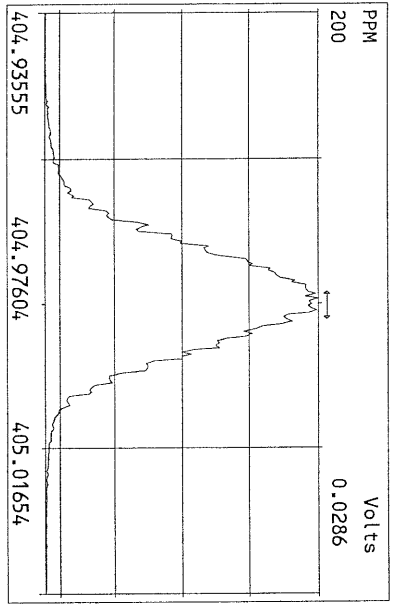
File:20APR11M #1-348 Acq:21-APR-2011 07:29:49 GC EI+ Voltage SIR Autospec-Ultima
513.6775 S:24 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04201IM3 File Text:Frontier Analytical Laboratory
100 %

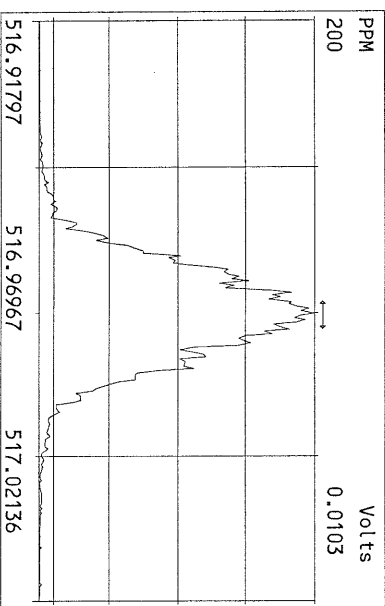
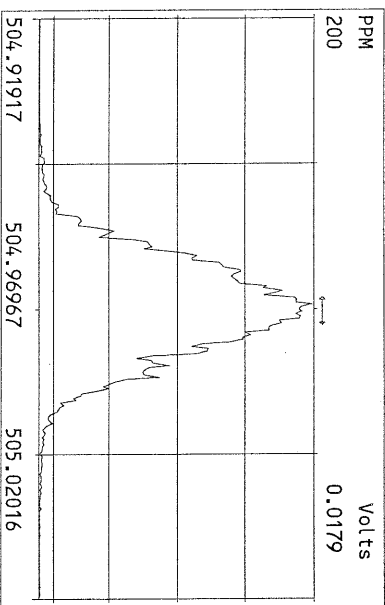
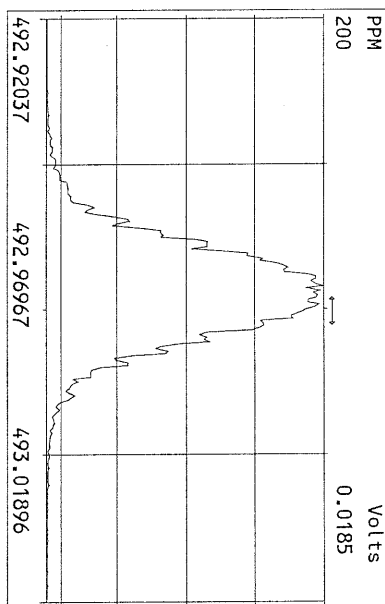
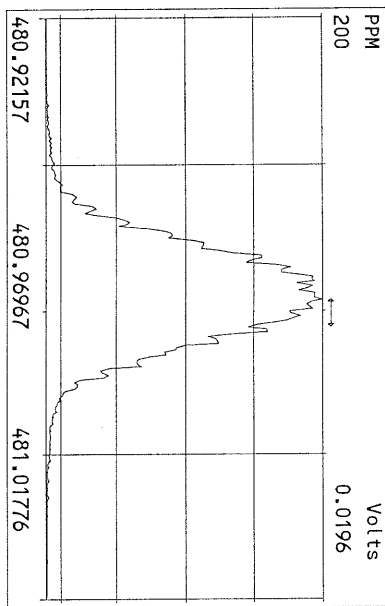
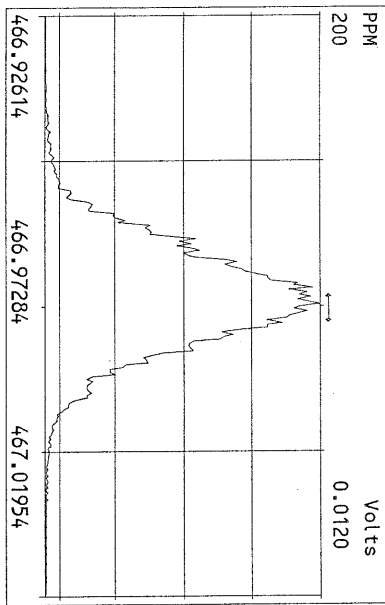
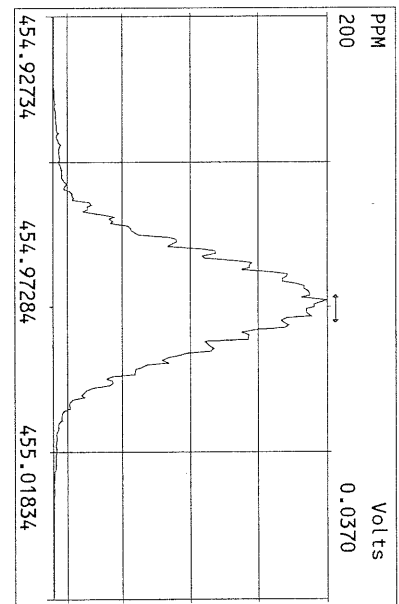
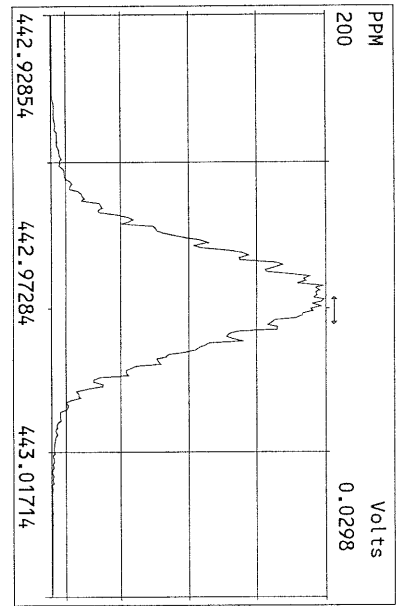
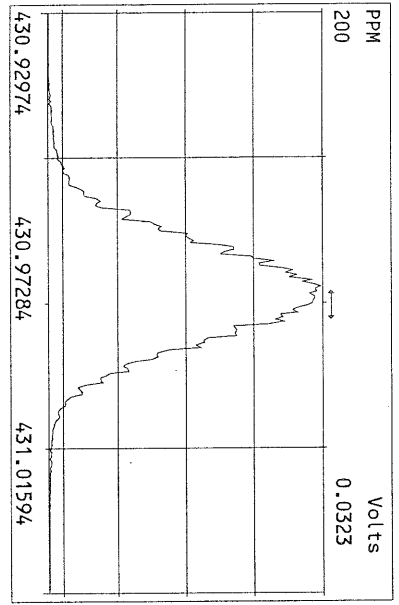












USEPA - ITD

FORM 4A
TCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date: 2/18/11

Instrument ID: FAL1 GC Column ID: DB225

VER Data Filename: 20APR11A Sam:1 Analysis Date: 20-APR-11 Time: 09:25:20

	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
NATIVE ANALYTES						
2,3,7,8-TCDF	M/M+2	0.75	0.65-0.89	y	8.59	8.40 - 12.0 ✓
LABELED COMPOUNDS						
13C-2,3,7,8-TCDF	M/M+2	0.79	0.65-0.89	y	73.1	71.0 - 140 ✓

(1) See Table 8, Method 1613, for m/z specifications.

(2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.



(3) Contract-required concentration range as specified in Table 6A, Method 1613

Analyst: 

Date: 4/20/11

FAL ID: ST042011A1 Filename: 20APR11A Sam:1 Acquired: 20-APR-11 09:25:20 ICal: TCDFFAL1-2-18-11
Client ID: 1613 CS3 100511J ConCal: ST042011A1 EndCal: ST042011A2
Results: 6701TCDF GC Column: DB225 Amount: 1.000


Name	Resp	RA	RT	RRF	Conc	Qual	Fac	Noise	DL	#Hom	Rec
2,3,7,8-TCDF	1.18e+07	0.75 y	19:19	1.16	8.59		2.50	-	-	1	
13C-2,3,7,8-TCDF	1.19e+08	0.79 y	19:18	1.05	73.1						73.1
13C-1,2,3,4-TCDF	1.55e+08	0.78 y	16:45	-	87.5						

Analyst:  Date: 

Frontier Analytical Laboratory - Acquisition Log

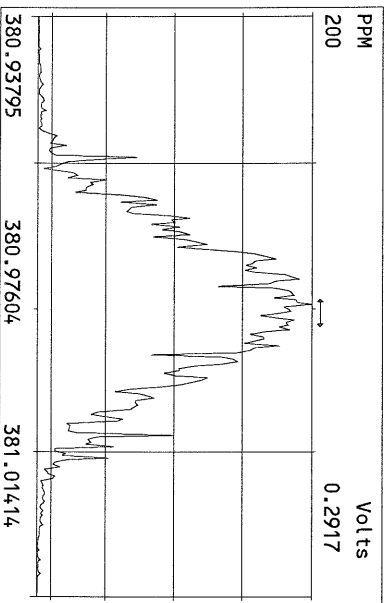
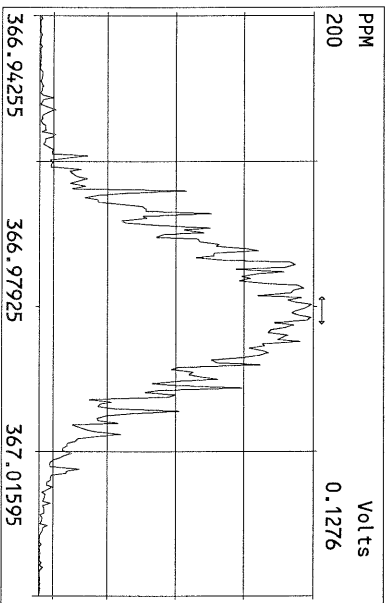
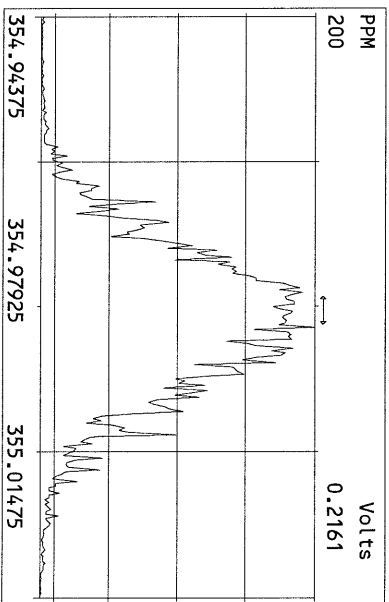
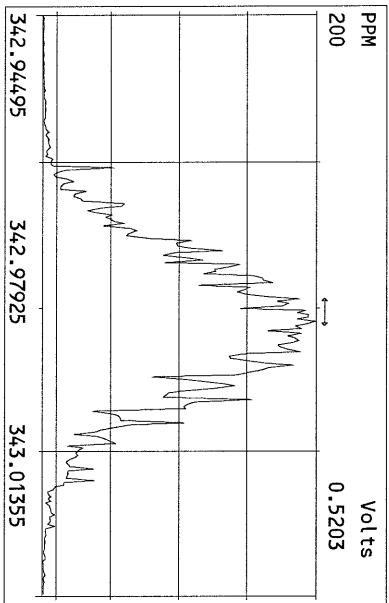
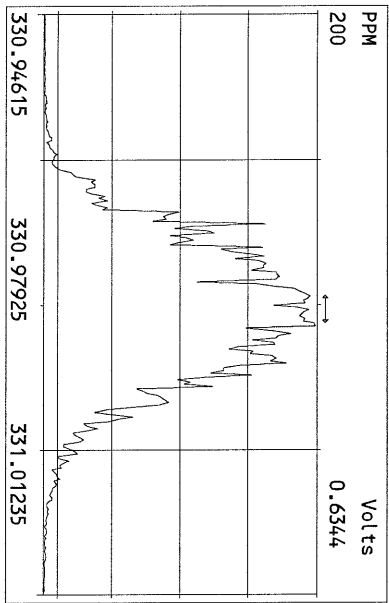
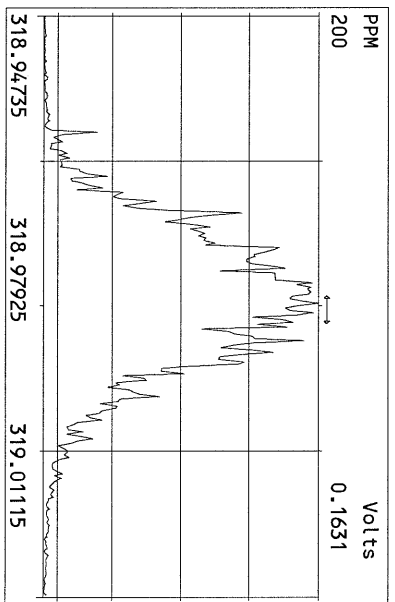
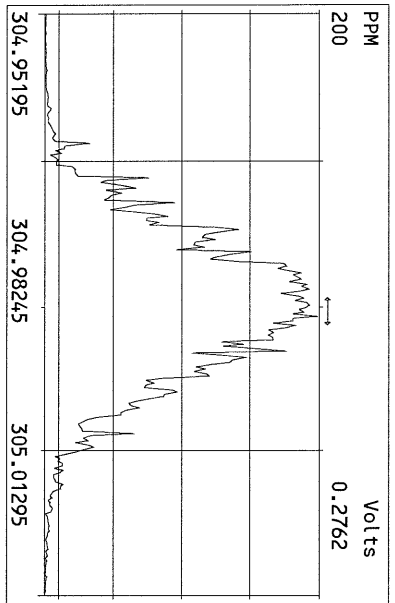
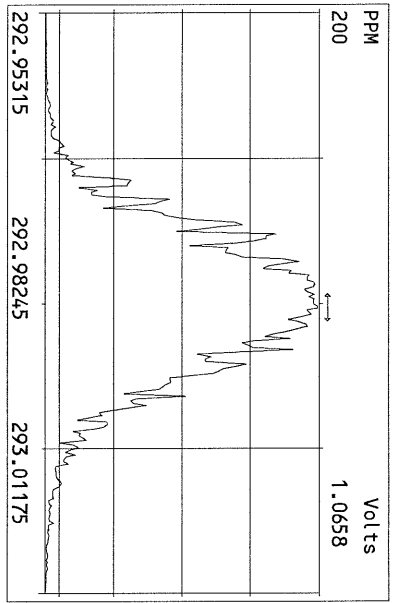
Run Name:20APR11A Instrument: FAL1 GC: DB225 Experiment:TCDF

Data File	S	FAL ID	Client ID	Acquired	ConCal	EndCal	Analyst
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20APR11A	2	6701-001-0001-SA	LL-SED1-0-15-032911	20-APR-11 10:00:23	ST042011A1	ST042011A2	TC
20APR11A	3	6701-002-0001-SA	LL-SED2-0-15-032911	20-APR-11 10:35:28	ST042011A1	ST042011A2	TC
20APR11A	4	6701-003-0001-SA	LL-SED3-0-15-032911	20-APR-11 11:10:31	ST042011A1	ST042011A2	TC
20APR11A	5	6701-004-0001-SA	LL-SED4-0-15-032911	20-APR-11 11:45:36	ST042011A1	ST042011A2	TC
20APR11A	6	6701-005-0001-SA	LL-SED1-0-15-032911-D	20-APR-11 12:20:39	ST042011A1	ST042011A2	TC
20APR11A	7	SB042011A1	Solvent Blank	20-APR-11 12:55:44	ST042011A1	ST042011A2	TC
20APR11A	8	SB042011A2	Solvent Blank	20-APR-11 13:30:47	ST042011A1	ST042011A2	TC
20APR11A	9	ST042011A2	1613 CS3 100511J	20-APR-11 14:06:38	ST042011A1	ST042011A2	TC

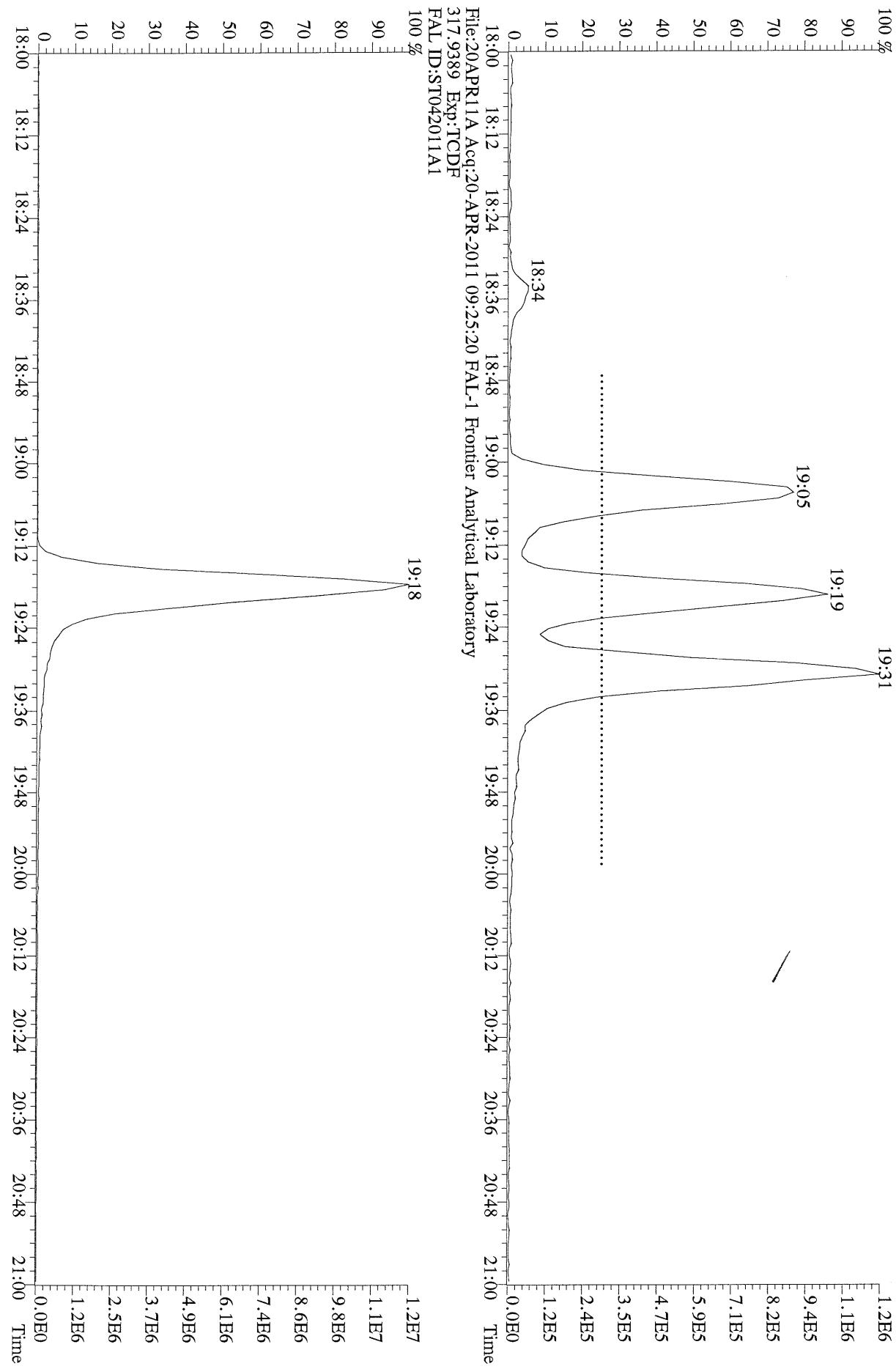
 4/20/11

Data Backed Up: _____

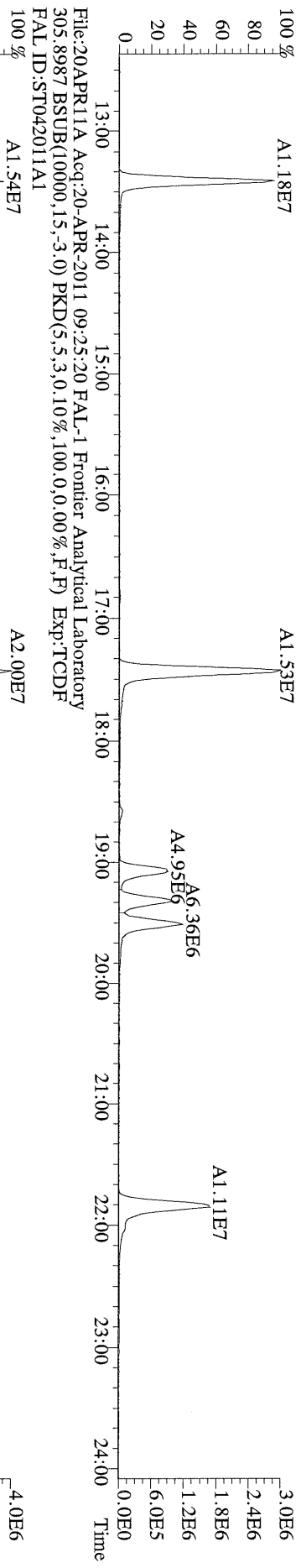
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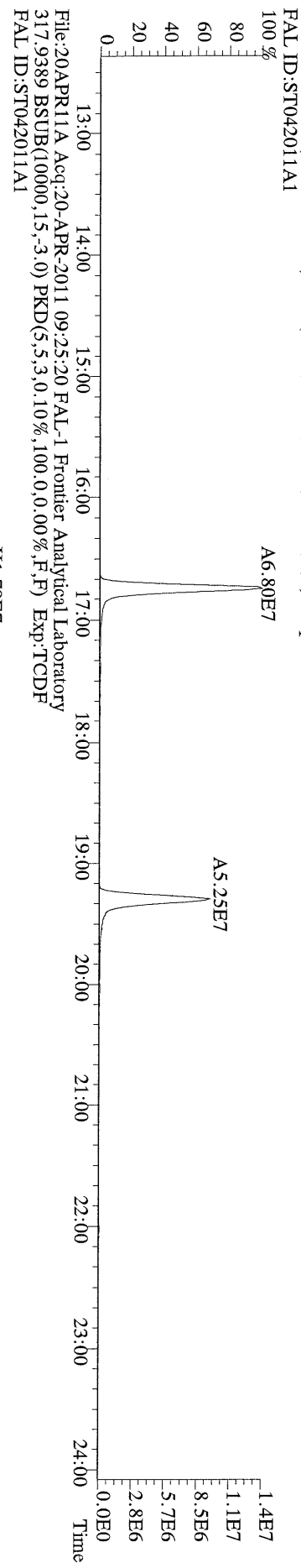
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303.9016 Exp:TCDF
FAL ID:ST042011A1
100%



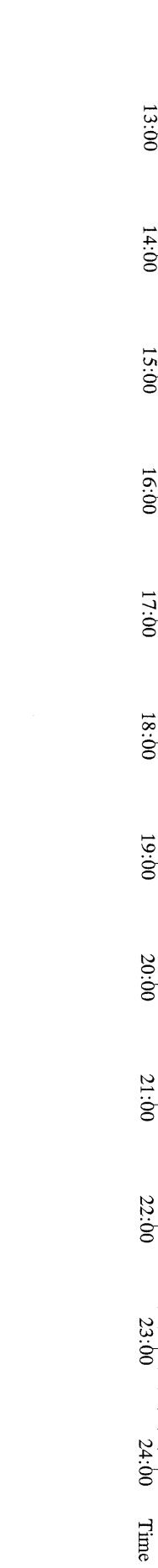
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FAL ID:ST042011A1
100%

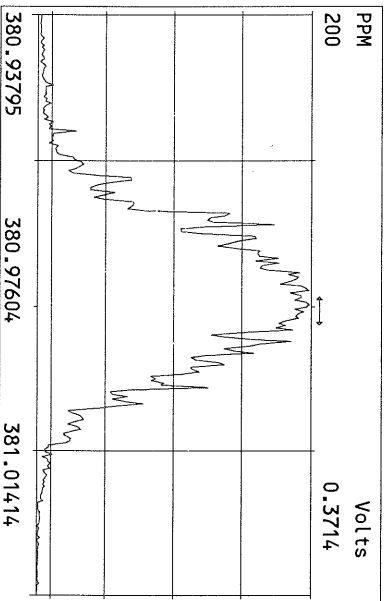
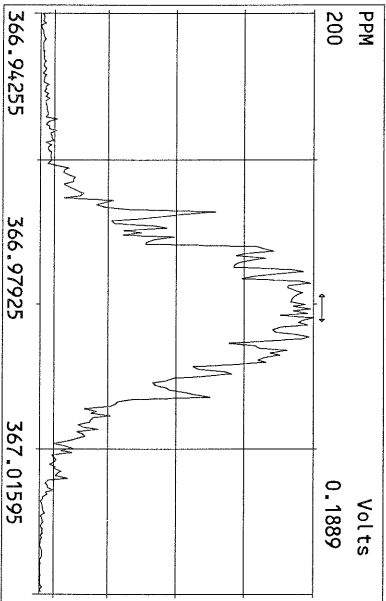
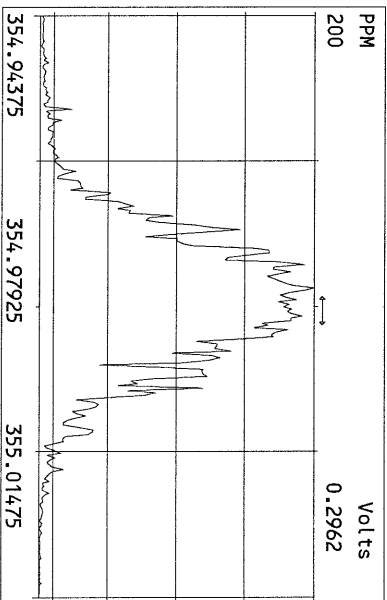
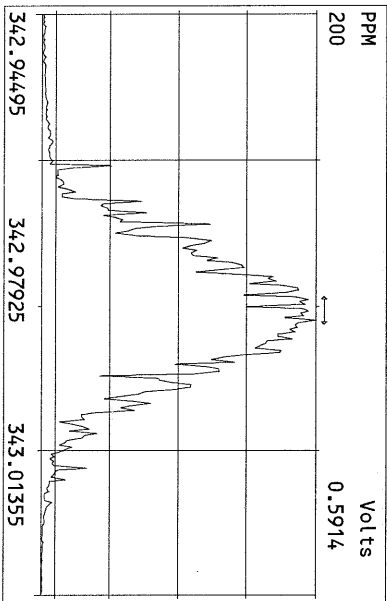
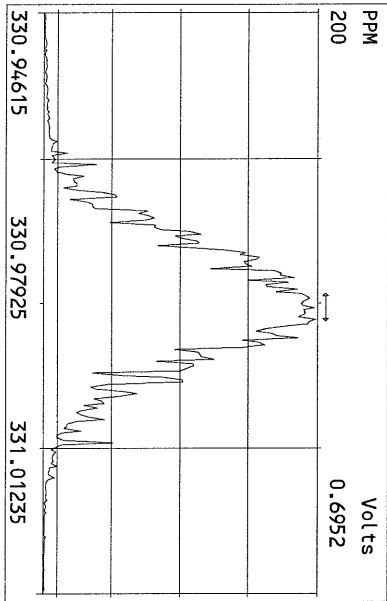
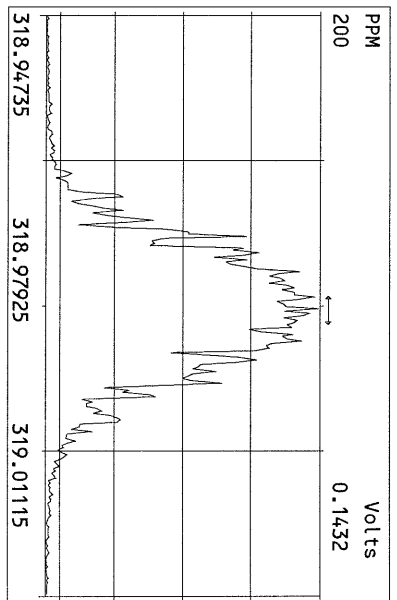
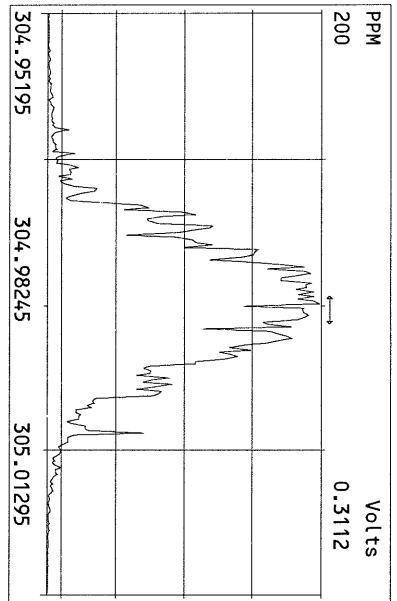
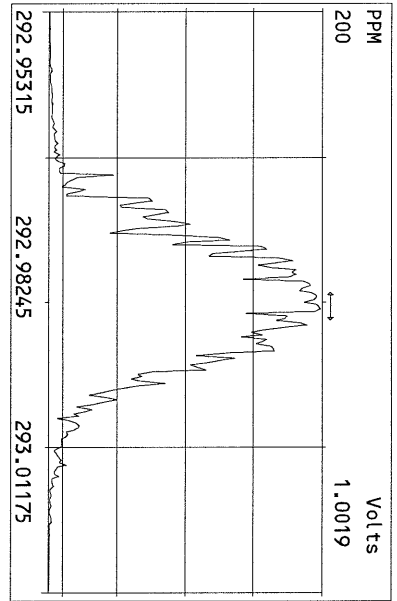


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FAL ID:ST042011A1
100%



File:20APR11A Acq:20-APR-2011 09:25:20 FAL-1 Frontier Analytical Laboratory
317.9389 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0,0,0) Exp:TCDF
FAL ID:ST042011A1
100%





USEPA - ITD

FORM 4A
TCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:


Initial Calibration Date: 2/18/11

Instrument ID: FAL1 GC Column ID: DB225

VER Data Filename: 20APR11A Sam:9 Analysis Date: 20-APR-11 Time: 14:06:38


	M/Z'S FORMING RATIO (1)	ION ABUND. RATIO	QC LIMITS (2)	ACCEPT	CONC. FOUND	CONC. RANGE (ng/mL) (3)
NATIVE ANALYTES						
2,3,7,8-TCDF	M/M+2	0.76	0.65-0.89	y	8.94	8.40 - 12.0 ✓
LABELED COMPOUNDS						
13C-2,3,7,8-TCDF	M/M+2	0.78	0.65-0.89	y	95.1	71.0 - 140 ✓

- (1) See Table 8, Method 1613, for m/z specifications.
- (2) Ion Abundance Ratio Control Limits as specified in Table 9, Method 1613.
- (3) Contract-required concentration range as specified in Table 6A, Method 1613

Analyst:  Date: 4/20/11

FAL ID: ST042011A2 Filename: 20APR11A Sam:9 Acquired: 20-APR-11 14:06:38 ICal: TCDFFAL1-2-18-11
Client ID: 1613 CS3 100511J ConCal: ST042011A1 EndCal: ST042011A2
Results: 6701tcdf GC Column: DB225 Amount: 1.000

Name	Resp	RA	RT	RRF	Conc	Qual	Fac	Noise	DL	#Hom	Rec
2,3,7,8-TCDF	1.45e+07	0.76 y	19:18	1.16	8.94		2.50	-	-	1	
13C-2,3,7,8-TCDF	1.40e+08	0.78 y	19:17	1.05	95.1						95.1
13C-1,2,3,4-TCDF	1.40e+08	0.79 y	16:43	-	79.3						

Analyst:  Date: 4/20/11

Frontier Analytical Laboratory - Acquisition Log

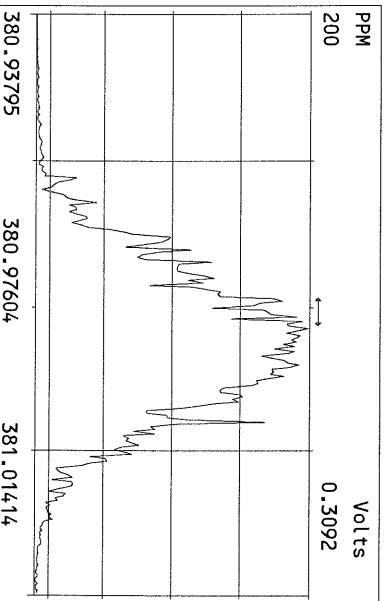
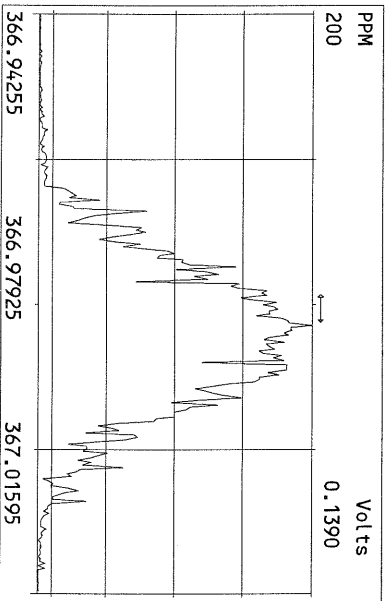
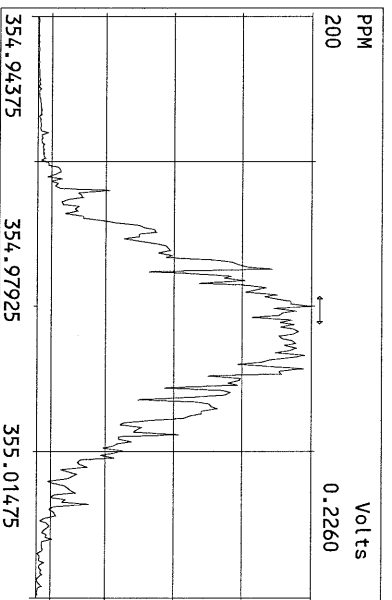
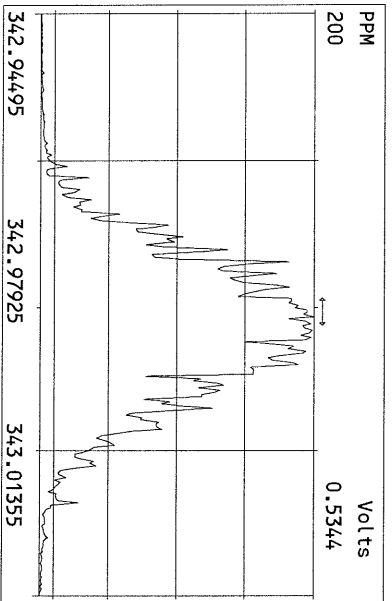
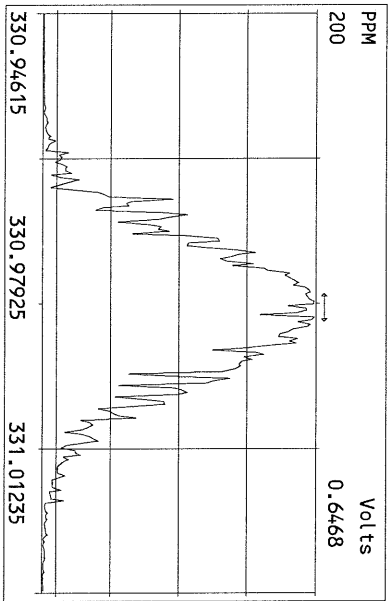
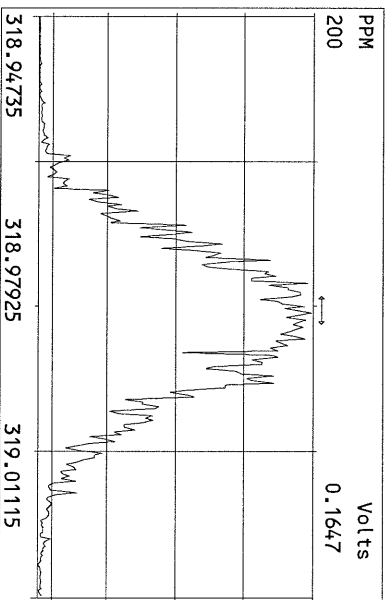
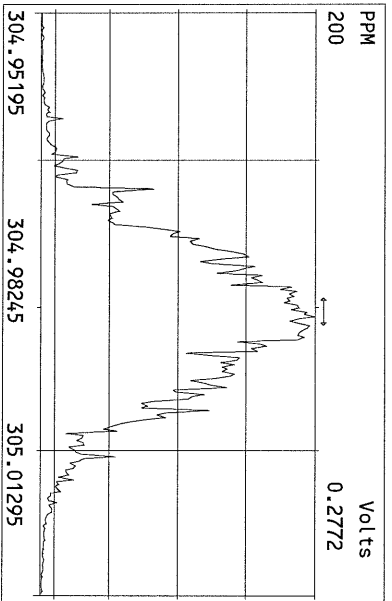
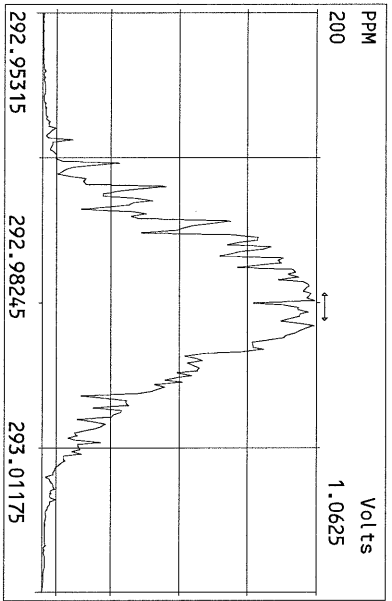
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Data File S	FAL ID	Client ID	Acquired	ConCal	EndCal	Analyst
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20APR11A 2	6701-001-0001-SA	LL-SED1-0-15-032911	20-APR-11 10:00:23	ST042011A1	ST042011A2	TC
20APR11A 3	6701-002-0001-SA	LL-SED2-0-15-032911	20-APR-11 10:35:28	ST042011A1	ST042011A2	TC
20APR11A 4	6701-003-0001-SA	LL-SED3-0-15-032911	20-APR-11 11:10:31	ST042011A1	ST042011A2	TC
20APR11A 5	6701-004-0001-SA	LL-SED4-0-15-032911	20-APR-11 11:45:36	ST042011A1	ST042011A2	TC
20APR11A 6	6701-005-0001-SA	LL-SED1-0-15-032911-D	20-APR-11 12:20:39	ST042011A1	ST042011A2	TC
20APR11A 7	SB042011A1	Solvent Blank	20-APR-11 12:55:44	ST042011A1	ST042011A2	TC
20APR11A 8	SB042011A2	Solvent Blank	20-APR-11 13:30:47	ST042011A1	ST042011A2	TC
20APR11A 9	ST042011A2	1613 CS3 100511J	20-APR-11 14:06:38	ST042011A1	ST042011A2	TC

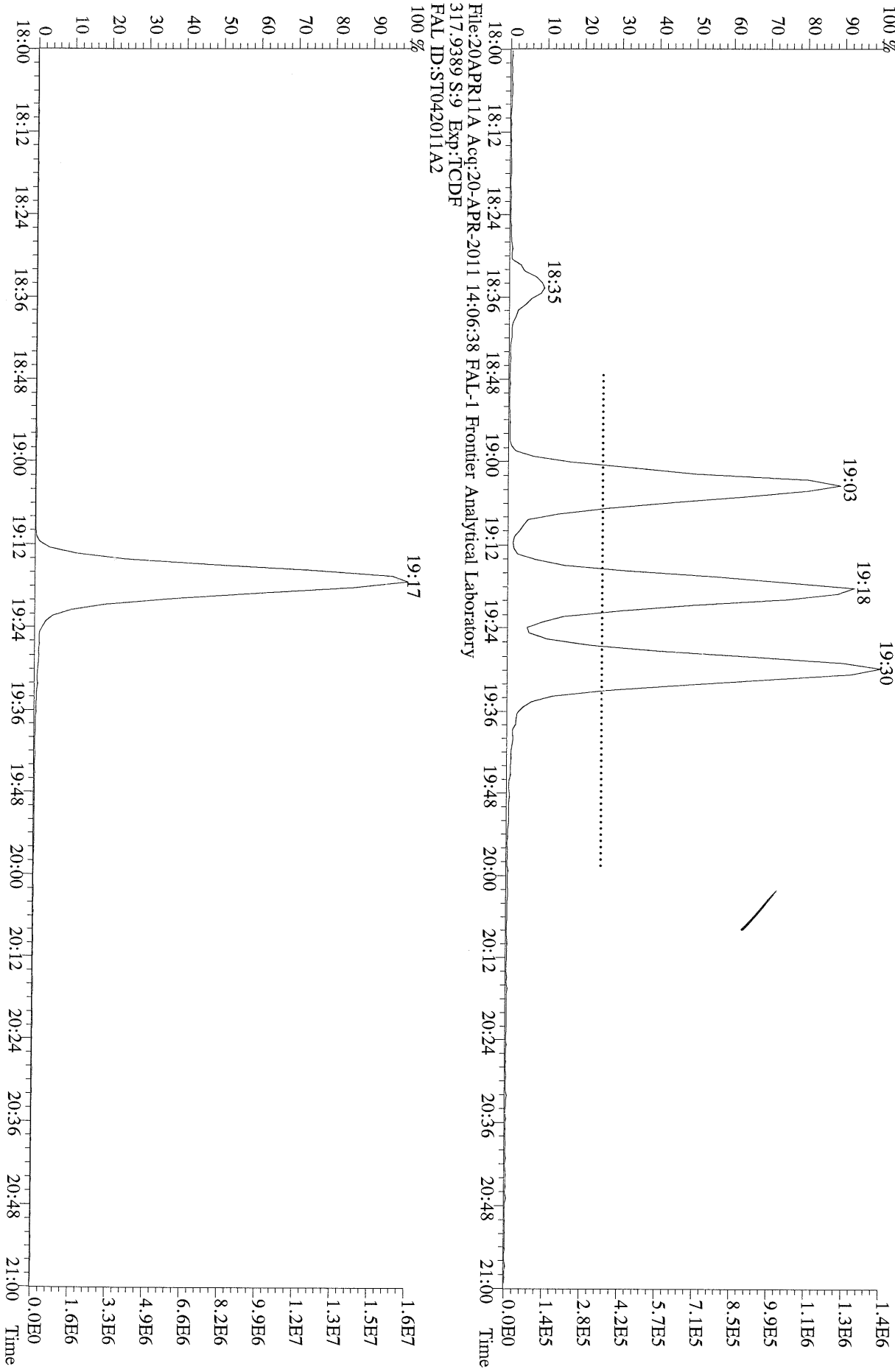
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Data Backed Up: _____

Date: _____

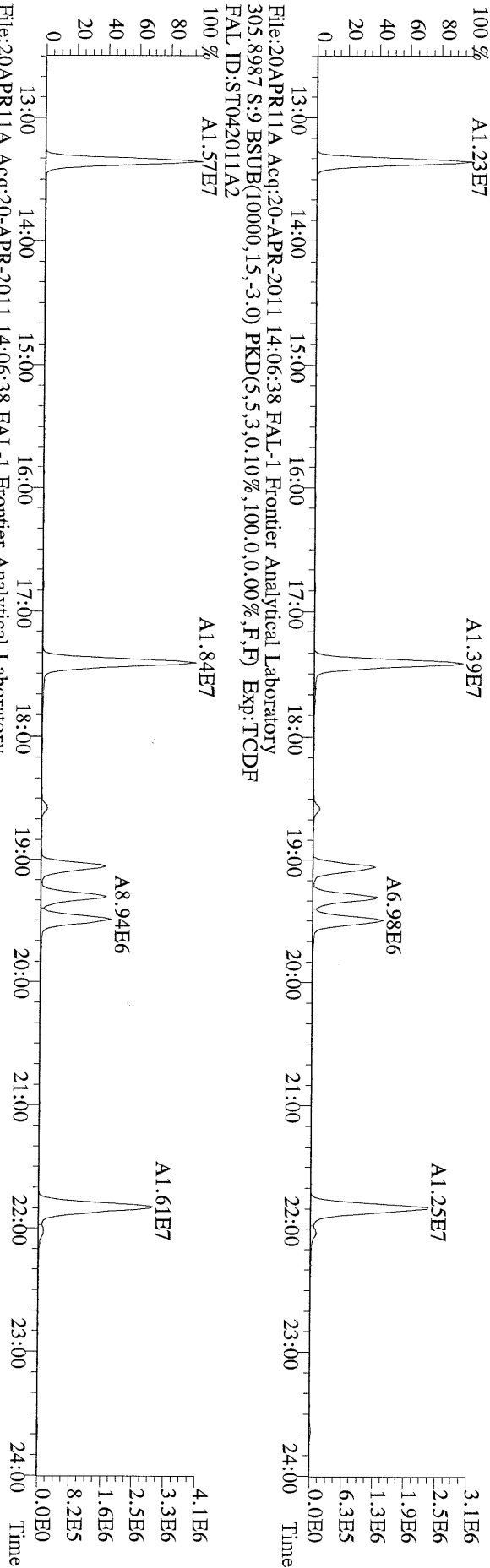


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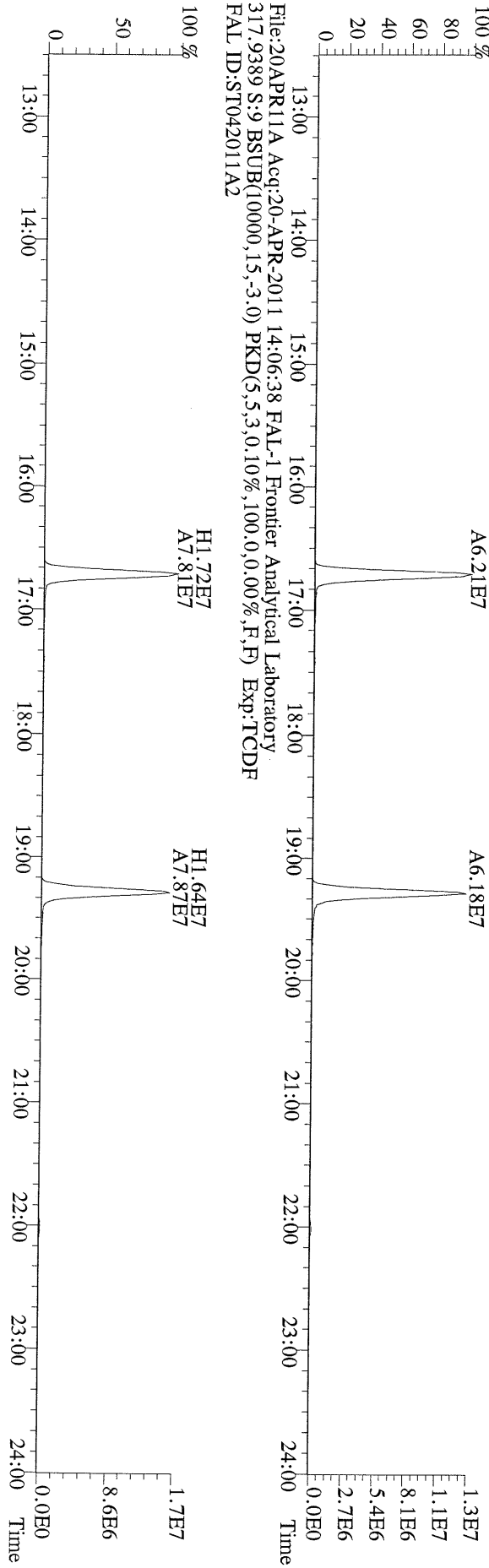


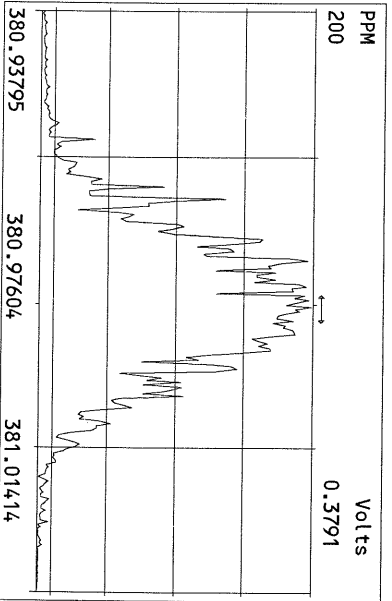
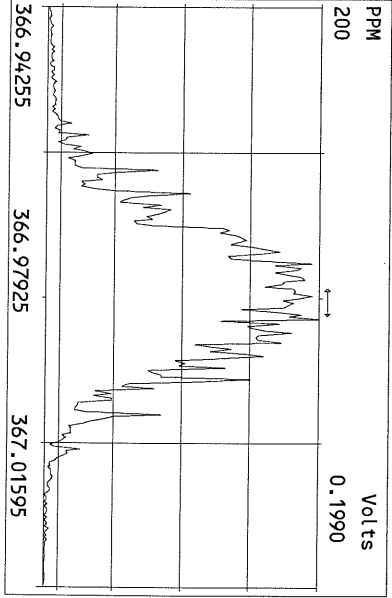
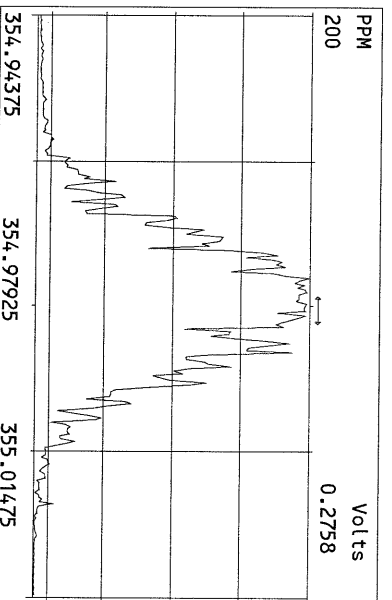
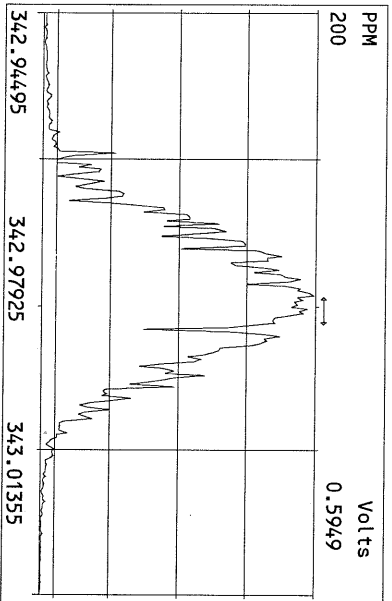
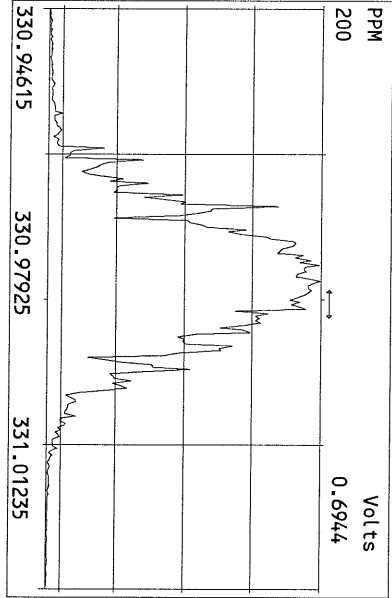
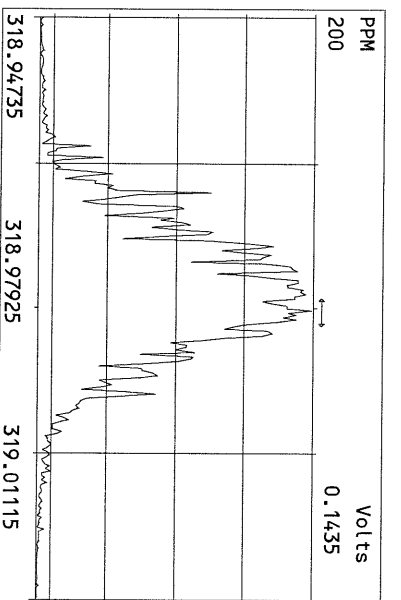
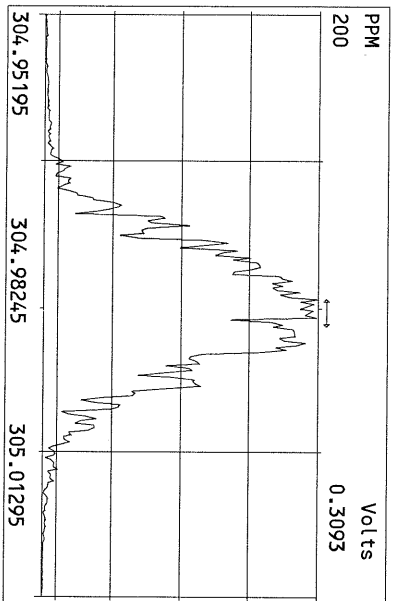
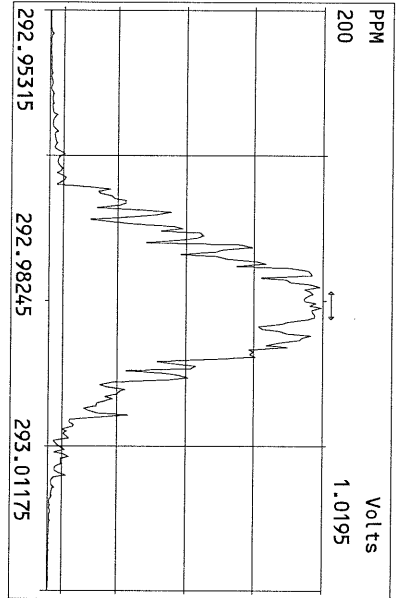
File: 20APR11A Acq: 20-APR-2011 14:06:38 FAL-1 Frontier Analytical Laboratory
317.9389 S:9 Exp: TCDF
FAL ID: ST042011A2

File:20APR11A Acq:20-APR-2011 14:06:38 FAL-1 Frontier Analytical Laboratory
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FAL ID:ST042011A2



File:20APR11A Acq:20-APR-2011 14:06:38 FAL-1 Frontier Analytical Laboratory
315.9419 S:9 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0,0.00%,F,F) Exp:TCDF
FAL ID:ST042011A2





**PCP/Chlorophenols Analysis
Report and Summary QC Forms**

ARI Job ID: SP34, SQ22

Sample ID: LL-SED1-0-15-032911-ER
SAMPLE

Lab Sample ID: SP34J
LIMS ID: 11-6959
Matrix: Water
Data Release Authorized: *AB*
Reported: 04/18/11

QC Report No: SP34-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/01/11
Date Analyzed: 04/15/11 07:31
Instrument/Analyst: ECD1/AAR

Sample Amount: 500 mL
Final Extract Volume: 50 mL
Dilution Factor: 1.00

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	0.25	< 0.25 U

Reported in µg/L (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	90.0%
----------------------	-------

SW8041 CHLOROPHENOLICS SURROGATE RECOVERY SUMMARY

Matrix: Water

QC Report No: SP34-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL

<u>Client ID</u>	<u>TBP</u>	<u>TOT OUT</u>
MB-040111	82.4%	0
LCS-040111	85.6%	0
LCSD-040111	87.6%	0
LL-SED1-0-15-032911-ER	90.0%	0

LCS/MB LIMITS QC LIMITS

(TBP) = 2,4,6-Tribromophenol

(40-130)

(11-156)

Prep Method: SW3510C
Log Number Range: 11-6959 to 11-6959

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LCS-040111
LCS/LCSD

Lab Sample ID: LCS-040111
 LIMS ID: 11-6959
 Matrix: Water
 Data Release Authorized: *[Signature]*
 Reported: 04/18/11

QC Report No: SP34-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted LCS/LCSD: 04/01/11

Sample Amount LCS: 500 mL

LCSD: 500 mL

Date Analyzed LCS: 04/15/11 05:43

Final Extract Volume LCS: 50 mL

LCSD: 04/15/11 06:19

LCSD: 50 mL

Instrument/Analyst LCS: ECD1/AAR

Dilution Factor LCS: 1.00

LCSD: ECD1/AAR

LCSD: 1.00

Analyte	Spike		LCS	LCSD	Spike		RPD
	LCS	Added-LCS	Recovery		Added-LCSD	Recovery	
Pentachlorophenol	2.30	2.50	92.0%	2.40	2.50	96.0%	4.3%

Chlorophenols Surrogate Recovery

	LCS	LCSD
2,4,6-Tribromophenol	85.6%	87.6%

Results reported in µg/L
 RPD calculated using sample concentrations per SW846.

4
CHLOROPHENOL METHOD BLANK SUMMARY

SAMPLE NO.

SP34MBW1

Lab Name: ANALYTICAL RESOURCES, INC
 ARI Job No.: SP34
 Lab Sample ID: SP34MBW1
 Matrix (soil/water) LIQUID
 Sulfur Cleanup (Y/N) Y
 Date Analyzed (1): 04/15/11
 Time Analyzed (1): 0506
 Instrument ID (1): ECD1
 GC Column (1): ZB5 ID: 0.53 (mm)


Client: FLOYD SNIDER
 Project: LORA LAKE SURFACE SEDIMENT
 Lab File ID: 0414A025
 Extraction: (SepF/Cont/Sonc) SW3510C
 Date Extracted: 04/01/11
 Date Analyzed (2): 04/15/11
 Time Analyzed (2): 0506
 Instrument ID (2): ECD1
 GC Column (2): ZB35 ID: 0.53 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE NO. =====	LAB SAMPLE ID =====	DATE ANALYZED 1 =====	DATE ANALYZED 2 =====
01	SP34LCSW1	SP34LCSW1	04/15/11	04/15/11
02	SP34LCSDW1	SP34LCSDW1	04/15/11	04/15/11
03	LL-SED1-0-15	SP34J	04/15/11	04/15/11

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: MB-040111
 METHOD BLANK

Lab Sample ID: MB-040111
 LIMS ID: 11-6959
 Matrix: Water
 Data Release Authorized: 
 Reported: 04/18/11

QC Report No: SP34-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: NA
 Date Received: NA

Date Extracted: 04/01/11
 Date Analyzed: 04/15/11 05:06
 Instrument/Analyst: ECD1/AAR

Sample Amount: 500 mL
 Final Extract Volume: 50 mL
 Dilution Factor: 1.00

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	0.25	< 0.25 U

Reported in µg/L (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	82.4%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
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Sample ID: MC-SED1-0-10-032911
SAMPLE

Lab Sample ID: SP34G
 LIMS ID: 11-6956
 Matrix: Sediment
 Data Release Authorized: *AS*
 Reported: 04/18/11

QC Report No: SP34-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/15/11 16:36
 Instrument/Analyst: ECD1/AAR

Sample Amount: 8.34 g-dry-wt
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: 17.2%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	7.5	< 7.5 U
Reported in µg/kg (ppb)			
Chlorophenol Surrogate Recovery			
	2,4,6-Tribromophenol	82.4%	

ORGANICS ANALYSIS DATA SHEET

PCP by GC/ECD Method SW8041

Page 1 of 1


Sample ID: MC-SED2-0-10-032911

SAMPLE

Lab Sample ID: SP34H

LIMS ID: 11-6957

Matrix: Sediment

Data Release Authorized: 

Reported: 04/18/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Date Extracted: 04/07/11

Date Analyzed: 04/15/11 17:12

Instrument/Analyst: ECD1/AAR

Sample Amount: 8.61 g-dry-wt

Final Extract Volume: 25 mL

Dilution Factor: 1.00

Percent Moisture: 14.8%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	7.3	< 7.3 U

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	65.2%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
Page 1 of 1



Sample ID: MC-SED3-0-10-032911
SAMPLE

Lab Sample ID: SP34I
LIMS ID: 11-6958
Matrix: Sediment
Data Release Authorized: *[Signature]*
Reported: 04/18/11

QC Report No: SP34-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/15/11 17:48
Instrument/Analyst: ECD1/AAR

Sample Amount: 9.62 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 8.1%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	6.5	< 6.5 U

Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	62.8%
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SW8041 CHLOROPHENOLICS SURROGATE RECOVERY SUMMARY

Matrix: Sediment

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling
POS-LL

<u>Client ID</u>	<u>TBP</u>	<u>TOT OUT</u>
MB-040711	80.0%	0
LCS-040711	84.4%	0
MC-SED1-0-10-032911	82.4%	0
MC-SED2-0-10-032911	65.2%	0
MC-SED3-0-10-032911	62.8%	0

LCS/MB LIMITS QC LIMITS

(TBP) = 2,4,6-Tribromophenol

(50-115)

(10-146)

Prep Method: SW3550B

Log Number Range: 11-6956 to 11-6958

ORGANICS ANALYSIS DATA SHEET

PCP by GC/ECD Method SW8041

Page 1 of 1

Sample ID: LCS-040711

LAB CONTROL

Lab Sample ID: LCS-040711

LIMS ID: 11-6956

Matrix: Sediment

Data Release Authorized: *AS*

Reported: 04/18/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Date Extracted: 04/07/11

Date Analyzed: 04/15/11 15:23

Instrument/Analyst: ECD1/AAR

Sample Amount: 10.0 g

Final Extract Volume: 25 mL

Dilution Factor: 1.00

Analyte	Lab Control	Spike Added	Recovery
Pentachlorophenol	58.6	62.5	93.8%

Chlorophenols Surrogate Recovery

2,4,6-Tribromophenol	84.4%
----------------------	-------

Results reported in µg/kg

4
CHLOROPHENOL METHOD BLANK SUMMARY

SAMPLE NO.

SP34MBS1

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

Lab Sample ID: SP34MBS1

Lab File ID: 0414A041

Matrix (soil/water) SOLID

Extraction: (SepF/Cont/Sonc) SW3550C

Sulfur Cleanup (Y/N) Y

Date Extracted: 04/07/11

Date Analyzed (1): 04/15/11

Date Analyzed (2): 04/15/11

Time Analyzed (1): 1447

Time Analyzed (2): 1447

Instrument ID (1): ECD1

Instrument ID (2): ECD1

GC Column (1): STX CLP1 ID: 0.53 (mm)

GC Column (2): STX CLP2 ID: 0.53 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED 1	DATE ANALYZED 2
	=====	=====	=====	=====
01	SP34LCSS1	SP34LCSS1	04/15/11	04/15/11
02	MC-SED1-0-10	SP34G	04/15/11	04/15/11
03	MC-SED2-0-10	SP34H	04/15/11	04/15/11
04	MC-SED3-0-10	SP34I	04/15/11	04/15/11

ORGANICS ANALYSIS DATA SHEET

PCP by GC/ECD Method SW8041

Page 1 of 1


Sample ID: MB-040711

METHOD BLANK

Lab Sample ID: MB-040711

LIMS ID: 11-6956

Matrix: Sediment

Data Release Authorized: 

Reported: 04/18/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: NA

Date Received: NA

Date Extracted: 04/07/11

Date Analyzed: 04/15/11 14:47

Instrument/Analyst: ECD1/AAR

Sample Amount: 10.0 g

Final Extract Volume: 25 mL

Dilution Factor: 1.00

Percent Moisture: NA

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	6.2	< 6.2 U

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	80.0%
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Sample ID: LL-SED1-0-15-032911
SAMPLE

Lab Sample ID: SQ22A
LIMS ID: 11-7355
Matrix: Sediment
Data Release Authorized: *AS*
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/18/11 22:14
Instrument/Analyst: ECD1/AAR

Sample Amount: 3.13 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 69.4%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	20	50 P

Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	105%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED2-0-15-032911
SAMPLE

Lab Sample ID: SQ22B
 LIMS ID: 11-7356
 Matrix: Sediment
 Data Release Authorized: *[Signature]*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/18/11 22:50
 Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	33	< 33 U
Reported in µg/kg (ppb)			
Chlorophenol Surrogate Recovery			
	2,4,6-Tribromophenol	86.4%	

Sample ID: LL-SED2-0-15-032911
REEXTRACT

Lab Sample ID: SQ22B
LIMS ID: 11-8696
Matrix: Sediment
Data Release Authorized: *B*
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface sediment
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/15/11 20:14
Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 10.0
Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	330	< 330 U

Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	83.6%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
Page 1 of 1



Sample ID: LL-SED3-0-15-032911
SAMPLE

Lab Sample ID: SQ22C
LIMS ID: 11-7357
Matrix: Sediment
Data Release Authorized: *AS*
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/19/11 00:39
Instrument/Analyst: ECD1/AAR

Sample Amount: 2.63 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 73.8%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	24	< 24 U

Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	88.0%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED4-0-15-032911
SAMPLE

Lab Sample ID: SQ22D
 LIMS ID: 11-7358
 Matrix: Sediment
 Data Release Authorized: *B*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/19/11 01:15
 Instrument/Analyst: ECD1/AAR

Sample Amount: 2.52 g-dry-wt
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: 75.0%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	25	< 25 U


Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	79.2%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED1-0-15-032911-D
SAMPLE

Lab Sample ID: SQ22E
 LIMS ID: 11-7359
 Matrix: Sediment
 Data Release Authorized: 
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/19/11 01:51
 Instrument/Analyst: ECD1/AAR

Sample Amount: 3.13 g-dry-wt
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: 69.0%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	20	< 20 U
Reported in µg/kg (ppb)			
Chlorophenol Surrogate Recovery			
	2,4,6-Tribromophenol	84.8%	



Sample ID: LL-SED5-0-15-032911
SAMPLE

Lab Sample ID: SQ22F
LIMS ID: 11-7360
Matrix: Sediment
Data Release Authorized: *[Signature]*
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/19/11 02:27
Instrument/Analyst: ECD1/AAR

Sample Amount: 8.60 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 17.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	7.3	33

Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	76.4%
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SW8041 CHLOROPHENOLICS SURROGATE RECOVERY SUMMARY

Matrix: Sediment

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL

<u>Client ID</u>	<u>TBP</u>	<u>TOT OUT</u>
LL-SED1-0-15-032911	105%	0
MB-040711	80.0%	0
LCS-040711	84.4%	0
LL-SED2-0-15-032911	86.4%	0
LL-SED2-0-15-032911 MS	46.8%	0
LL-SED2-0-15-032911 MSD	90.0%	0
LL-SED3-0-15-032911	88.0%	0
LL-SED4-0-15-032911	79.2%	0
LL-SED1-0-15-032911-D	84.8%	0
LL-SED5-0-15-032911	76.4%	0
LL-SED2-0-15-032911-RE	83.6%	0
LL-SED2-0-15-032911-RE MS	68.6%	0
LL-SED2-0-15-032911-RE MSD	122%	0

LCS/MB LIMITS QC LIMITS

(TBP) = 2,4,6-Tribromophenol (50-115) (10-146)

Prep Method: SW3550B
Log Number Range: 11-7355 to 11-8696

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED2-0-15-032911
MS/MSD

Lab Sample ID: SQ22B
 LIMS ID: 11-8696
 Matrix: Sediment
 Data Release Authorized: *[Signature]*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface sediment
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11


Date Extracted MS/MSD: 04/07/11
 Date Analyzed MS: 04/15/11 20:50
 MSD: 04/15/11 21:26
 Instrument/Analyst MS: ECD1/AAR
 MSD: ECD1/AAR
 Percent Moisture: 81.3%

Sample Amount MS: 1.90 g-dry-wt
 MSD: 1.90 g-dry-wt
 Final Extract Volume MS: 25 mL
 MSD: 25 mL
 Dilution Factor MS: 10.0
 MSD: 10.0

Analyte	Sample	MS	Spike Added-MS	MS Recovery	MSD	Spike Added-MSD	MSD Recovery	RPD
Pentachlorophenol	< 329	272	329	82.7%	455	329	138%	50.3%

Results reported in µg/kg
 RPD calculated using sample concentrations per SW846.

Sample ID: LL-SED2-0-15-032911
MATRIX SPIKE

Lab Sample ID: SQ22B
LIMS ID: 11-8696
Matrix: Sediment
Data Release Authorized: 
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface sediment
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/15/11 20:50
Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 10.0
Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	330	---
	Reported in $\mu\text{g}/\text{kg}$ (ppb)		
	Chlorophenol Surrogate Recovery		
	2,4,6-Tribromophenol	68.6%	

ORGANICS ANALYSIS DATA SHEET

PCP by GC/ECD Method SW8041

Page 1 of 1


Sample ID: LL-SED2-0-15-032911

MATRIX SPIKE DUP

Lab Sample ID: SQ22B

LIMS ID: 11-8696

Matrix: Sediment

Data Release Authorized: 

Reported: 04/20/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface sediment

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Date Extracted: 04/07/11

Date Analyzed: 04/15/11 21:26

Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt

Final Extract Volume: 25 mL

Dilution Factor: 10.0

Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	330	---


Reported in $\mu\text{g}/\text{kg}$ (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	122%
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ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED2-0-15-032911
MS/MSD

Lab Sample ID: SQ22B
 LIMS ID: 11-7356
 Matrix: Sediment
 Data Release Authorized: 
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11


Date Extracted MS/MSD: 04/07/11
 Date Analyzed MS: 04/18/11 23:27
 MSD: 04/19/11 00:03
 Instrument/Analyst MS: ECD1/AAR
 MSD: ECD1/AAR
 Percent Moisture: 81.3%

Sample Amount MS: 1.90 g-dry-wt
 MSD: 1.90 g-dry-wt
 Final Extract Volume MS: 25 mL
 MSD: 25 mL
 Dilution Factor MS: 1.00
 MSD: 1.00

Analyte	Sample	MS	Spike		MSD	Spike		RPD
			Added-MS	MS Recovery		Added-MSD	MSD Recovery	
Pentachlorophenol	< 32.9	134	329	40.7%	281	329	85.4%	70.8%

Results reported in µg/kg
 RPD calculated using sample concentrations per SW846.

Sample ID: LL-SED2-0-15-032911
MATRIX SPIKE

Lab Sample ID: SQ22B
LIMS ID: 11-7356
Matrix: Sediment
Data Release Authorized: 
Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Date Extracted: 04/07/11
Date Analyzed: 04/18/11 23:27
Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	33	---
Reported in $\mu\text{g}/\text{kg}$ (ppb)			
Chlorophenol Surrogate Recovery			
	2,4,6-Tribromophenol	46.8%	

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LL-SED2-0-15-032911
MATRIX SPIKE DUP

Lab Sample ID: SQ22B
 LIMS ID: 11-7356
 Matrix: Sediment
 Data Release Authorized: *AS*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/19/11 00:03
 Instrument/Analyst: ECD1/AAR

Sample Amount: 1.90 g-dry-wt
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: 81.3%

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	33	---
Reported in µg/kg (ppb)			
Chlorophenol Surrogate Recovery			
	2,4,6-Tribromophenol	90.0%	

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
 Page 1 of 1

Sample ID: LCS-040711
LAB CONTROL

Lab Sample ID: LCS-040711
 LIMS ID: 11-7356
 Matrix: Sediment
 Data Release Authorized: *RB*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

Date Extracted: 04/07/11
 Date Analyzed: 04/15/11 15:23
 Instrument/Analyst: ECD1/AAR

Sample Amount: 10.0 g
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00

Analyte	Lab Control	Spike Added	Recovery
Pentachlorophenol	58.6	62.5	93.8%

Chlorophenols Surrogate Recovery

2,4,6-Tribromophenol	84.4%
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Results reported in µg/kg

4
CHLOROPHENOL METHOD BLANK SUMMARY

SAMPLE NO.

SP34MBS1

Lab Name: ANALYTICAL RESOURCES, INC	Client: FLOYD SNIDER
ARI Job No.: SQ22	Project: LORA LAKE SURFACE SEDIMENT
Lab Sample ID: SP34MBS1	Lab File ID: 0414A041
Matrix (soil/water) SOLID	Extraction: (SepF/Cont/Sonc) SW3550C
Sulfur Cleanup (Y/N) Y	Date Extracted: 04/07/11
Date Analyzed (1): 04/15/11	Date Analyzed (2): 04/15/11
Time Analyzed (1): 1447	Time Analyzed (2): 1447
Instrument ID (1): ECD1	Instrument ID (2): ECD1
GC Column (1): STX CLP1 ID: 0.53 (mm)	GC Column (2): STX CLP2 ID: 0.53 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED 1	DATE ANALYZED 2
=====				
01	SP34LCSS1	SP34LCSS1	04/15/11	04/15/11
02	LL-SED1-0-15	SQ22A	04/18/11	04/19/11
03	LL-SED2-0-15	SQ22B	04/18/11	04/18/11
04	LL-SED2-0-15	SQ22BMS	04/18/11	04/18/11
05	LL-SED2-0-15	SQ22BMSD	04/19/11	04/19/11
06	LL-SED3-0-15	SQ22C	04/19/11	04/19/11
07	LL-SED4-0-15	SQ22D	04/19/11	04/19/11
08	LL-SED1-0-15	SQ22A	04/18/11	04/19/11
09	LL-SED2-0-15	SQ22F	04/19/11	04/19/11

ORGANICS ANALYSIS DATA SHEET
PCP by GC/ECD Method SW8041
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Sample ID: MB-040711
METHOD BLANK

Lab Sample ID: MB-040711
 LIMS ID: 11-7356
 Matrix: Sediment
 Data Release Authorized: *AB*
 Reported: 04/20/11

QC Report No: SQ22-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: NA
 Date Received: NA

Date Extracted: 04/07/11
 Date Analyzed: 04/15/11 14:47
 Instrument/Analyst: ECD1/AAR

Sample Amount: 10.0 g
 Final Extract Volume: 25 mL
 Dilution Factor: 1.00
 Percent Moisture: NA

CAS Number	Analyte	RL	Result
87-86-5	Pentachlorophenol	6.2	< 6.2 U

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

2,4,6-Tribromophenol	80.0%
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6D
 CHLOROPHENOL INITIAL CALIBRATION
 RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB5 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/14/11

COMPOUND	RT OF STANDARDS						MEAN RT	RT WINDOW	
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		FROM	TO
Pentachlorophenol	20.99	20.99	20.99	20.99	20.99	20.99	20.99	20.92	21.06
2,4,6-Trichloropheno	13.10	13.10	13.10	13.09	13.09	13.09	13.10	13.02	13.16
2,3,6-Trichloropheno	14.09	14.09	14.09	14.09	14.09	14.09	14.09	14.02	14.16
2,4,5-Trichloropheno	15.84	15.84	15.84	15.84	15.84	15.84	15.84	15.77	15.91
2,3,4-Trichloropheno	17.35	17.35	17.35	17.35	17.35	17.34	17.35	17.28	17.42
2,3,5,6-Tetrachlorop	17.15	17.15	17.15	17.15	17.15	17.15	17.15	17.08	17.22
2,3,4,5-Tetrachlorop	20.15	20.15	20.15	20.15	20.15	20.15	20.15	20.08	20.22
2,4-Dichlorophenol	12.55	12.55	12.55	12.55	12.55	12.55	12.55	12.48	12.62
2,4,6-Tribromophenol	18.59	18.59	18.59	18.59	18.59	18.59	18.59	18.52	18.66

6D
 CHLOROPHENOL INITIAL CALIBRATION
 RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB35 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/14/11

COMPOUND	RT OF STANDARDS						MEAN RT	RT WINDOW	
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		FROM	TO
Pentachlorophenol	22.96	22.96	22.96	22.96	22.96	22.96	22.96	22.89	23.03
2,4,6-Trichloropheno	14.31	14.31	14.31	14.30	14.30	14.30	14.30	14.23	14.37
2,3,6-Trichloropheno	15.55	15.55	15.55	15.55	15.55	15.55	15.55	15.48	15.62
2,4,5-Trichloropheno	17.47	17.47	17.47	17.47	17.47	17.47	17.47	17.40	17.54
2,3,4-Trichloropheno	19.02	19.02	19.02	19.02	19.02	19.02	19.02	18.95	19.09
2,3,5,6-Tetrachlorop	18.81	18.81	18.81	18.81	18.81	18.81	18.81	18.74	18.88
2,3,4,5-Tetrachlorop	22.07	22.08	22.08	22.07	22.08	22.07	22.08	22.00	22.14
2,4-Dichlorophenol	13.81	13.82	13.82	13.81	13.81	13.81	13.81	13.74	13.88
2,4,6-Tribromophenol	20.93	20.93	20.93	20.93	20.93	20.93	20.93	20.86	21.00

6E
 CHLOROPHENOL INITIAL CALIBRATION
 CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB5 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/14/11

COMPOUND	CALIBRATION FACTORS						R ² / %RSD	CT
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		
Pentachlorophenol	18330	17330	16468	15138	14544	13342	11.7	A
2,4,6-Trichlorophenol	11745	11226	10426	9411	8773	7987	14.7	A
2,3,6-Trichlorophenol	11524	10670	9838	8910	8304	7526	15.8	A
2,4,5-Trichlorophenol	6896	6582	6014	5318	4751	4147	19.0	A
2,3,4-Trichlorophenol	8312	7672	7053	6262	5821	5184	17.5	A
2,3,5,6-Tetrachloroph	15657	14796	14050	13099	12511	11466	11.3	A
2,3,4,5-Tetrachloroph	12930	11414	10586	9492	8889	8103	17.3	A
2,4-Dichlorophenol	890	791	692	575	492	419	0.9994	Q
2,4,6-Tribromophenol	13828	13279	12729	11921	11832	11073	8.2	A
AVE RSD							16.0	

CT stands for Curve Types:

- A Indicates an Average Response Factor Curve
- L Indicates a Linear Curve
- Q Indicates a Quadratic Curve

CALIBRATION FILES

LVL 1: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A010.d
 LVL 2: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A011.d
 LVL 3: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A012.d
 LVL 4: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A009.d
 LVL 5: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A013.d
 LVL 6: /chem2/ecd1.i/PCP20110414.b/ical-1.b/0414A014.d

6E
 CHLOROPHENOL INITIAL CALIBRATION
 CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB35 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/14/11

COMPOUND	CALIBRATION FACTORS						R ² / %RSD	CT
	LVL 1	LVL 2	LVL 3	LVL 4	LVL 5	LVL 6		
Pentachlorophenol	34545	28057	26029	23780	22068	19893	0.9999	Q
2,4,6-Trichlorophenol	16228	14895	13845	12744	11452	10173	16.9	A
2,3,6-Trichlorophenol	16178	14843	13791	12572	11496	10026	17.1	A
2,4,5-Trichlorophenol	9097	8460	7618	6667	6237	5207	0.9996	Q
2,3,4-Trichlorophenol	11335	10280	9322	8303	7534	6539	0.9999	Q
2,3,5,6-Tetrachloroph	23914	22454	20497	18869	17642	15789	15.3	A
2,3,4,5-Tetrachloroph	21489	17304	15765	13972	12707	11236	0.9998	Q
2,4-Dichlorophenol	1085	896	764	644	548	460	0.9996	Q
2,4,6-Tribromophenol	21999	20210	19161	17970	17054	15618	12.2	A
AVE RSD							19.7	

CT stands for Curve Types:

- A Indicates an Average Response Factor Curve
- L Indicates a Linear Curve
- Q Indicates a Quadratic Curve

CALIBRATION FILES

- LVL 1: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A010.d
- LVL 2: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A011.d
- LVL 3: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A012.d
- LVL 4: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A009.d
- LVL 5: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A013.d
- LVL 6: /chem2/ecd1.i/PCP20110414.b/ical-2.b/0414A014.d

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB5 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :0430

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	20.99	20.92	21.06	26.2	25.0	4.8
2,4,6-Trichlorophenol	13.10	13.02	13.16	26.0	25.0	4.0
2,3,6-Trichlorophenol	14.09	14.02	14.16	25.8	25.0	3.2
2,4,5-Trichlorophenol	15.84	15.77	15.91	25.8	25.0	3.2
2,3,4-Trichlorophenol	17.35	17.28	17.42	25.7	25.0	2.8
2,3,5,6-Tetrachlorophenol	17.15	17.08	17.22	26.4	25.0	5.6
2,3,4,5-Tetrachlorophenol	20.15	20.08	20.22	25.3	25.0	1.2
2,4-Dichlorophenol	12.55	12.48	12.62	277	250	10.8
2,4,6-Tribromophenol (surr	18.59	18.52	18.66	26.5	25.0	6.0

AVERAGE %D = 4.6

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB35 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :0430

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	22.96	22.89	23.03	26.4	25.0	5.6
2,4,6-Trichlorophenol	14.31	14.23	14.37	25.3	25.0	1.2
2,3,6-Trichlorophenol	15.55	15.48	15.62	25.0	25.0	0.0
2,4,5-Trichlorophenol	17.47	17.40	17.54	25.3	25.0	1.2
2,3,4-Trichlorophenol	19.02	18.95	19.09	26.1	25.0	4.4
2,3,5,6-Tetrachlorophenol	18.81	18.74	18.88	24.9	25.0	-0.4
2,3,4,5-Tetrachlorophenol	22.08	22.00	22.14	26.1	25.0	4.4
2,4-Dichlorophenol	13.82	13.74	13.88	26.4	25.0	5.6
2,4,6-Tribromophenol (surr)	20.93	20.86	21.00	25.2	25.0	0.8

AVERAGE %D = 2.6

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB5 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1032

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	20.99	20.92	21.06	28.4	25.0	13.6
2,4,6-Trichlorophenol	13.09	13.02	13.16	27.9	25.0	11.6
2,3,6-Trichlorophenol	14.09	14.02	14.16	27.6	25.0	10.4
2,4,5-Trichlorophenol	15.83	15.77	15.91	27.4	25.0	9.6
2,3,4-Trichlorophenol	17.34	17.28	17.42	27.4	25.0	9.6
2,3,5,6-Tetrachlorophenol	17.14	17.08	17.22	28.4	25.0	13.6
2,3,4,5-Tetrachlorophenol	20.14	20.08	20.22	27.3	25.0	9.2
2,4-Dichlorophenol	12.55	12.48	12.62	304	250	21.6
2,4,6-Tribromophenol (surr)	18.58	18.52	18.66	28.4	25.0	13.6

AVERAGE %D = 12.5

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB35 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1032

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	22.96	22.89	23.03	27.7	25.0	10.8
2,4,6-Trichlorophenol	14.30	14.23	14.37	26.3	25.0	5.2
2,3,6-Trichlorophenol	15.55	15.48	15.62	25.7	25.0	2.8
2,4,5-Trichlorophenol	17.46	17.40	17.54	27.8	25.0	11.2
2,3,4-Trichlorophenol	19.01	18.95	19.09	27.5	25.0	10.0
2,3,5,6-Tetrachlorophenol	18.80	18.74	18.88	25.8	25.0	3.2
2,3,4,5-Tetrachlorophenol	22.07	22.00	22.14	27.5	25.0	10.0
2,4-Dichlorophenol	13.81	13.74	13.88	278	250	11.2
2,4,6-Tribromophenol (surr	20.92	20.86	21.00	26.7	25.0	6.8

AVERAGE %D = 7.9

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP1 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1410

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	20.99	20.92	21.06	23.6	25.0	-5.6
2,4,6-Trichlorophenol	13.09	13.02	13.16	23.5	25.0	-6.0
2,3,6-Trichlorophenol	14.08	14.02	14.16	23.1	25.0	-7.6
2,4,5-Trichlorophenol	15.83	15.77	15.91	23.0	25.0	-8.0
2,3,4-Trichlorophenol	17.34	17.28	17.42	23.2	25.0	-7.2
2,3,5,6-Tetrachlorophenol	17.14	17.08	17.22	23.8	25.0	-4.8
2,3,4,5-Tetrachlorophenol	20.14	20.08	20.22	23.0	25.0	-8.0
2,4-Dichlorophenol	12.55	12.48	12.62	259	250	3.6
2,4,6-Tribromophenol (surr	18.58	18.52	18.66	23.7	25.0	-5.2

AVERAGE %D = 6.2

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP2 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1410

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	22.95	22.89	23.03	22.6	25.0	-9.6
2,4,6-Trichlorophenol	14.30	14.23	14.37	21.7	25.0	-13.2
2,3,6-Trichlorophenol	15.55	15.48	15.62	21.1	25.0	-15.6
2,4,5-Trichlorophenol	17.46	17.40	17.54	22.1	25.0	-11.6
2,3,4-Trichlorophenol	19.01	18.95	19.09	22.8	25.0	-8.8
2,3,5,6-Tetrachlorophenol	18.80	18.74	18.88	21.9	25.0	-12.4
2,3,4,5-Tetrachlorophenol	22.07	22.00	22.14	22.5	25.0	-10.0
2,4-Dichlorophenol	13.81	13.74	13.88	228	250	-8.8
2,4,6-Tribromophenol (surr)	20.92	20.86	21.00	22.2	25.0	-11.2

AVERAGE %D = 11.2

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP1 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1901

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	20.98	20.92	21.06	24.9	25.0	-0.4
2,4,6-Trichlorophenol	13.09	13.02	13.16	24.5	25.0	-2.0
2,3,6-Trichlorophenol	14.08	14.02	14.16	24.0	25.0	-4.0
2,4,5-Trichlorophenol	15.83	15.77	15.91	24.1	25.0	-3.6
2,3,4-Trichlorophenol	17.34	17.28	17.42	24.6	25.0	-1.6
2,3,5,6-Tetrachlorophenol	17.14	17.08	17.22	25.1	25.0	0.4
2,3,4,5-Tetrachlorophenol	20.14	20.08	20.22	24.1	25.0	-3.6
2,4-Dichlorophenol	12.54	12.48	12.62	274	250	9.6
2,4,6-Tribromophenol (surr	18.58	18.52	18.66	25.0	25.0	0.0

AVERAGE %D = 2.8

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP2 ID: 0.53 (mm)

Init. Calib. Date(s): 04/14/11 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/15/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1901

PCP MIX COMPOUND	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Pentachlorophenol	22.95	22.89	23.03	23.7	25.0	-5.2
2,4,6-Trichlorophenol	14.30	14.23	14.37	22.6	25.0	-9.6
2,3,6-Trichlorophenol	15.54	15.48	15.62	22.0	25.0	-12.0
2,4,5-Trichlorophenol	17.46	17.40	17.54	24.5	25.0	-2.0
2,3,4-Trichlorophenol	19.01	18.95	19.09	23.6	25.0	-5.6
2,3,5,6-Tetrachlorophenol	18.80	18.74	18.88	22.6	25.0	-9.6
2,3,4,5-Tetrachlorophenol	22.07	22.00	22.14	24.2	25.0	-3.2
2,4-Dichlorophenol	13.81	13.74	13.88	237	250	-5.2
2,4,6-Tribromophenol (surr	20.92	20.86	21.00	23.3	25.0	-6.8

AVERAGE %D = 6.6

7E
 CHLOROPHENOLS CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC. Client: FLOYD SNIDER
 ARI Job No.: SQ22 Project: LORA LAKE SURFACE SEDIMENT
 GC Column: STX CLP1 ID# 100000
 Init. Calib. Date(s): 04/14/11

Client Sample No. (PCP): Date Analyzed :04/18/11
 Lab Sample ID (PCP): PCP-100 Time Analyzed :2138

PCP STX C.I. Name	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
1,2,4-trichloropheno	13.02	20.92	21.06	28.0	25.0	12.0
2,4,6-trichloropheno	13.16	13.02	13.16	27.4	25.0	9.6
1,2,4,6-tetrachloropheno	14.02	14.02	14.16	27.2	25.0	8.8
2,4,6-trichloropheno	15.77	15.77	15.91	27.6	25.0	10.4
2,4,6-trichloropheno	17.28	17.28	17.42	27.3	25.0	9.2
2,3,4,6-tetrachlorop	17.08	17.08	17.22	28.1	25.0	12.4
1,2,4-trichloropheno	20.08	20.08	20.22	27.2	25.0	8.8
2,4-Dichlorophenol	12.48	12.48	12.62	295	250	18.0
2,4-Dichloropheno	18.52	18.52	18.66	28.1	25.0	12.4

AVERAGE %D = 11.3

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESEARCH SERVICES, INC

Client: FLOYD SNIDER

ART Job No.: S922

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP2 150m 0.53mm ID

Init. Cal. Date(s): 04/11/11

Client Sample No. (PCP):

Date Analyzed : 04/18/11

Lab Sample ID (PCP):

Time Analyzed : 2138

PCP NAME	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
Phenol	22.89	22.89	23.03	25.6	25.0	2.4
2,3-dichlorophenol	14.23	14.23	14.37	25.0	25.0	0.0
2,4-dichlorophenol	15.48	15.48	15.62	24.9	25.0	-0.4
2,5-dichlorophenol	17.40	17.40	17.54	25.3	25.0	1.2
2,3,4-trichlorophenol	19.02	18.95	19.09	26.0	25.0	4.0
2,3,5-trichlorophenol	18.81	18.74	18.88	24.7	25.0	-1.2
2,4,5-trichlorophenol	22.03	22.00	22.14	25.2	25.0	0.8
2,4-dichlorophenol	13.74	13.74	13.88	265	250	6.0
2,6-dichlorophenol	20.86	20.86	21.00	25.4	25.0	1.6

AVERAGE %D = 2.0

7E
 CHLOROPHENOLS CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL LABORATORIES, INC

Client: FLOYD SNIDER

ART # : 0014 SQ22

Project: LORA LAKE SURFACE SEDIMENT

GC Column: HP-1 CLP1 1.00E+05

Init. Calib. Date(s): 04/14/11

Client Sample No. (PCP):

Date Analyzed :04/19/11

Lab Sample ID (PCP):

Time Analyzed :0340

	RT	RT WINDOW		CALC AMOUNT	NOM AMOUNT	%D
		FROM	TO			
1,2,4-trichloropheno	13.02	20.91	21.06	24.6	25.0	-1.6
2,3,4-trichloropheno	14.02	13.02	13.16	27.4	25.0	9.6
2,3,5-trichloropheno	15.77	14.02	14.16	27.2	25.0	8.8
2,3,4-trichloropheno	17.35	15.77	15.91	27.2	25.0	8.8
2,3,5-trichloropheno	17.35	17.28	17.42	26.2	25.0	4.8
2,3,4,6-tetrachlorop	17.08	17.08	17.22	27.3	25.0	9.2
2,3,5-trichlorop	20.08	20.08	20.22	24.0	25.0	-4.0
2,4-dichloropheno	12.42	12.42	12.62	293	250	17.2
2,4-dichloropheno	18.52	18.52	18.66	26.1	25.0	4.4

AVERAGE %D = 7.6

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL PRODUCTS, INC. Client: FLOYD SNIDER
 ARI Job No.: S922 Project: LORA LAKE SURFACE SEDIMENT
 GC Column: RTX CLP2 (100% Ph) (53 ft x 0.25 mm)
 Init. Cal. Date(s): 11/14/11

Client Sample No. (PCP): Date Analyzed : 04/19/11
 Lab Sample ID (PCP): P-100 Time Analyzed : 0340

PCP Name	RT	RT FROM	RT TO	CALC AMOUNT	NOM AMOUNT	%D
2,4,6-trichlorophenol	22.87	22.89	23.03	23.1	25.0	-7.6
2,4,5-trichlorophenol	14.21	14.21	14.37	24.6	25.0	-1.6
2,4,6-trichlorophenol	15.48	15.48	15.62	24.7	25.0	-1.2
2,4,5-trichlorophenol	17.40	17.40	17.54	24.6	25.0	-1.6
2,3,4-trichlorophenol	19.03	18.95	19.09	23.9	25.0	-4.4
2,3,4,6-tetrachlorophenol	18.80	18.74	18.88	23.8	25.0	-4.8
2,3,4,5-tetrachlorophenol	22.00	22.00	22.14	22.1	25.0	-11.6
2,4,6-trichlorophenol	13.7	13.7	13.88	257	250	2.8
2,4,6-trichlorophenol	20.84	20.8	21.00	23.5	25.0	-6.0

AVERAGE %D = 4.6

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB5

ID: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION				
S1 : 18.59				
CLIENT	LAB	DATE	TIME	S1
SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT #
=====	=====	=====	=====	=====
01	PCPD	04/14/11	1927	18.59
02	PCPA	04/14/11	2003	18.59
03	PCPB	04/14/11	2040	18.59
04	PCPC	04/14/11	2116	18.59
05	PCPE	04/14/11	2152	18.59
06	PCPF	04/14/11	2228	18.59
07	ZZZZZ	04/14/11	2304	18.59
08	PCP CCAL	04/15/11	0430	18.59
09	SP34MBW1	04/15/11	0506	18.59
10	SP34LCSW1	04/15/11	0543	18.59
11	SP34LCSDW1	04/15/11	0619	18.59
12	ZZZZZ	04/15/11	0655	18.59
13	LL-SED1-0-15	04/15/11	0731	18.59
14	PCP CCAL	04/15/11	1032	18.58

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: ZB35

ID: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION					
S1 : 20.93					
CLIENT	LAB	DATE	TIME	S1	
SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT	#
=====					
01	PCPD	04/14/11	1927	20.93	
02	PCPA	04/14/11	2003	20.93	
03	PCPB	04/14/11	2040	20.93	
04	PCPC	04/14/11	2116	20.93	
05	PCPE	04/14/11	2152	20.93	
06	PCPF	04/14/11	2228	20.93	
07	ZZZZZ	04/14/11	2304	20.93	
08	PCP CCAL	04/15/11	0430	20.93	
09	SP34MBW1	04/15/11	0506	20.93	
10	SP34LCSW1	04/15/11	0543	20.93	
11	SP34LCSDW1	04/15/11	0619	20.93	
12	ZZZZZ	04/15/11	0655	20.93	
13	LL-SED1-0-15	04/15/11	0731	20.93	
14	PCP CCAL	04/15/11	1032	20.92	

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP1 ID: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION					
S1 : 18.59					
CLIENT	LAB	DATE	TIME	SI	
SAMPLE NO.	SAMPLE ID	ANALYZED	ANALYZED	RT	#
=====	=====	=====	=====	=====	=====
01	PCPD	04/14/11	1927	18.59	
02	PCPA	04/14/11	2003	18.59	
03	PCPB	04/14/11	2040	18.59	
04	PCPC	04/14/11	2116	18.59	
05	PCPE	04/14/11	2152	18.59	
06	PCPF	04/14/11	2228	18.59	
07	ZZZZZ	04/14/11	2304	18.59	
08	PCP CCAL	04/15/11	1410	18.58	
09	SP34MBS1	04/15/11	1447	18.58	
10	SP34LCSS1	04/15/11	1523	18.58	
11	ZZZZZ	04/15/11	1559	18.58	
12	MC-SED1-0-10	04/15/11	1636	18.58	
13	MC-SED2-0-10	04/15/11	1712	18.58	
14	MC-SED3-0-10	04/15/11	1748	18.58	
15	ZZZZZ	04/15/11	1825	18.58	
16	PCP CCAL	04/15/11	1901	18.58	

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SP34

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP2 ID: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION S1 : 20.93					
	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT #
	=====	=====	=====	=====	=====
01		PCPD	04/14/11	1927	20.93
02		PCPA	04/14/11	2003	20.93
03		PCPB	04/14/11	2040	20.93
04		PCPC	04/14/11	2116	20.93
05		PCPE	04/14/11	2152	20.93
06		PCPF	04/14/11	2228	20.93
07	ZZZZZ	ZZZZZ	04/14/11	2304	20.93
08		PCP CCAL	04/15/11	1410	20.92
09	SP34MBS1	SP34MBS1	04/15/11	1447	20.92
10	SP34LCSS1	SP34LCSS1	04/15/11	1523	20.92
11	ZZZZZ	ZZZZZ	04/15/11	1559	20.92
12	MC-SED1-0-10	SP34G	04/15/11	1636	20.92
13	MC-SED2-0-10	SP34H	04/15/11	1712	20.92
14	MC-SED3-0-10	SP34I	04/15/11	1748	20.92
15	ZZZZZ	ZZZZZ	04/15/11	1825	20.92
16		PCP CCAL	04/15/11	1901	20.92

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SQ22

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP1 ID: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION RT : 18.59					
CURRENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT	#
01		04/14/11	1927	18.59	
02		04/14/11	2003	18.59	
03		04/14/11	2040	18.59	
04		04/14/11	2116	18.59	
05		04/14/11	2152	18.59	
06		04/14/11	2228	18.59	
07	ZZZZZ	04/14/11	2304	18.59	
08	INTERNAL	04/15/11	1410	18.58	
09	ST-MBS1	04/15/11	1447	18.58	
10	34LCSS1	04/15/11	1523	18.58	
11	INTERNAL	04/15/11	1901	18.58	
12	ZZZZZ	04/18/11	2025	18.59	
13	ZZZZZ	04/18/11	2102	18.59	
14	INTERNAL	04/18/11	2138	18.59	
15	LI-SED1-0-15	04/18/11	2214	18.59	
16	LI-SED2-0-15	04/18/11	2250	18.60	
17	LI-SED2-0-15	04/18/11	2327	18.60	
18	LI-SED2-0-15	04/19/11	0003	18.60	
19	LI-SED3-0-15	04/19/11	0039	18.60	
20	LI-SED4-0-15	04/19/11	0115	18.60	
21	LI-SED1-0-15	04/19/11	0151	18.60	
22	LI-SED3-0-15	04/19/11	0227	18.60	
23	ZZZZZ	04/19/11	0304	18.60	
24	INTERNAL	04/19/11	0340	18.60	

QC LIMITS
S1 = 2,4,6-Trichlorophenol (+/- 0.07 MINUTES)

Values outside QC limits.

CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SQ22

Project: LORA LAKE SURFACE SEDIMENT

GC Column: STX CLP2 I.D.: 0.53 (mm)

Instrument ID: ECD1

Init. Calib. Date(s): 04/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE FROM INITIAL CALIBRATION					
SI : 20.93					
CLIENT SAMPLE NO.	LAB FILE ID	DATE ANALYZED	TIME ANALYZED	SI RT	#
01		04/14/11	1927	20.93	
02		04/14/11	2003	20.93	
03		04/14/11	2040	20.93	
04		04/14/11	2116	20.93	
05		04/14/11	2152	20.93	
06		04/14/11	2228	20.93	
07	ZZZZZ	04/14/11	2304	20.93	
08	INTERNAL	04/14/11	1410	20.92	
09	SP34MBS1	04/14/11	1447	20.92	
10	SP34LCS1	04/14/11	1523	20.92	
11	INTERNAL	04/14/11	1901	20.92	
12	ZZZZZ	04/18/11	2025	20.93	
13	ZZZZZ	04/18/11	2102	20.93	
14	INTERNAL	04/18/11	2138	20.93	
15	LL-SED1-0-15	04/18/11	2214	20.93	
16	LL-SED2-0-15	04/18/11	2250	20.94	
17	LL-SED3-0-15	04/18/11	2327	20.94	
18	LL-SED2-0-15	04/18/11	0003	20.94	
19	LL-SED3-0-15	04/18/11	0039	20.94	
20	LL-SED4-0-15	04/18/11	0115	20.94	
21	LL-SED1-0-15	04/18/11	0151	20.94	
22	LL-SED3-0-15	04/18/11	0227	20.94	
23	ZZZZZ	04/18/11	0304	20.94	
24	INTERNAL	04/18/11	0340	20.94	

QC LIMITS

SI = 2,4,6-Trichlorophenol (4 - 0.07 MINUTES)

Values outside QC limits

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SQ22 Project: LORA LAKE SURFACE SEDIMENT
 GC Column: STX CLP1 ID: 0.53 (mm) Instrument ID: ECD1
 Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
 SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION					
S1 : 18.59					
	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT #
	=====	=====	=====	=====	=====
01	LL-SED2-0-15	SQ22B	04/15/11	2014	18.58
02	LL-SED2-0-15	SQ22BMS	04/15/11	2050	18.58
03	LL-SED2-0-15	SQ22BMSD	04/15/11	2126	18.58
04		PCP CCAL	04/16/11	0104	18.59

QC LIMITS

S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SQ22 Project: LORA LAKE SURFACE SEDIMENT
 GC Column: STX CLP2 ID: 0.53 (mm) Instrument ID: ECD1
 Init. Calib. Date(s): 04/14/11 04/14/11

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS,
 SAMPLES, AND STANDARDS IS GIVEN BELOW:

MEAN SURROGATE RT FROM INITIAL CALIBRATION					
S1 : 20.93					
	CLIENT SAMPLE NO.	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	S1 RT #
	=====	=====	=====	=====	=====
01	LL-SED2-0-15	SQ22B	04/15/11	2014	20.92
02	LL-SED2-0-15	SQ22BMS	04/15/11	2050	20.92
03	LL-SED2-0-15	SQ22BMSD	04/15/11	2126	20.92
04		PCP CCAL	04/16/11	0104	20.93

QC LIMITS
 S1 = 2,4,6-Tribromophenol (+/- 0.07 MINUTES)

* Values outside of QC limits.

**Metals Analysis
Report and Summary QC Forms**

ARI Job ID: SP34, SQ22

Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
MC-SED1-0-10-03291	SP34G	11-6956	
MC-SED1-0-10-03291D	SP34GDUP	11-6956	
MC-SED1-0-10-03291S	SP34GSPK	11-6956	
MC-SED2-0-10-03291	SP34H	11-6957	
PBS	SP34MB1	11-6957	
LCSS	SP34MB1SPK	11-6957	
MC-SED3-0-10-03291	SP34I	11-6958	
LL-SED1-0-15-03291	SP34J	11-6959	
LL-SED1-0-15-03291D	SP34JDUP	11-6959	
LL-SED1-0-15-03291S	SP34JSPK	11-6959	
PBW	SP34MB2	11-6959	
LCSW	SP34MB2SPK	11-6959	

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

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

Comments: _____

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:  Name: Jay Kuhn

Date: 4/8/11 Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: MC-SED1-0-10-032911

SAMPLE

Lab Sample ID: SP34G

LIMS ID: 11-6956

Matrix: Sediment

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 79.2%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	6	8	
3050B	04/01/11	6010B	04/07/11	7439-92-1	Lead	2	12	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: MC-SED2-0-10-032911

SAMPLE

Lab Sample ID: SP34H

LIMS ID: 11-6957

Matrix: Sediment

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 74.8%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	6	6	U
3050B	04/01/11	6010B	04/07/11	7439-92-1	Lead	3	11	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: MC-SED3-0-10-032911

SAMPLE

Lab Sample ID: SP34I


QC Report No: SP34-Floyd Snider

LIMS ID: 11-6958

Project: Lora Lake Surface Sediment Sampling

Matrix: Sediment

POS-LL

Data Release Authorized 

Date Sampled: 03/29/11

Reported: 04/08/11

Date Received: 03/30/11

Percent Total Solids: 79.9%


Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	6	6	U
3050B	04/01/11	6010B	04/07/11	7439-92-1	Lead	2	4	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS
 Page 1 of 1

Sample ID: MC-SED1-0-10-032911
MATRIX SPIKE

Lab Sample ID: SP34G
 LIMS ID: 11-6956
 Matrix: Sediment
 Data Release Authorized: 
 Reported: 04/08/11

QC Report No: SP34-Floyd Snider
 Project: Lora Lake Surface Sediment Sampling
 POS-LL
 Date Sampled: 03/29/11
 Date Received: 03/30/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010B	8	235	246	92.3%	
Lead	6010B	12	235	246	90.7%	

Reported in mg/kg-dry

N-Control Limit Not Met
 H-% Recovery Not Applicable, Sample Concentration Too High
 NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: MC-SED1-0-10-032911

DUPLICATE

Lab Sample ID: SP34G

LIMS ID: 11-6956

Matrix: Sediment

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010B	8	7	13.3%	+/- 6	L
Lead	6010B	12	9	28.6%	+/- 2	L*

Reported in mg/kg-dry

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: SP34LCS


QC Report No: SP34-Floyd Snider

LIMS ID: 11-6957

Project: Lora Lake Surface Sediment Sampling

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: NA

Reported: 04/08/11

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010B	190	200	95.0%	
Lead	6010B	185	200	92.5%	

Reported in mg/kg-dry

N-Control limit not met

NA-Not Applicable, Analyte Not Spiked

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Sample ID: METHOD BLANK

Page 1 of 1

Lab Sample ID: SP34MB


QC Report No: SP34-Floyd Snider

LIMS ID: 11-6957

Project: Lora Lake Surface Sediment Sampling

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: NA

Reported: 04/08/11

Date Received: NA

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	5	5	U
3050B	04/01/11	6010B	04/07/11	7439-92-1	Lead	2	2	U

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LL-SED1-0-15-032911-ER
SAMPLE

Lab Sample ID: SP34J

LIMS ID: 11-6959

Matrix: Water

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/L	Q
3010A	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	0.05	0.05	U
3010A	04/01/11	6010B	04/07/11	7439-92-1	Lead	0.02	0.02	U

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED1-0-15-032911-ER
MATRIX SPIKE

Lab Sample ID: SP34J


QC Report No: SP34-Floyd Snider

LIMS ID: 11-6959

Project: Lora Lake Surface Sediment Sampling

Matrix: Water

POS-LL

Data Release Authorized: 

Date Sampled: 03/29/11

Reported: 04/08/11

Date Received: 03/30/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010B	0.05 U	2.06	2.00	103%	
Lead	6010B	0.02 U	2.00	2.00	100%	

Reported in mg/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

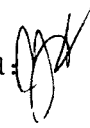
Page 1 of 1

Sample ID: LL-SED1-0-15-032911-ER
DUPLICATE

Lab Sample ID: SP34J

LIMS ID: 11-6959

Matrix: Water

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010B	0.05 U	0.05 U	0.0%	+/- 0.05	L
Lead	6010B	0.02 U	0.02 U	0.0%	+/- 0.02	L

Reported in mg/L

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: SP34LCS

LIMS ID: 11-6959

Matrix: Water

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: NA

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010B	2.02	2.00	101%	
Lead	6010B	1.96	2.00	98.0%	

Reported in mg/L

N-Control limit not met

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: METHOD BLANK

Lab Sample ID: SP34MB

LIMS ID: 11-6959

Matrix: Water

Data Release Authorized: 

Reported: 04/08/11

QC Report No: SP34-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: NA

Date Received: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/L	Q
3010A	04/01/11	6010B	04/07/11	7440-38-2	Arsenic	0.05	0.05	U
3010A	04/01/11	6010B	04/07/11	7439-92-1	Lead	0.02	0.02	U

U-Analyte undetected at given RL

RL-Reporting Limit

Calibration Verification



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34

UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IP040771	2000.0	1925.04	96.3	2000.0	1953.46	97.7	1910.15	95.5	1909.68	95.5	1897.18	94.9	1876.94	93.8
Lead	PB	ICP	IP040771	2000.0	1930.51	96.5	2000.0	1938.85	96.9	1917.58	95.9	1915.58	95.8	1899.83	95.0	1887.03	94.4

Control Limits: Mercury 80-120; Other Metals 90-110

CRDL Standard

CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34



UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IP040771	50.0		50.13	100.3	52.32	104.6								
Lead	PB	ICP	IP040771	20.0		20.14	100.7	20.19	101.0								

Control Limits: no control limits have been established by the EPA at this time.

Calibration Blanks

CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34



UNITS: ug/L

ANALYTE	EL METH	RUN	CRDL	IDL	ICB	CCB1	CCB2	CCB3	CCB4	CCB5
Arsenic	AS ICP	IP040771	10.0	50.0	50.0	50.0	50.0	50.0	50.0	50.0
Lead	PB ICP	IP040771	3.0	20.0	20.0	20.0	20.0	20.0	20.0	20.0

ICP Interference Check Sample



CLIENT: Floyd Snider

ICS SOURCE: I.V.

PROJECT: Lora Lake Surface Se

RUNID: IP040771

SDG: SP34

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	198664.3	198922.0	99.5	196129.7	196390.6	98.2			
Antimony	1000	1000	3.2	1011.1	101.1	4.2	990.6	99.1			
Arsenic	1000	1000	15.9	988.4	98.8	17.3	969.9	97.0			
Barium	1000	1000	-0.5	975.5	97.6	-0.5	959.6	96.0			
Beryllium	1000	1000	0.0	969.0	96.9	-0.1	951.0	95.1			
Boron			-12.2	-9.2		-8.4	-12.2				
Cadmium	1000	1000	1.3	988.3	98.8	1.2	978.5	97.9			
Calcium	100000	100000	98936.4	98630.9	98.6	97565.3	97876.7	97.9			
Chromium	1000	1000	0.8	996.9	99.7	-0.9	991.0	99.1			
Cobalt	1000	1000	1.0	949.3	94.9	1.0	933.3	93.3			
Copper	1000	1000	-0.7	1024.4	102.4	-0.7	1019.4	101.9			
Iron	200000	200000	195500.2	195209.8	97.6	191329.5	191635.2	95.8			
Lead	1000	1000	-4.1	928.9	92.9	-3.3	908.5	90.9			
Magnesium	100000	100000	102960.4	98140.2	98.1	100895.9	97209.1	97.2			
Manganese	1000	1000	0.3	932.1	93.2	0.4	911.1	91.1			
Molybdenum			4.5	3.7		3.5	3.1				
Nickel	1000	1000	0.3	962.8	96.3	2.2	946.7	94.7			
Potassium			44.2	585.7		50.4	578.6				
Selenium	1000	1000	20.6	994.4	99.4	22.3	960.8	96.1			
Silicon			-5.4	-13.0		-8.7	-10.8				
Silver	1000	1000	0.1	1013.3	101.3	0.5	996.6	99.7			
Sodium			-1.9	-4.8		-12.2	-14.4				
Strontium			3.7	3.7		3.6	3.6				
Thallium	1000	1000	4.7	920.3	92.0	5.1	898.3	89.8			
Tin			1.6	1.8		0.8	1.2				
Titanium			-3.7	-2.9		-3.9	-3.1				
Vanadium	1000	1000	-0.3	960.1	96.0	-0.2	955.4	95.5			
Zinc	1000	1000	2.3	956.0	95.6	1.5	953.1	95.3			

IDLs and ICP Linear Ranges



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2011	30000.0	2/3/2011
Lead	PB	ICP	OPTIMA ICP 2	220.35		3	20.0	4/1/2011	300000.0	2/3/2011

ICP Interelement Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34

IEC DATE: 3/14/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	9.9066900	0.000000	0.000000
Arsenic	188.98	0.000000	0.000000	0.000000	0.000000	0.0893242	0.000000	-1.0280600	0.9896930	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.1423420	0.000000	0.000000	0.0649797
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Cadmium	228.80	0.000000	3.6086500	0.000000	0.000000	0.000000	0.000000	0.1351930	0.000000	0.000000	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.000000	0.000000	0.000000	0.000000	0.0148832	0.000000	0.000000	0.000000	0.000000	-0.0451581
Cobalt	228.62	0.000000	0.000000	0.0227510	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Copper	324.75	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.1969920	-0.0283867	0.000000	-0.0499474
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Lead	220.35	-0.1952990	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-1.9460200	1.1789000	0.0588763
Magnesium	279.08	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-1.3945700	-0.8349460	0.000000	0.4579600
Manganese	257.61	0.0056707	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-0.0082982
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.0185703	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.000000	0.000000	0.000000	0.000000	0.1513640	0.000000	0.000000	0.000000	0.000000	0.000000
Silicon	288.16	0.000000	0.000000	0.000000	0.000000	0.000000	-3.7058000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Thallium	190.80	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	2.4468400	0.3572340	0.000000	-0.1350510
Tin	189.93	0.000000	0.000000	0.000000	0.000000	-0.2650000	0.000000	0.000000	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	0.000000	0.000000	0.1735400	0.000000	0.000000	0.1546720	0.000000	0.000000
Vanadium	292.40	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	-4.7348500	0.000000	0.0820500
Zinc	206.20	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0805698	0.000000	0.000000

ICP Inter-element Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SP34

IEC DATE: 3/14/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.000000	1.878880	12.130700	0.000000	0.000000	0.000000	2.350330	0.000000	18.036400	0.000000
Antimony	206.84	0.000000	0.000000	0.000000	-0.385518	0.000000	0.000000	-1.514710	0.000000	-3.206310	0.000000
Arsenic	188.98	0.000000	0.000000	1.362000	0.000000	0.000000	0.000000	-8.144460	0.000000	0.000000	0.000000
Barium	233.53	0.000000	0.000000	0.000000	0.075903	0.000000	0.000000	0.000000	0.000000	0.490078	0.000000
Beryllium	313.04	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.537185	0.000000
Cadmium	228.80	0.000000	0.000000	0.000000	-0.621208	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Calcium	317.93	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Chromium	267.72	0.0685552	0.000000	0.219900	0.000000	0.000000	0.000000	0.000000	0.000000	0.268954	0.000000
Cobalt	228.62	0.000000	0.000000	-0.253629	0.158455	0.000000	0.000000	1.615940	0.000000	0.000000	0.000000
Copper	324.75	0.0040015	0.000000	0.155843	0.000000	0.000000	0.000000	0.303169	0.000000	0.000000	0.000000
Iron	273.96	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	5.275560	0.000000
Lead	220.35	0.000000	0.000000	-0.372923	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Magnesium	279.08	0.000000	0.000000	-2.695300	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Manganese	257.61	0.0058832	0.000000	0.000000	0.000000	-0.265904	0.000000	0.000000	0.000000	-0.0245885	0.000000
Molybdenum	202.03	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Nickel	231.60	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Potassium	766.49	0.000000	0.000000	0.000000	0.000000	0.000000	-0.589798	0.000000	0.000000	0.000000	0.000000
Selenium	196.03	0.0859101	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silicon	288.16	-0.119790	0.000000	-1.847410	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Silver	328.07	0.000000	0.177307	0.106727	0.000000	0.000000	0.000000	0.000000	0.000000	-0.222475	0.000000
Sodium	589.59	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	37.740000	0.000000	0.000000	62.929000
Thallium	190.80	0.000000	0.000000	-3.913880	0.000000	0.000000	0.000000	0.000000	0.000000	1.622560	0.000000
Tin	189.93	0.000000	0.000000	0.000000	0.000000	0.000000	-0.636504	-0.351611	0.000000	0.000000	0.000000
Titanium	334.90	0.000000	0.000000	1.236370	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000
Vanadium	252.40	0.000000	-0.161968	-0.951633	0.000000	0.000000	0.000000	0.620937	0.000000	0.000000	0.000000
Zinc	206.20	0.000000	0.000000	0.254730	0.000000	-0.0673589	0.000000	0.000000	0.000000	0.000000	0.000000

Preparation Log



CLIENT: Floyd Snider

ANALYSIS METHOD: ICP

PROJECT: Lora Lake Surface Se

ARI PREP CODE: SWC

SDG: SP34

PREPDATE: 4/1/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
MC-SED1-0-10-03291	SP34G	1.028	0.0	50.0
MC-SED1-0-10-03291D	SP34GDUP	1.027	0.0	50.0
MC-SED1-0-10-03291S	SP34GSPK	1.025	0.0	50.0
MC-SED2-0-10-03291	SP34H	1.061	0.0	50.0
MC-SED3-0-10-03291	SP34I	1.043	0.0	50.0
PBS	SP34MB1	1.000	0.0	50.0
LCSS	SP34MB1SPK	1.000	0.0	50.0

Preparation Log



CLIENT: Floyd Snider

ANALYSIS METHOD: ICP

PROJECT: Lora Lake Surface Se

ARI PREP CODE: TWC

SDG: SP34

PREPDATE: 4/1/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
LL-SED1-0-15-03291	SP34J	0.000	50.0	50.0
LL-SED1-0-15-03291D	SP34JDUP	0.000	50.0	50.0
LL-SED1-0-15-03291S	SP34JSPK	0.000	50.0	50.0
PBW	SP34MB2	0.000	50.0	50.0
LCSW	SP34MB2SPK	0.000	50.0	50.0



Analysis Run Log

CLIENT: Floyd Snider
 PROJECT: Lora Lake Surface Se
 SDG: SP34
 INSTRUMENT ID: OPTIMA ICP 2
 RUNID: IP040771
 METHOD: ICP
 START DATE: 4/7/2011
 END DATE: 4/7/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN		
S0		1.00	11151		X																												X		
S2		1.00	11191																															X	
S3		1.00	11205		X																													X	
S4		1.00	11232																																
S5		1.00	11253																																
ICV		1.00	11375		X																													X	
ICB		1.00	11405		X																													X	
CRI		1.00	11450		X																													X	
ICSA		1.00	11491		X																													X	
ICSAB		1.00	11532		X																													X	
CCV		1.00	11570		X																													X	
CCB		1.00	12002		X																													X	
ZZZZZZ	SP69MB1		2.00	12043																															
ZZZZZZ	SP69A-L		10.00	12084																															
ZZZZZZ	SP69A		2.00	12123																															
ZZZZZZ	SP69ADUP		2.00	12165																															
ZZZZZZ	SP69ASPK		2.00	12210																															
ZZZZZZ	ZZZZZZ		2.00	12244																															
ZZZZZZ	SP69B		2.00	12283																															
ZZZZZZ	SP69C		2.00	12324																															
ZZZZZZ	SP69D		2.00	12365																															
ZZZZZZ	SP69MB1SPK		2.00	12410																															
CCV	CCV2		1.00	12445					X																									X	
CCB	CCB2		1.00	12475					X																									X	
CRI	CRIF		1.00	12520					X																									X	
ICSA	ICSAF		1.00	12562					X																									X	
ICSAB	ICSABF		1.00	13003					X																									X	
CCV	CCV3		1.00	13041					X																									X	
CCB	CCB3		1.00	13071					X																									X	
PBS	SP34MB1		2.00	13112					X																									X	
PBW	SP34MB2		1.00	13153					X																									X	
LL-SED1-0-15-03291D	SP34JDUP		1.00	13194					X																									X	
LL-SED1-0-15-03291	SP34J		1.00	13234					X																									X	
LL-SED1-0-15-03291S	SP34JSPK		1.00	13275					X																									X	
MC-SED2-0-10-03291	SP34H		2.00	13314					X																									X	



Analysis Run Log

CLIENT: Floyd Snider
 PROJECT: Lora Lake Surface Se INSTRUMENT ID: OPTIMA ICP 2 START DATE: 4/7/2011
 SDG: SP34 RUNID: IP040771 METHOD: ICP END DATE: 4/7/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN			
MC-SED1-0-10-03291D	SP34GDUP	2.00	13354																															X		
MC-SED1-0-10-03291	SP34G	2.00	13393																															X		
MC-SED1-0-10-03291S	SP34GSPK	2.00	13433																															X		
LCSW	SP34MB2SPK	1.00	13463																															X		
CCV	CCV4	1.00	13502																															X		
CCB	CCB4	1.00	13532																															X		
ZZZZZZ	SP50MB	1.00	13573																																	
ZZZZZZ	SP50A	1.00	14014																																	
ZZZZZZ	SP50B	1.00	14055																																	
ZZZZZZ	SP50C	1.00	14095																																	
ZZZZZZ	SP50D	1.00	14134																																	
ZZZZZZ	SP50E	1.00	14174																																	
ZZZZZZ	SP50F	1.00	14213																																	
MC-SED3-0-10-03291	SP34I	2.00	14253																																X	
LCSS	SP34MB1SPK	2.00	14292																																X	
ZZZZZZ	SP50MBSPK	1.00	14332																																	
CCV	CCV5	1.00	14371																																	X
CCB	CCB5	1.00	14401																																	X

Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
LL-SED1-0-15-03291	SQ22A	11-7355	
PBS	SQ22MB1	11-7355	
LCSS	SQ22MB1SPK	11-7355	
LL-SED2-0-15-03291	SQ22B	11-7356	
LL-SED2-0-15-03291D	SQ22BDUP	11-7356	
LL-SED2-0-15-03291S	SQ22BSPK	11-7356	
LL-SED3-0-15-03291	SQ22C	11-7357	
LL-SED4-0-15-03291	SQ22D	11-7358	
LL-SED1-0-15-03291	SQ22E	11-7359	
LL-SED5-0-15-03291	SQ22F	11-7360	

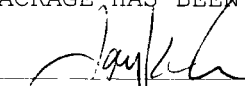
Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

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

Comments: _____

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature: 

Name: Jay Kuhn

Date: 4/12/11

Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: LL-SED1-0-15-032911

SAMPLE

Lab Sample ID: SQ22A

LIMS ID: 11-7355

Matrix: Sediment

Data Release Authorized: 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 31.1%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	20	20	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	6	319	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LL-SED2-0-15-032911
SAMPLE

Lab Sample ID: SQ22B

LIMS ID: 11-7356

Matrix: Sediment

Data Release Authorized: 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 19.6%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	20	50	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	10	390	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED3-0-15-032911

SAMPLE

Lab Sample ID: SQ22C

LIMS ID: 11-7357

Matrix: Sediment

Data Release Authorized: *[Signature]*

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 27.0%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	20	70	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	7	361	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED4-0-15-032911

SAMPLE

Lab Sample ID: SQ22D


QC Report No: SQ22-Floyd Snider

LIMS ID: 11-7358

Project: Lora Lake Surface Sediment Sampling

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/29/11

Reported: 04/12/11

Date Received: 03/30/11

Percent Total Solids: 24.6%


Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	20	40	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	8	492	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS
Page 1 of 1

Sample ID: LL-SED1-0-15-032911-D
SAMPLE

Lab Sample ID: SQ22E
LIMS ID: 11-7359
Matrix: Sediment
Data Release Authorized: 
Reported: 04/12/11

QC Report No: SQ22-Floyd Snider
Project: Lora Lake Surface Sediment Sampling
POS-LL
Date Sampled: 03/29/11
Date Received: 03/30/11

Percent Total Solids: 31.2%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	20	20	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	6	281	

U-Analyte undetected at given RL
RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: LL-SED5-0-15-032911

SAMPLE

Lab Sample ID: SQ22F

LIMS ID: 11-7360

Matrix: Sediment

Data Release Authorized 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

Percent Total Solids: 80.0%

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	6	7	
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	2	48	

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1


Sample ID: LL-SED2-0-15-032911

MATRIX SPIKE

Lab Sample ID: SQ22B

LIMS ID: 11-7356

Matrix: Sediment

Data Release Authorized: 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Arsenic	6010B	50	1,030	982	99.8%	
Lead	6010B	390	1,370	982	99.8%	

Reported in mg/kg-dry

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LL-SED2-0-15-032911
DUPLICATE

Lab Sample ID: SQ22B

LIMS ID: 11-7356

Matrix: Sediment

Data Release Authorized: 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: 03/29/11

Date Received: 03/30/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Arsenic	6010B	50	40	22.2%	+/- 20	L
Lead	6010B	390	370	5.3%	+/- 20%	

Reported in mg/kg-dry

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: SQ22LCS

LIMS ID: 11-7355

Matrix: Sediment

Data Release Authorized: 

Reported: 04/12/11

QC Report No: SQ22-Floyd Snider

Project: Lora Lake Surface Sediment Sampling

POS-LL

Date Sampled: NA

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Arsenic	6010B	165	200	82.5%	
Lead	6010B	161	200	80.5%	

Reported in mg/kg-dry

N-Control limit not met

NA-Not Applicable, Analyte Not Spiked

Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Sample ID: METHOD BLANK

Page 1 of 1

Lab Sample ID: SQ22MB


QC Report No: SQ22-Floyd Snider

LIMS ID: 11-7355

Project: Lora Lake Surface Sediment Sampling

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: NA

Reported: 04/12/11

Date Received: NA

Percent Total Solids: NA

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	RL	mg/kg-dry	Q
3050B	04/06/11	6010B	04/11/11	7440-38-2	Arsenic	5	5	U
3050B	04/06/11	6010B	04/11/11	7439-92-1	Lead	2	2	U

U-Analyte undetected at given RL
RL-Reporting Limit



Calibration Verification

CLIENT: Floyd Snider
 PROJECT: Lora Lake Surface Se
 SDG: SQ22

UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	ICP	IP041171	2000.0	1965.34	98.3	2000.0	1955.49	97.8	1940.91	97.0	1925.89	96.3	1951.19	97.6		
Lead	PB	ICP	IP041171	2000.0	1976.21	98.8	2000.0	1959.94	98.0	1945.07	97.3	1930.41	96.5	1957.40	97.9		

Control Limits: Mercury 80-120; Other Metals 90-110

CRDI Standard

CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22



UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Arsenic	AS	ICP	IP041171	50.0		48.42	96.8	48.25	96.5								
Lead	PB	ICP	IP041171	20.0		20.39	102.0	19.66	98.3								

Control Limits: no control limits have been established by the EPA at this time.

Calibration Blanks

CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22



UNITS: ug/L

ANALYTE	EL METH	RUN	CRDL	IDL	ICB	CCB1	CCB2	CCB3	CCB4	CCB5
				C	C	C	C	C	C	C
Arsenic	AS ICP	IP041171	10.0	50.0	50.0	50.0	50.0	50.0	50.0	50.0
Lead	PB ICP	IP041171	3.0	20.0	20.0	20.0	20.0	20.0	20.0	20.0

ICP Interference Check Sample



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22

ICS SOURCE: I.V.

RUNID: IP041171

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Aluminum	200000	200000	198769.6	199716.4	99.9	194155.7	196828.1	98.4			
Antimony	1000	1000	4.7	1044.4	104.4	1.7	1020.3	102.0			
Arsenic	1000	1000	12.4	1012.7	101.3	12.5	991.8	99.2			
Barium	1000	1000	0.2	1006.0	100.6	0.8	982.0	98.2			
Beryllium	1000	1000	0.0	997.1	99.7	0.0	970.0	97.0			
Boron			-6.3	-9.0		-10.3	-10.7				
Cadmium	1000	1000	1.3	1013.4	101.3	1.2	995.3	99.5			
Calcium	100000	100000	100429.9	100857.2	100.9	98712.6	99904.6	99.9			
Chromium	1000	1000	-1.5	1026.6	102.7	-1.0	1014.0	101.4			
Cobalt	1000	1000	0.9	975.4	97.5	1.3	938.5	93.9			
Copper	1000	1000	0.5	1046.5	104.7	0.6	1036.8	103.7			
Iron	200000	200000	197988.3	200347.0	100.2	189498.7	193000.8	96.5			
Lead	1000	1000	-7.1	948.2	94.8	-5.8	926.9	92.7			
Magnesium	100000	100000	103448.7	100137.4	100.1	102443.1	98914.3	98.9			
Manganese	1000	1000	0.1	954.6	95.5	0.3	920.5	92.1			
Molybdenum			3.8	3.5		3.0	3.5				
Nickel	1000	1000	3.1	982.8	98.3	3.5	949.5	95.0			
Potassium			43.1	547.8		68.1	552.1				
Selenium	1000	1000	19.9	1021.2	102.1	32.4	995.7	99.6			
Silicon			-12.3	-13.4		-15.2	-10.5				
Silver	1000	1000	0.3	1025.0	102.5	0.5	988.1	98.8			
Sodium			11.7	8.8		13.6	18.8				
Strontium			3.8	3.7		3.7	3.7				
Thallium	1000	1000	6.7	955.8	95.6	7.4	919.4	91.9			
Tin			-2.0	-0.6		-3.6	-3.0				
Titanium			-3.4	-2.5		-4.7	-2.9				
Vanadium	1000	1000	-4.0	973.3	97.3	-4.3	965.7	96.6			
Zinc	1000	1000	3.6	982.6	98.3	3.3	973.3	97.3			

IDLs and ICP Linear Ranges



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Arsenic	AS	ICP	OPTIMA ICP 2	197.20		10	50.0	4/1/2011	30000.0	8/19/2010
Lead	PB	ICP	OPTIMA ICP 2	220.35		3	20.0	4/1/2011	300000.0	8/19/2010

ICP Interelement Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22

IEC DATE: 3/14/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AL	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Antimony	206.84	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	9.9066900	0.0000000	0.0000000
Arsenic	188.98	0.0000000	0.0000000	0.0000000	0.0000000	0.0893242	0.0000000	-1.0280600	0.9896930	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.1423420	0.0000000	0.0000000	0.0649797
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Cadmium	228.80	0.0000000	3.6086500	0.0000000	0.0000000	0.0000000	0.0000000	0.1351930	0.0000000	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Chromium	267.72	0.0000000	0.0000000	0.0000000	0.0000000	0.0148832	0.0000000	0.0000000	0.0000000	0.0000000	-0.0451581
Cobalt	228.62	0.0000000	0.0000000	0.0227510	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Copper	324.75	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.1969920	-0.0283867	0.0000000	-0.0499474
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Lead	220.35	-0.1952990	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Magnesium	279.08	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.9460200	1.1789000	0.0588763
Manganese	257.61	0.0056707	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.3945700	-0.8349460	0.0000000	0.4579600
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0185703	0.0000000	0.0000000	0.0000000	0.0000000	-0.0082982
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0000000	0.0000000	0.0000000	0.0000000	0.1513640	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silicon	288.16	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.7058000	0.0000000	0.0000000	0.0000000	0.0000000
Silver	328.07	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	-0.2650000	0.0000000	0.0000000	0.0000000	0.0000000	-0.1350510
Titanium	334.90	0.0000000	0.0000000	0.0000000	0.0000000	0.1735400	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Vanadium	292.40	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.1546720	0.0000000	0.0000000
Zinc	206.20	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-4.7348500	0.0000000	0.0820500
									0.0805698	0.0000000	0.0000000

ICP Interelement Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake Surface Se

SDG: SQ22

IEC DATE: 3/14/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PB	SB	TI	TL	V	ZN
Aluminum	308.22	0.0000000	1.8788800	12.1307000	0.0000000	0.0000000	0.0000000	2.3503300	0.0000000	18.0364000	0.0000000
Antimony	206.84	0.0000000	0.0000000	0.0000000	-0.3855180	0.0000000	0.0000000	-1.5147100	0.0000000	-3.2063100	0.0000000
Arsenic	188.98	0.0000000	0.0000000	1.3620000	0.0000000	0.0000000	0.0000000	-8.1444600	0.0000000	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0759003	0.0000000	0.0000000	0.0000000	0.0000000	0.4900780	0.0000000
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.5371850	0.0000000
Cadmium	228.80	0.0000000	0.0000000	0.0000000	-0.6212080	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Chromium	267.72	0.0685552	0.0000000	0.2199000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.2689540	0.0000000
Cobalt	228.62	0.0000000	0.0000000	-0.2536290	0.1584550	0.0000000	0.0000000	1.6159400	0.0000000	0.0000000	0.0000000
Copper	324.75	0.0040015	0.0000000	0.1558430	0.0000000	0.0000000	0.0000000	0.3031690	0.0000000	0.0000000	0.0000000
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	5.2755600	0.0000000
Lead	220.35	0.0000000	0.0000000	-0.3729230	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Magnesium	279.08	0.0000000	0.0000000	-2.6953000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Manganese	257.61	0.0058832	0.0000000	0.0000000	0.0000000	-0.2659040	0.0000000	0.0000000	0.0000000	-0.0245885	0.0000000
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.5897980	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0859101	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silicon	288.16	-0.1197900	0.0000000	-1.8474100	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Silver	328.07	0.0000000	0.1773070	0.1067270	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.2224750	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.0000000	0.0000000	-3.9138800	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	1.6225600	0.0000000
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.6365040	-0.3516110	0.0000000	0.0000000	0.0000000
Titanium	334.90	0.0000000	0.0000000	1.2363700	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Vanadium	292.40	0.0000000	-0.1619680	-0.9516330	0.0000000	0.0000000	0.0000000	0.6209970	0.0000000	0.0000000	0.0000000
Zinc	206.20	0.0000000	0.0000000	0.2547300	0.0000000	-0.0673589	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000

Preparation Log



CLIENT: Floyd Snider

ANALYSIS METHOD: ICP

PROJECT: Lora Lake Surface Se

ARI PREP CODE: SWC

SDG: SQ22

PREPDATE: 4/6/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
LL-SED1-0-15-03291	SQ22A	1.056	0.0	50.0
LL-SED2-0-15-03291	SQ22B	1.041	0.0	50.0
LL-SED2-0-15-03291D	SQ22BDUP	1.045	0.0	50.0
LL-SED2-0-15-03291S	SQ22BSPK	1.039	0.0	50.0
LL-SED3-0-15-03291	SQ22C	1.066	0.0	50.0
LL-SED4-0-15-03291	SQ22D	1.078	0.0	50.0
LL-SED1-0-15-03291	SQ22E	1.024	0.0	50.0
LL-SED5-0-15-03291	SQ22F	1.059	0.0	50.0
PBS	SQ22MB1	1.000	0.0	50.0
LCSS	SQ22MB1SPK	1.000	0.0	50.0