

USEPA - ITD

FORM 4A
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory Episode No.:

Contract No.: SAS No.:

Initial Calibration Date:

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 07APR11M Sam:15

Analysis Date: 8-APR-11 04:08:18

| 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 | 0.81 | 0.65-0.89 | y | 12.3 | 7.80 - 12.9 |
| 1,2,3,7,8-PeCDD | M+2/M+4 | 1.52 | 1.32-1.78 | y | 50.7 | 39.0 - 65.0 |
| 1,2,3,4,7,8-HxCDD | M+2/M+4 | 1.27 | 1.05-1.43 | y | 47.2 | 39.0 - 64.0 |
| 1,2,3,6,7,8-HxCDD | M+2/M+4 | 1.30 | 1.05-1.43 | y | 48.6 | 39.0 - 64.0 |
| 1,2,3,7,8,9-HxCDD | M+2/M+4 | 1.26 | 1.05-1.43 | y | 51.7 | 41.0 - 61.0 |
| 1,2,3,4,6,7,8-HpCDD | M+2/M+4 | 0.90 | 0.88-1.20 | y | 51.7 | 43.0 - 58.0 |
| OCDD | M+2/M+4 | 0.92 | 0.76-1.02 | y | 95.7 | 79.0 - 126 |
| 2,3,7,8-TCDF | M/M+2 | 0.72 | 0.65-0.89 | y | 9.82 | 8.40 - 12.0 |
| 1,2,3,7,8-PeCDF | M+2/M+4 | 1.61 | 1.32-1.78 | y | 50.5 | 41.0 - 60.0 |
| 2,3,4,7,8-PeCDF | M+2/M+4 | 1.57 | 1.32-1.78 | y | 47.9 | 41.0 - 60.0 |
| 1,2,3,4,7,8-HxCDF | M+2/M+4 | 1.22 | 1.05-1.43 | y | 49.7 | 45.0 - 56.0 |
| 1,2,3,6,7,8-HxCDF | M+2/M+4 | 1.25 | 1.05-1.43 | y | 51.0 | 44.0 - 57.0 |
| 2,3,4,6,7,8-HxCDF | M+2/M+4 | 1.25 | 1.05-1.43 | y | 51.1 | 44.0 - 57.0 |
| 1,2,3,7,8,9-HxCDF | M+2/M+4 | 1.26 | 1.05-1.43 | y | 49.8 | 45.0 - 56.0 |
| 1,2,3,4,6,7,8-HpCDF | M+2/M+4 | 1.09 | 0.88-1.20 | y | 49.4 | 45.0 - 55.0 |
| 1,2,3,4,7,8,9-HpCDF | M+2/M+4 | 1.03 | 0.88-1.20 | y | 49.4 | 43.0 - 58.0 |
| OCDF | M+2/M+4 | 0.94 | 0.76-1.02 | y | 99.9 | 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: 4/8/11

USEPA - ITD

FORM 4B
PCDD/PCDF CALIBRATION VERIFICATION

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Initial Calibration Date:

Instrument ID: FAL3

GC Column ID: DB5

VER Data Filename: 07APR11M

Sam:15

Analysis Date: 8-APR-11 04:08:18


| 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 | 0.71 | 0.65-0.89 | y | 97.8 | 82.0 - 121 |
| 13C-1,2,3,7,8-PeCDD | M+2/M+4 | 1.78 | 1.32-1.78 | y | 102 | 62.0 - 160 |
| 13C-1,2,3,4,7,8-HxCDD | M+2/M+4 | 1.27 | 1.05-1.43 | y | 86.4 | 85.0 - 117 |
| 13C-1,2,3,6,7,8-HxCDD | M+2/M+4 | 1.32 | 1.05-1.43 | y | 103 | 85.0 - 118 |
| 13C-1,2,3,4,6,7,8-HpCDD | M+2/M+4 | 1.01 | 0.88-1.20 | y | 105 | 72.0 - 138 |
| 13C-OCDD | M+2/M+4 | 0.96 | 0.76-1.02 | y | 198 | 96.0 - 415 |
| 13C-2,3,7,8-TCDF | M/M+2 | 0.88 | 0.65-0.89 | y | 97.3 | 71.0 - 140 |
| 13C-1,2,3,7,8-PeCDF | M+2/M+4 | 1.63 | 1.32-1.78 | y | 106 | 76.0 - 130 |
| 13C-2,3,4,7,8-PeCDF | M+2/M+4 | 1.65 | 1.32-1.78 | y | 108 | 77.0 - 130 |
| 13C-1,2,3,4,7,8-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 87.5 | 76.0 - 131 |
| 13C-1,2,3,6,7,8-HxCDF | M/M+2 | 0.48 | 0.43-0.59 | y | 93.5 | 70.0 - 143 |
| 13C-2,3,4,6,7,8-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 92.9 | 73.0 - 137 |
| 13C-1,2,3,7,8,9-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 94.5 | 74.0 - 135 |
| 13C-1,2,3,4,6,7,8-HpCDF | M/M+2 | 0.49 | 0.37-0.51 | y | 98.1 | 78.0 - 129 |
| 13C-1,2,3,4,7,8,9-HpCDF | M/M+2 | 0.48 | 0.37-0.51 | y | 104 | 77.0 - 129 |
| 13C-OCDF | M+2/M+4 | 0.96 | 0.76-1.02 | y | 182 | 96.0 - 415 |
| CLEANUP STANDARD (4) | | | | | | |
| 37Cl-2,3,7,8-TCDD | | | | | 9.24 | 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: 4/8/11

USEPA - ITD

FORM 6A

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date:

Instrument ID: FAL3

GC Column ID: DB5

Analysis Date: 8-APR-11 04:08:18

CS3 or VER Data Filename: 07APR11M

Sam:15

| 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 | | 1.023 | 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: 4/8/11

USEPA - ITD

FORM 6B

PCDD/PCDF RELATIVE RETENTION TIMES

Lab Name: Frontier Analytical Laboratory

Episode No.:

Contract No.:

SAS No.:

Init. Cal. Date:

Instrument ID: FAL3

GC Column ID: DB5


Analysis Date: 8-APR-11 04:08:18

CS3 or VER Data Filename: 07APR11M


Sam:15

| 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 | 1.000 | 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 | | 1.127 | 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 | | 1.151 | 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: Date: 

| Name | Resp | RA | RT | RRF | Conc | Qual | Fac Noise-1 | Noise-2 | DL | Rec | #Hom |
|--------------------------|----------|--------|-------|------|------|------|-------------|---------|----|-------|------|
| 2,3,7,8-TCDD | 3.52e+06 | 0.81 y | 27:38 | 1.13 | 12.3 | | 2.50 | - | * | | |
| 1,2,3,7,8-PeCDD | 1.34e+07 | 1.52 y | 33:29 | 1.02 | 50.7 | | 2.50 | - | * | | |
| 1,2,3,4,7,8-HxCDD | 1.45e+07 | 1.27 y | 38:52 | 1.45 | 47.2 | | 2.50 | - | * | | |
| 1,2,3,6,7,8-HxCDD | 1.40e+07 | 1.30 y | 39:02 | 1.45 | 48.6 | | 2.50 | - | * | | |
| 1,2,3,7,8,9-HxCDD | 1.56e+07 | 1.26 y | 39:28 | 1.47 | 51.7 | | 2.50 | - | * | | |
| 1,2,3,4,6,7,8-HpCDD | 1.21e+07 | 0.90 y | 44:29 | 1.30 | 51.7 | | 2.50 | - | * | | |
| OCDD | 1.72e+07 | 0.92 y | 50:07 | 1.45 | 95.7 | | 2.50 | - | * | | |
| 2,3,7,8-TCDF | 4.88e+06 | 0.72 y | 26:51 | 1.15 | 9.82 | | 2.50 | - | * | | |
| 1,2,3,7,8-PeCDF | 1.79e+07 | 1.61 y | 31:45 | 0.89 | 50.5 | | 2.50 | - | * | | |
| 2,3,4,7,8-PeCDF | 1.69e+07 | 1.57 y | 33:04 | 0.89 | 47.9 | | 2.50 | - | * | | |
| 1,2,3,4,7,8-HxCDF | 1.67e+07 | 1.22 y | 37:28 | 1.01 | 49.7 | | 2.50 | - | * | | |
| 1,2,3,6,7,8-HxCDF | 2.01e+07 | 1.25 y | 37:41 | 0.89 | 51.0 | | 2.50 | - | * | | |
| 2,3,4,6,7,8-HxCDF | 1.86e+07 | 1.25 y | 38:36 | 1.02 | 51.1 | | 2.50 | - | * | | |
| 1,2,3,7,8,9-HxCDF | 2.12e+07 | 1.26 y | 40:04 | 1.10 | 49.8 | | 2.50 | - | * | | |
| 1,2,3,4,6,7,8-HpCDF | 1.46e+07 | 1.09 y | 42:35 | 1.48 | 49.4 | | 2.50 | - | * | | |
| 1,2,3,4,7,8,9-HpCDF | 1.16e+07 | 1.03 y | 45:25 | 1.43 | 49.4 | | 2.50 | - | * | | |
| OCDF | 1.84e+07 | 0.94 y | 50:31 | 0.84 | 99.9 | | 2.50 | - | * | | |
| 13C-2,3,7,8-TCDD | 2.53e+07 | 0.71 y | 27:37 | 1.03 | 97.8 | | | | | 97.8 | |
| 13C-1,2,3,7,8-PeCDD | 2.60e+07 | 1.78 y | 33:28 | 1.01 | 102 | | | | | 102 | |
| 13C-1,2,3,4,7,8-HxCDD | 2.12e+07 | 1.27 y | 38:50 | 1.19 | 86.4 | | | | | 86.4 | |
| 13C-1,2,3,6,7,8-HxCDD | 1.98e+07 | 1.32 y | 39:00 | 0.94 | 103 | | | | | 103 | |
| 13C-1,2,3,4,6,7,8-HpCDD | 1.80e+07 | 1.01 y | 44:28 | 0.83 | 105 | | | | | 105 | |
| 13C-OCDD | 2.48e+07 | 0.96 y | 50:06 | 0.61 | 198 | | | | | 98.8 | |
| 13C-2,3,7,8-TCDF | 4.33e+07 | 0.88 y | 26:50 | 0.98 | 97.3 | | | | | 97.3 | |
| 13C-1,2,3,7,8-PeCDF | 4.00e+07 | 1.63 y | 31:43 | 0.83 | 106 | | | | | 106 | |
| 13C-2,3,4,7,8-PeCDF | 3.94e+07 | 1.65 y | 33:03 | 0.80 | 108 | | | | | 108 | |
| 13C-1,2,3,4,7,8-HxCDF | 3.33e+07 | 0.47 y | 37:27 | 1.84 | 87.5 | | | | | 87.5 | |
| 13C-1,2,3,6,7,8-HxCDF | 4.42e+07 | 0.48 y | 37:39 | 2.29 | 93.5 | | | | | 93.5 | |
| 13C-2,3,4,6,7,8-HxCDF | 3.57e+07 | 0.47 y | 38:36 | 1.86 | 92.9 | | | | | 92.9 | |
| 13C-1,2,3,7,8,9-HxCDF | 3.87e+07 | 0.47 y | 40:02 | 1.98 | 94.5 | | | | | 94.5 | |
| 13C-1,2,3,4,6,7,8-HpCDF | 2.00e+07 | 0.49 y | 42:34 | 0.99 | 98.1 | | | | | 98.1 | |
| 13C-1,2,3,4,7,8,9-HpCDF | 1.64e+07 | 0.48 y | 45:24 | 0.77 | 104 | | | | | 104 | |
| 13C-OCDF | 4.37e+07 | 0.96 y | 50:29 | 1.17 | 182 | | | | | 90.9 | |
| 37Cl-2,3,7,8-TCDD | 1.70e+06 | | 27:37 | 0.73 | 9.24 | | | | | 92.4 | |
| 13C-1,2,3,4-TCDD | 2.52e+07 | 0.70 y | 27:00 | - | 66.3 | | | | | | |
| 13C-1,2,3,4-TCDF | 4.54e+07 | 0.87 y | 25:44 | - | 63.1 | | | | | | |
| 13C-1,2,3,7,8,9-HxCDD | 2.07e+07 | 1.28 y | 39:27 | - | 83.2 | | | | | | |
| Total Tetra-Dioxins | 1.64e+07 | | 23:12 | 1.13 | 57.2 | | 2.50 | - | * | | 21 |
| Total Penta-Dioxins | 2.90e+07 | | 30:29 | 1.02 | 110 | | 2.50 | - | * | | 6 |
| Total Hexa-Dioxins | 5.00e+07 | | 36:24 | 1.46 | 167 | | 2.50 | - | * | | 10 |
| Total Hepta-Dioxins | 2.57e+07 | | 43:06 | 1.30 | 110 | | 2.50 | - | * | | 20 |
| Total Tetra-Furans | 2.35e+07 | | 23:13 | 1.15 | 47.3 | | 2.50 | - | * | | 22 |
| 1st Fn. Tot Penta-Furans | 1.89e+07 | | 28:39 | 0.89 | 53.6 | | 2.50 | - | * | PeCDF | 1 |
| Total Penta-Furans | 4.95e+07 | | 30:24 | 0.89 | 140 | | 2.50 | - | * | 194 | 8 |
| Total Hexa-Furans | 8.85e+07 | | 35:31 | 1.00 | 233 | | 2.50 | - | * | | 12 |
| Total Hepta-Furans | 2.68e+07 | | 42:35 | 1.46 | 101 | | 2.50 | - | * | | 14 |

Analyst: 

Date: 

Frontier Analytical Laboratory - Acquisition Log

Run Name:07APR11M

Instrument: FAL3

GC: DB5

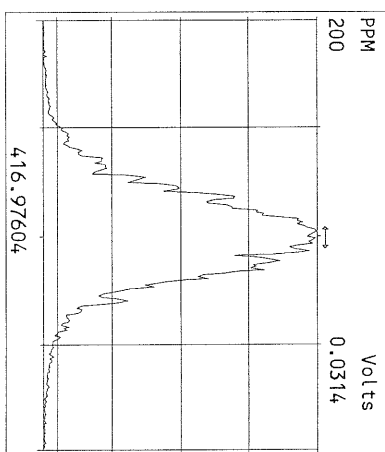
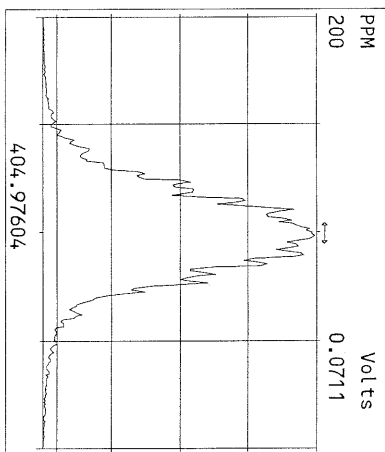
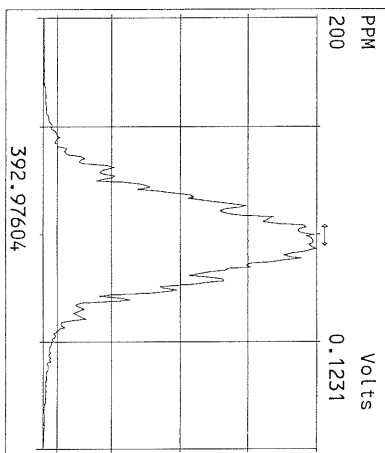
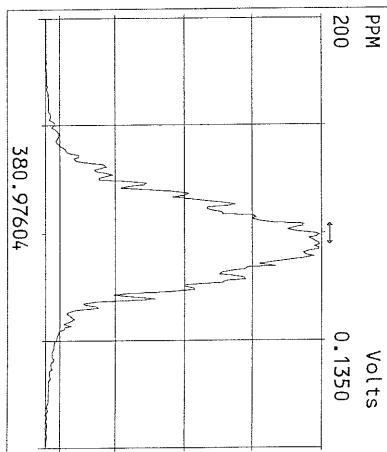
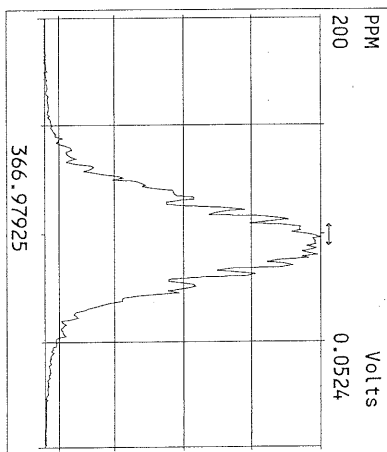
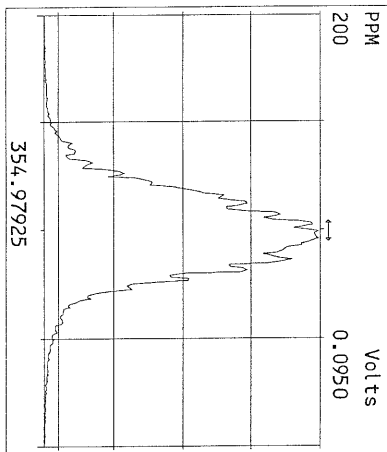
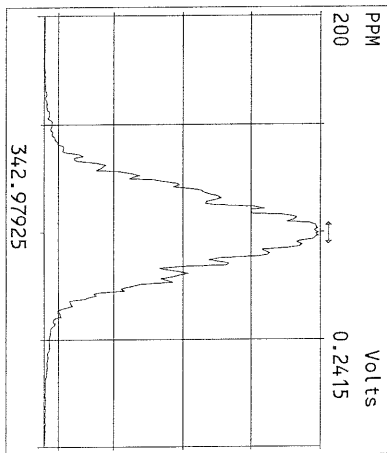
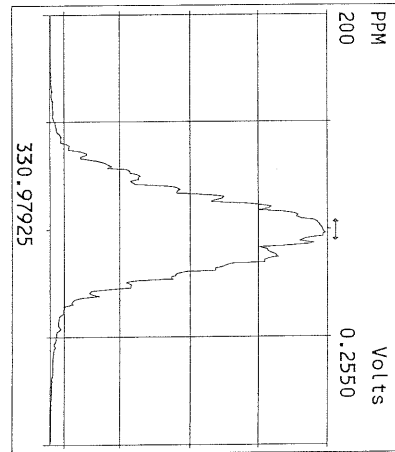
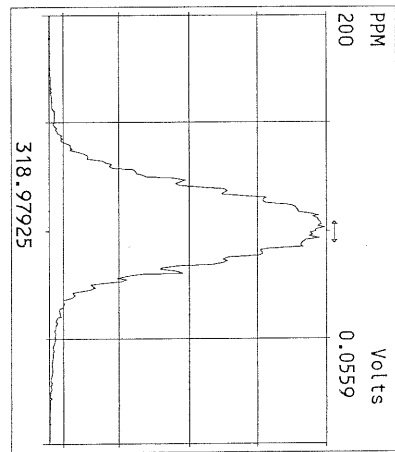
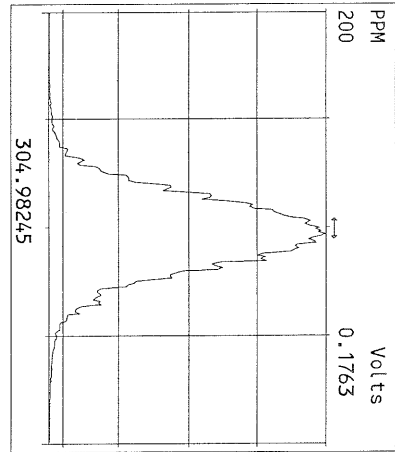
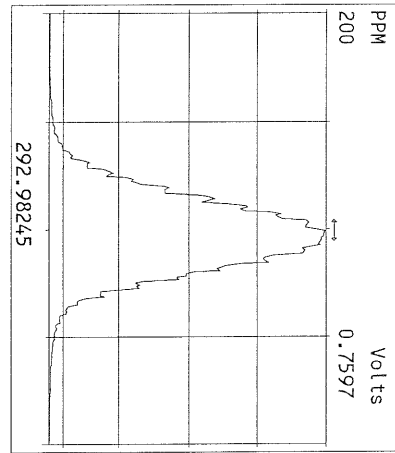
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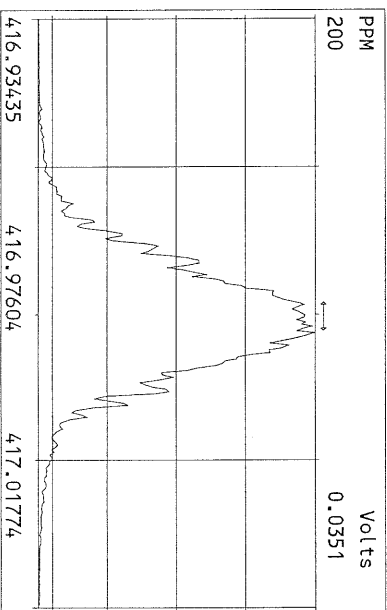
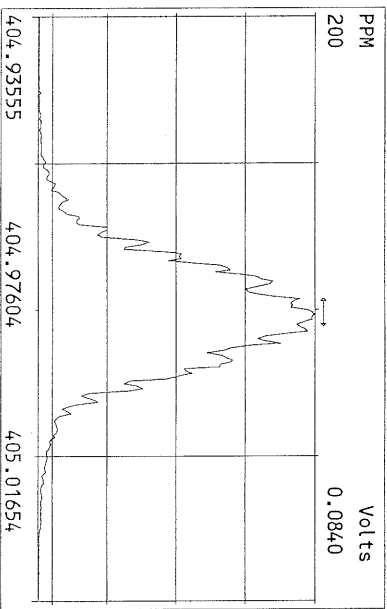
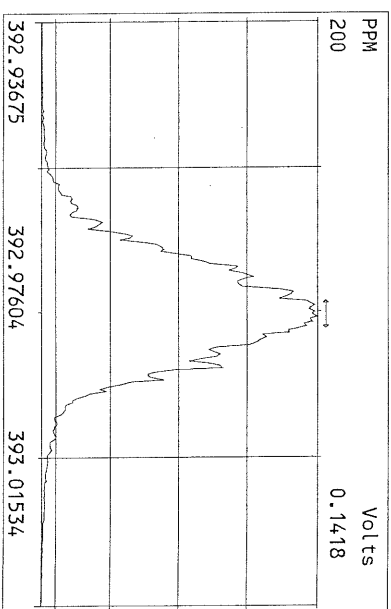
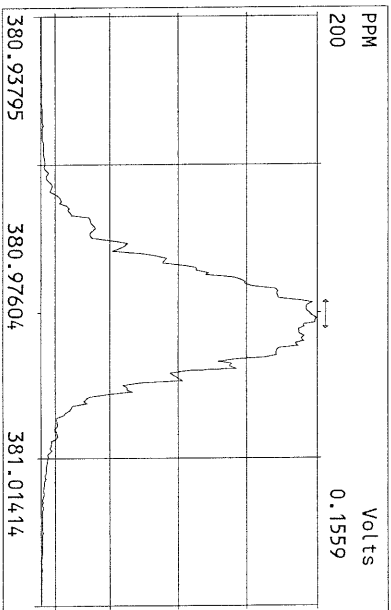
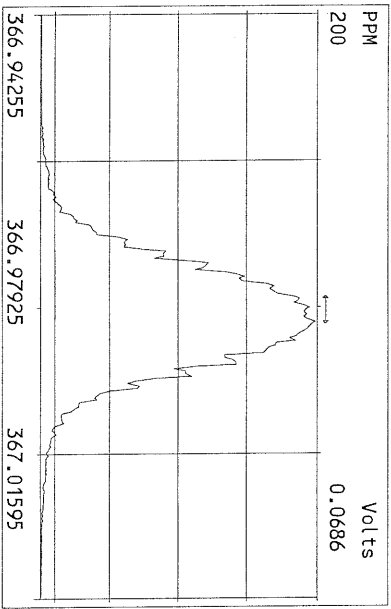
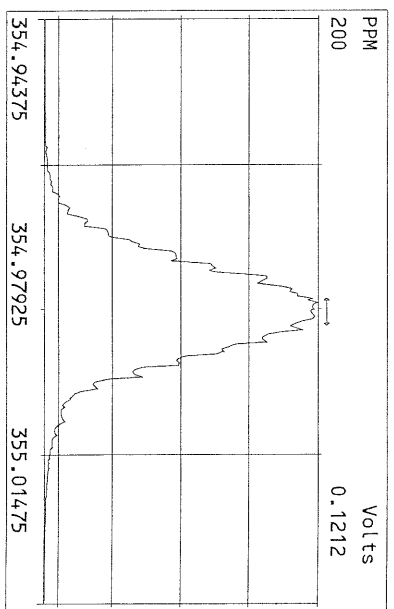
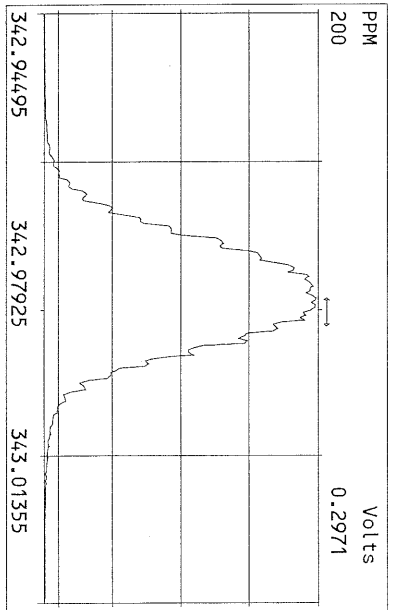
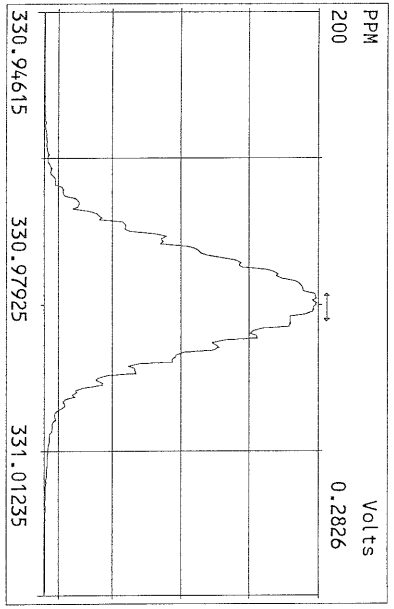
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| 07APR11M 2 | 2264-001-0001-MB | Method Blank | 7-APR-11 16:08:34 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 3 | 2264-001-0001-OPR | OPR | 7-APR-11 17:03:57 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 4 | 6678-008-0001-SA | LL-SED1-0-56-031511 | 7-APR-11 17:59:16 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 5 | 6678-007-0001-SA | LL-SED2-0-56-031511-D | 7-APR-11 18:54:38 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 6 | 6678-006-0001-SA | LL-SED2-112-168-031511 | 7-APR-11 19:49:58 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 7 | 6678-005-0001-SA | LL-SED2-56-112-031511 | 7-APR-11 20:45:17 | ST040711M1 | ST040711M2 | TC |
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| 07APR11M 9 | 6678-003-0001-SA | LL-SED3-141-167-031511 | 7-APR-11 22:36:00 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 10 | 6678-002-0001-SA | LL-SED3-36-141-031511 | 7-APR-11 23:31:19 | ST040711M1 | ST040711M2 | TC |
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| 07APR11M 14 | SB040711M1 | Solvent Blank | 8-APR-11 03:12:55 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 15 | ST040711M2 | 1613 CS3 100511J | 8-APR-11 04:08:18 | ST040711M2 | ST040711M3 | TC |
| 07APR11M 16 | SB040711M2 | Solvent Blank | 8-APR-11 05:03:38 | ST040711M2 | ST040711M3 | TC |
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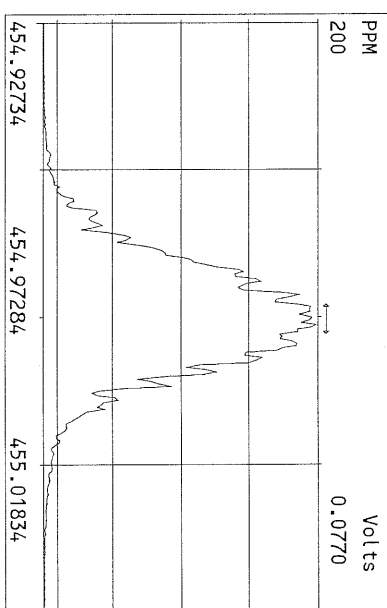
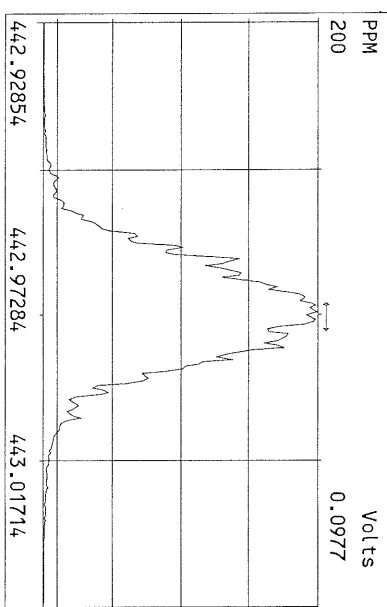
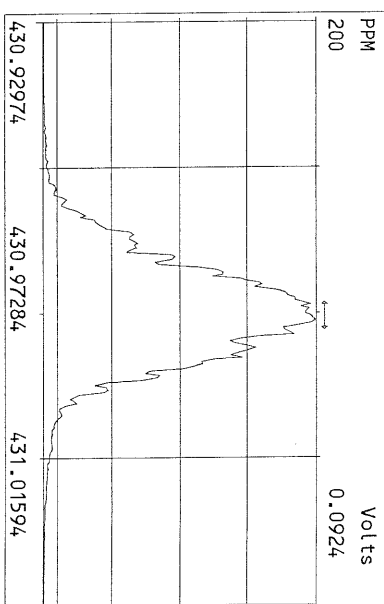
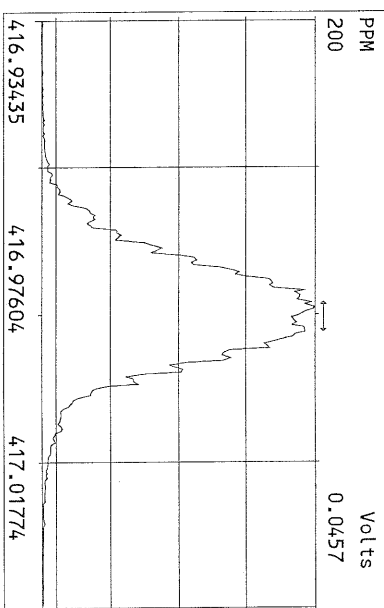
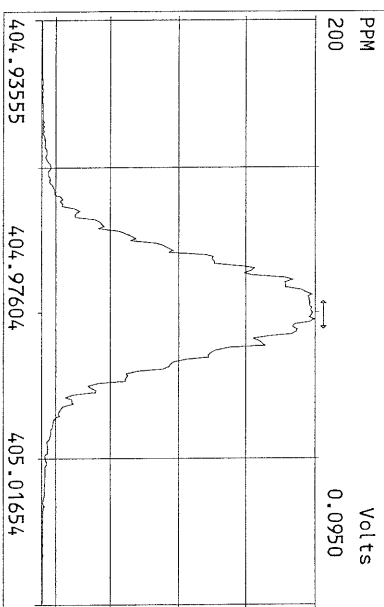
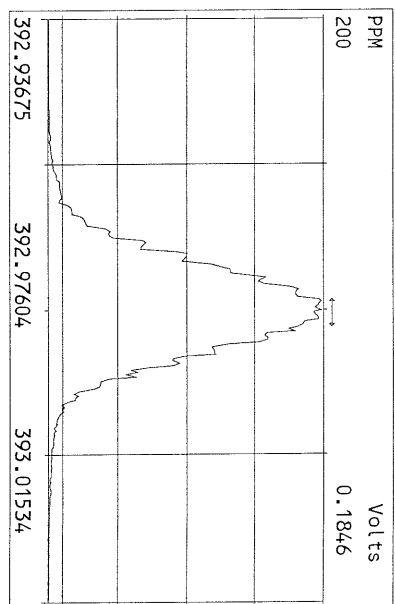
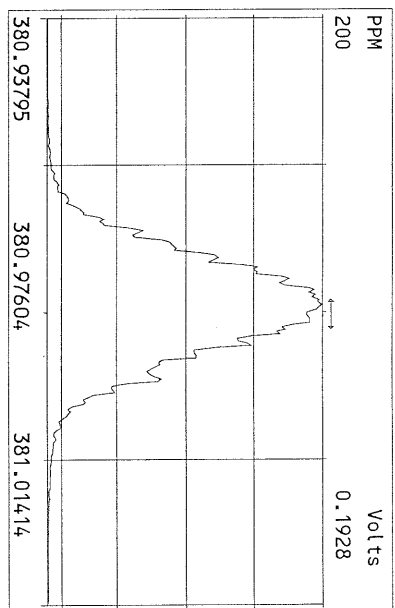
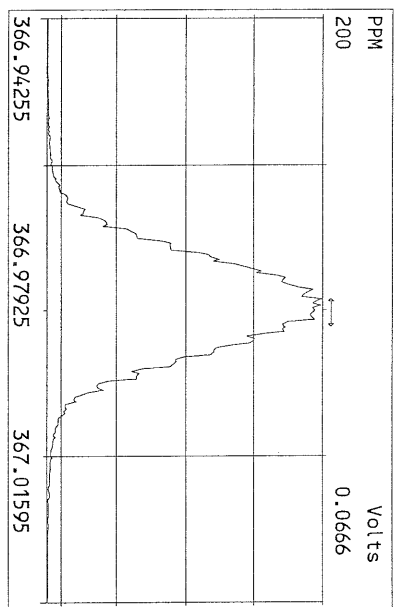
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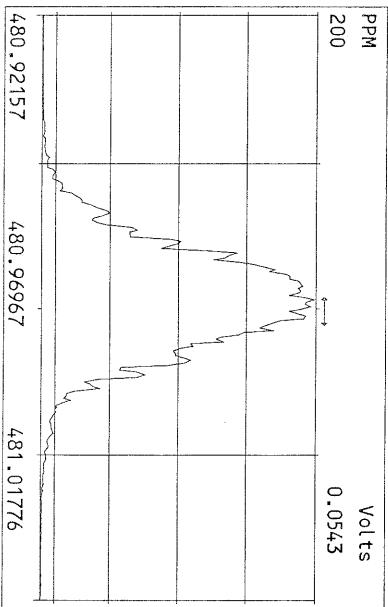
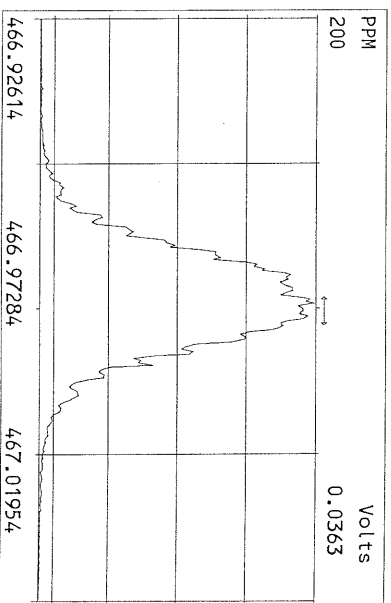
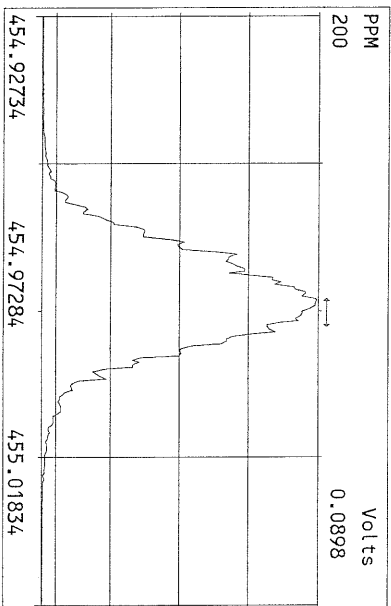
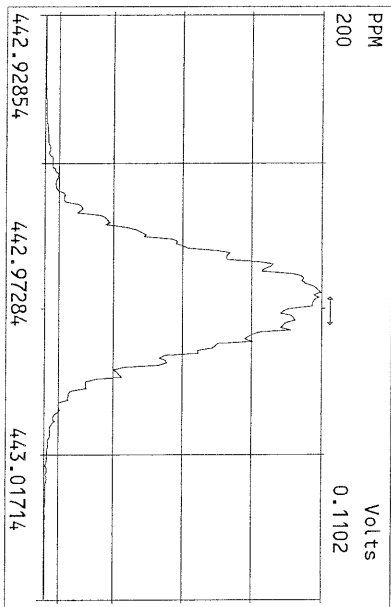
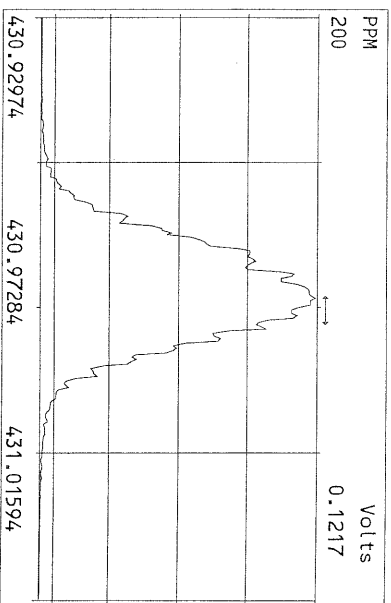
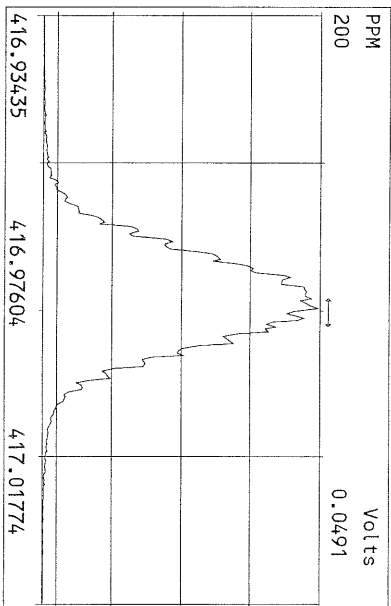
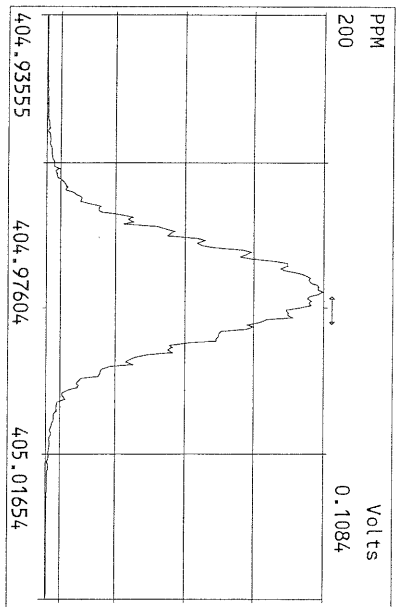
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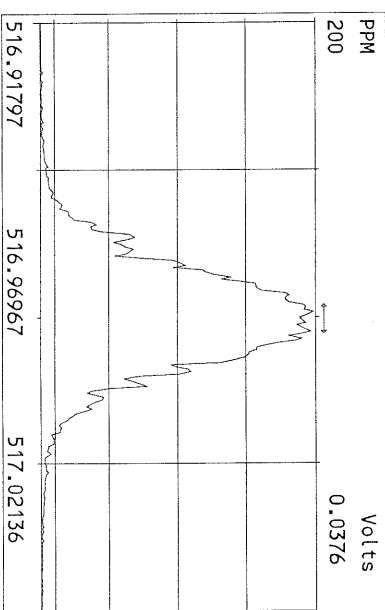
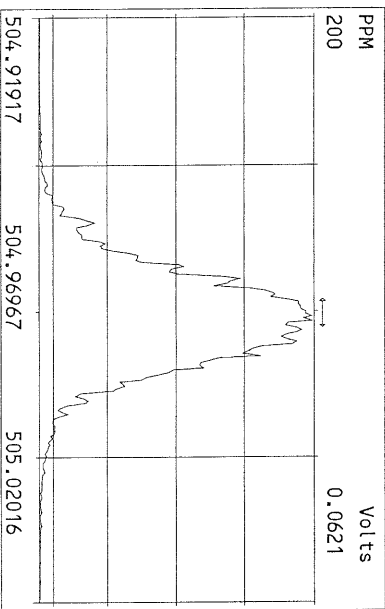
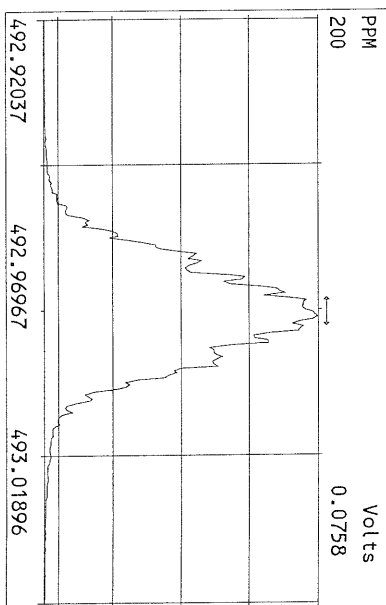
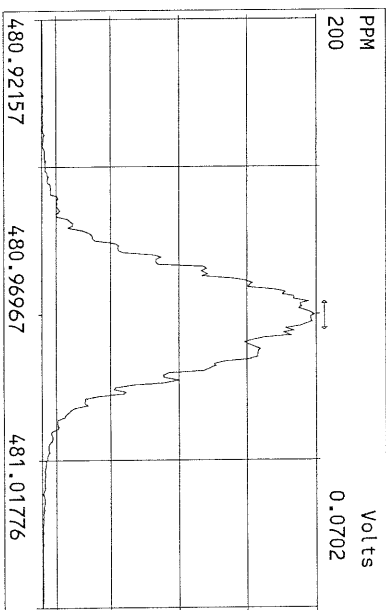
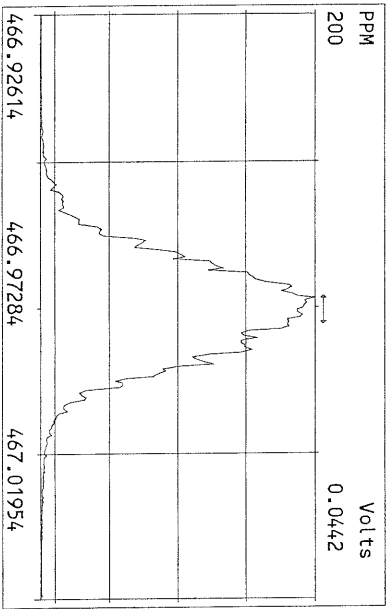
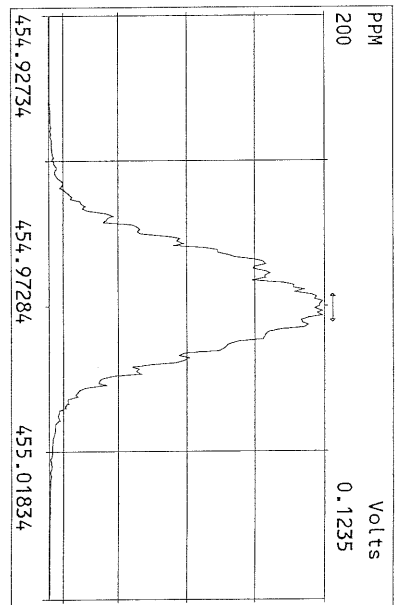
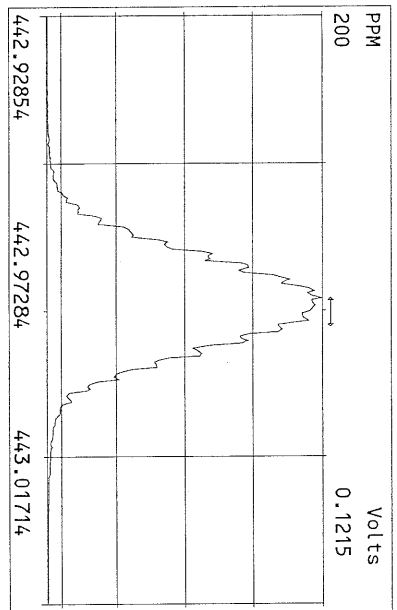
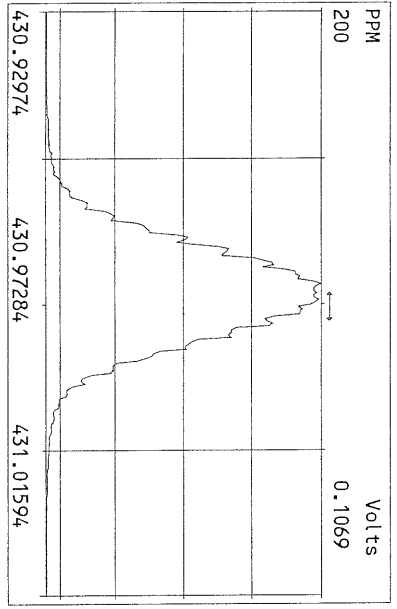
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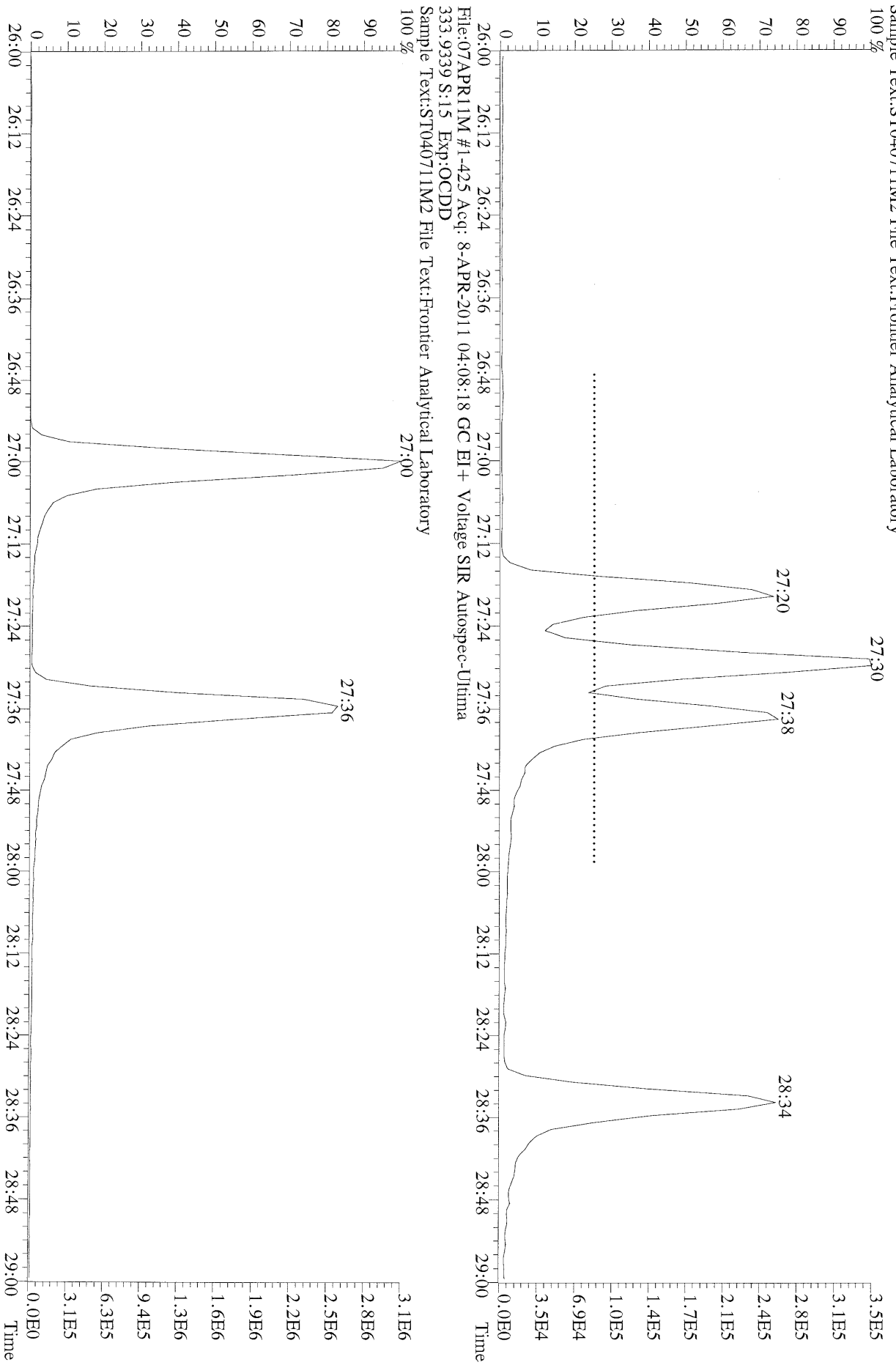




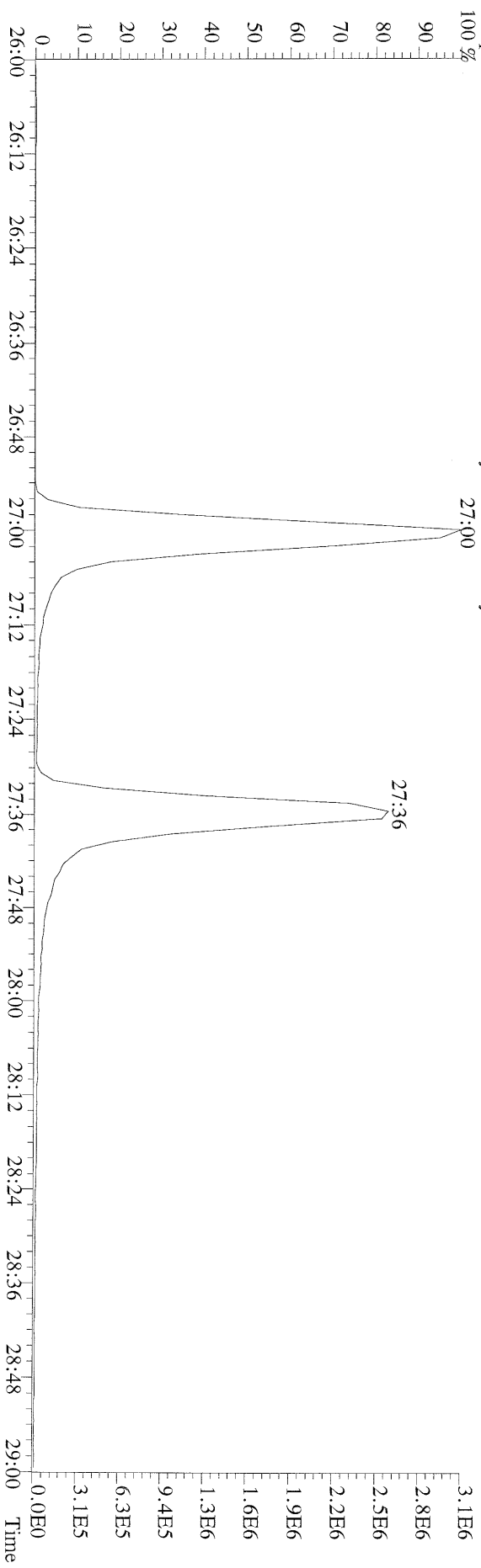




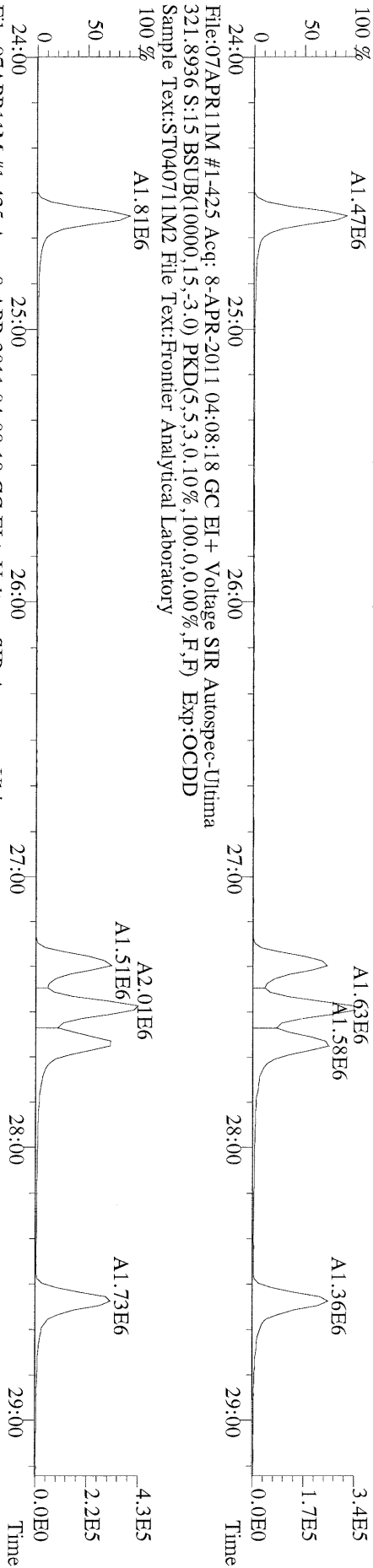
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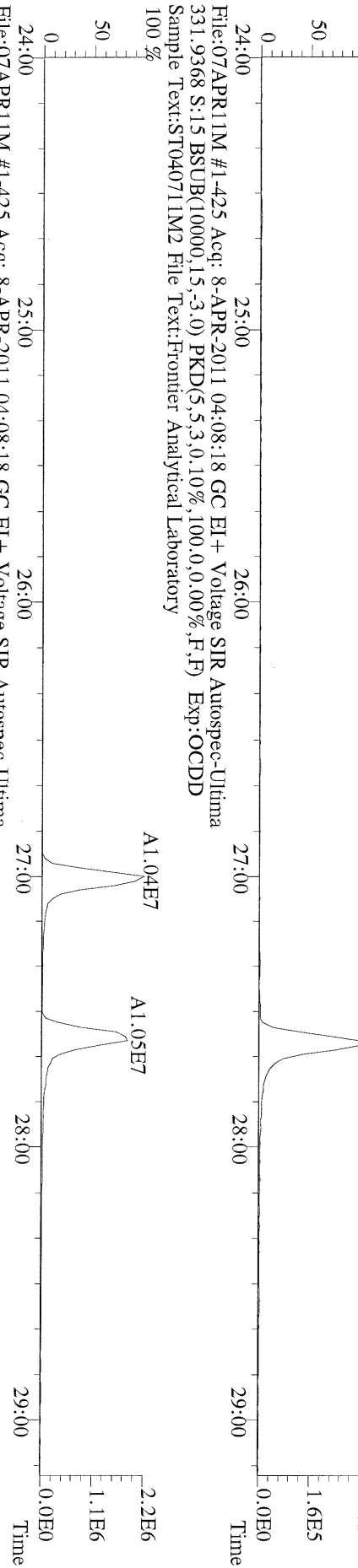
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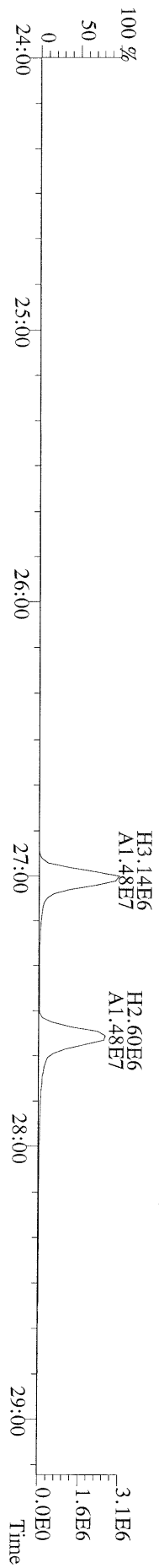
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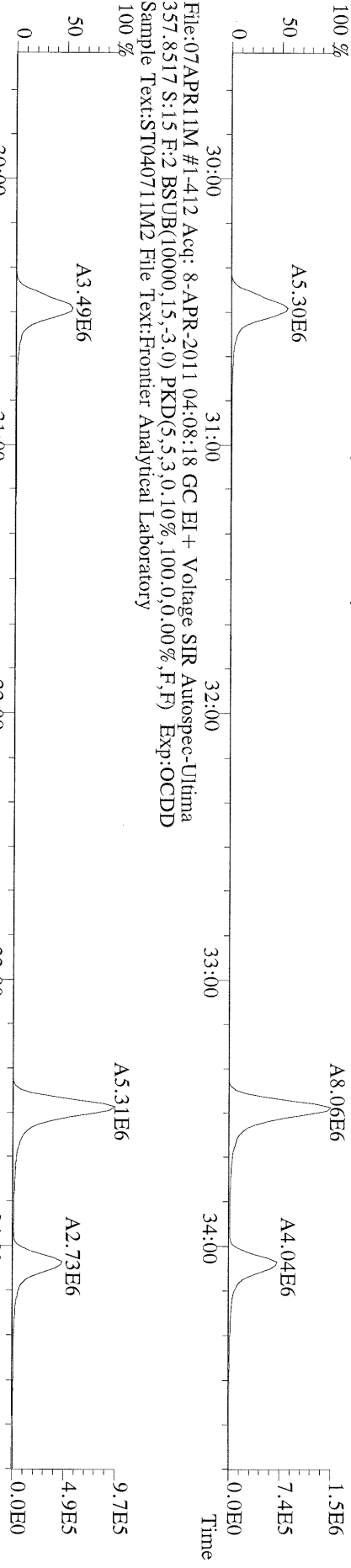
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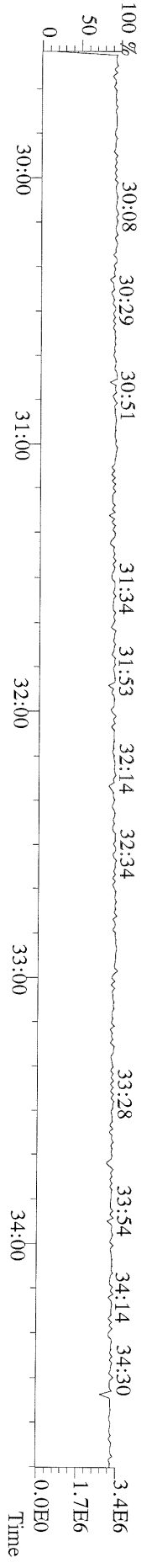
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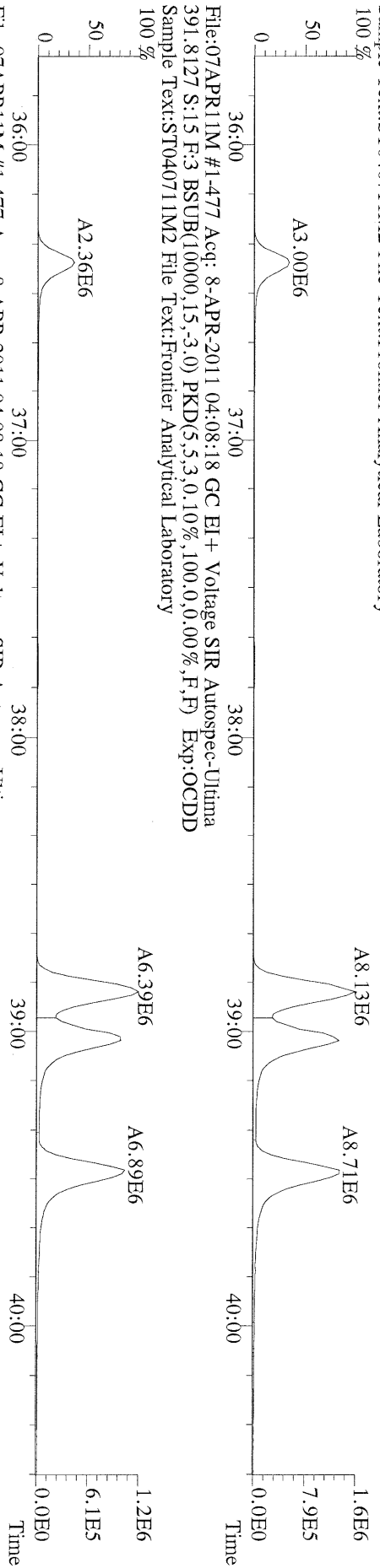
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100 %



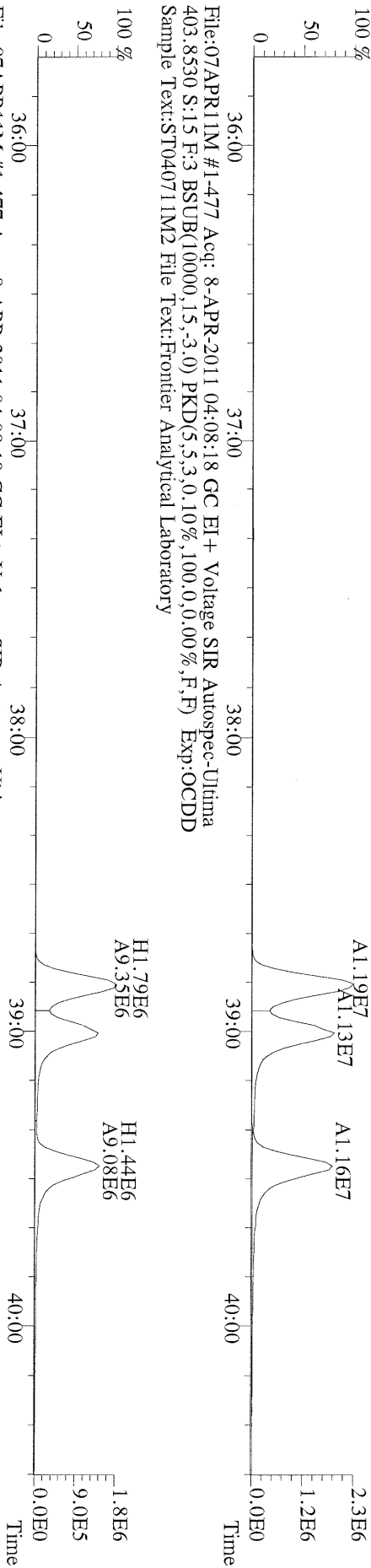
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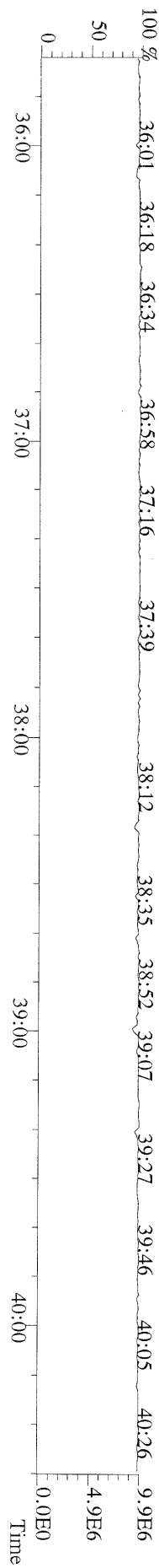
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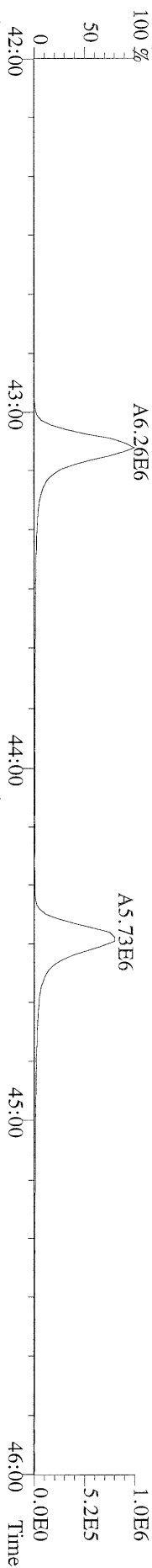
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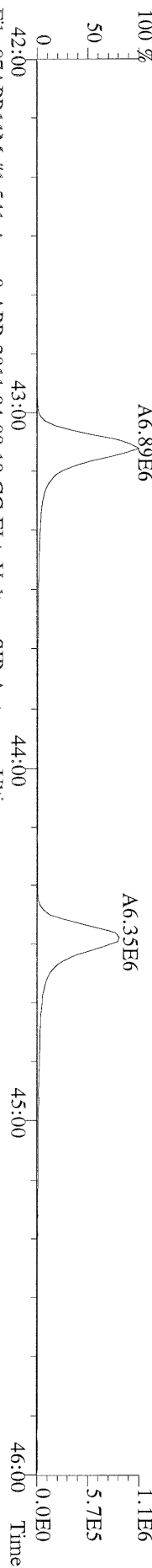
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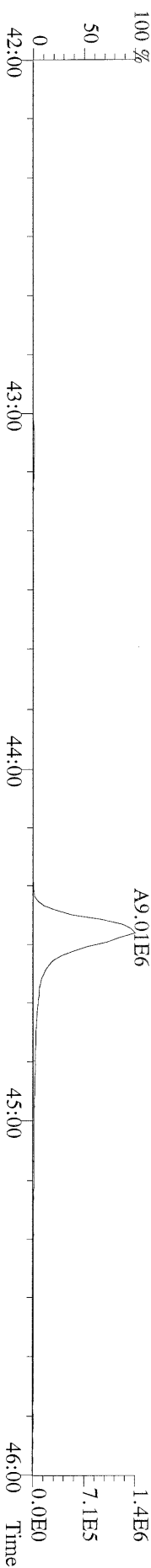
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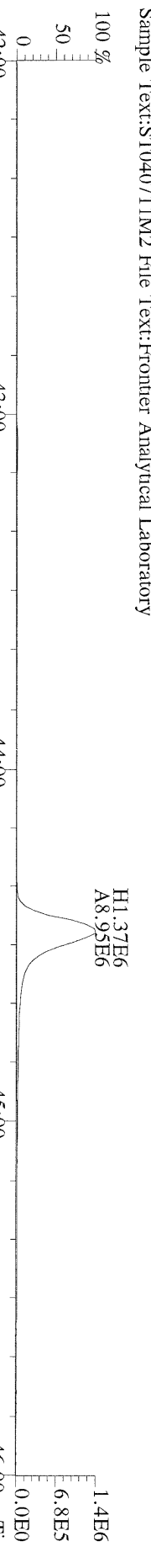
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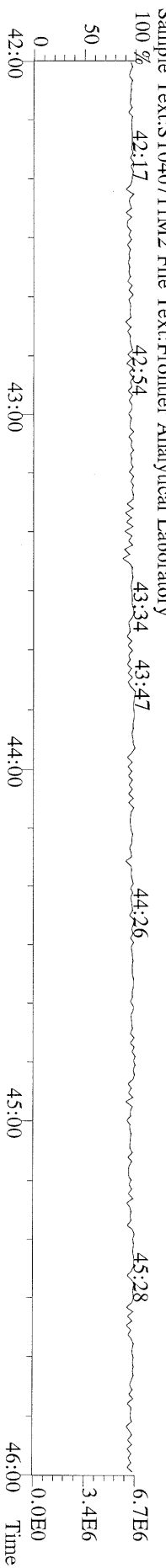
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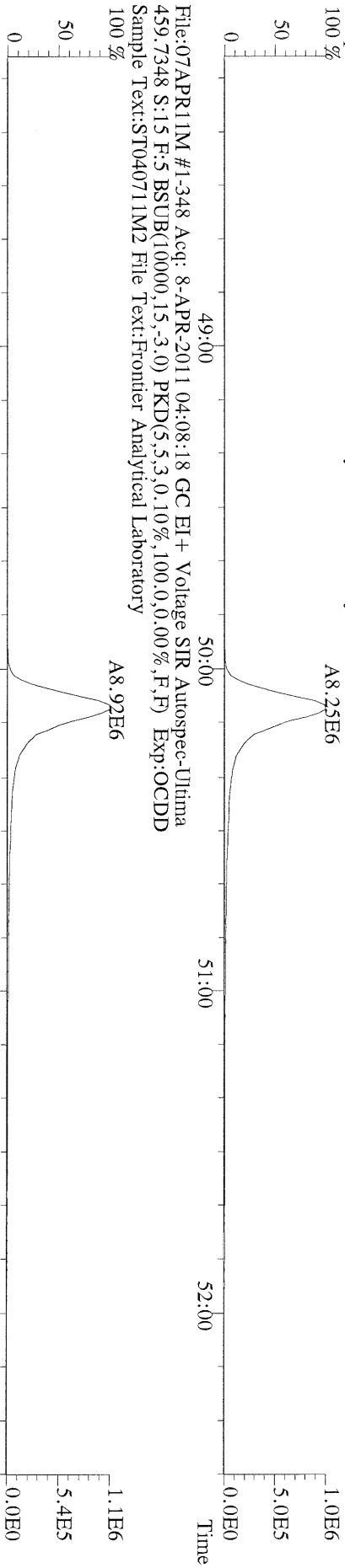
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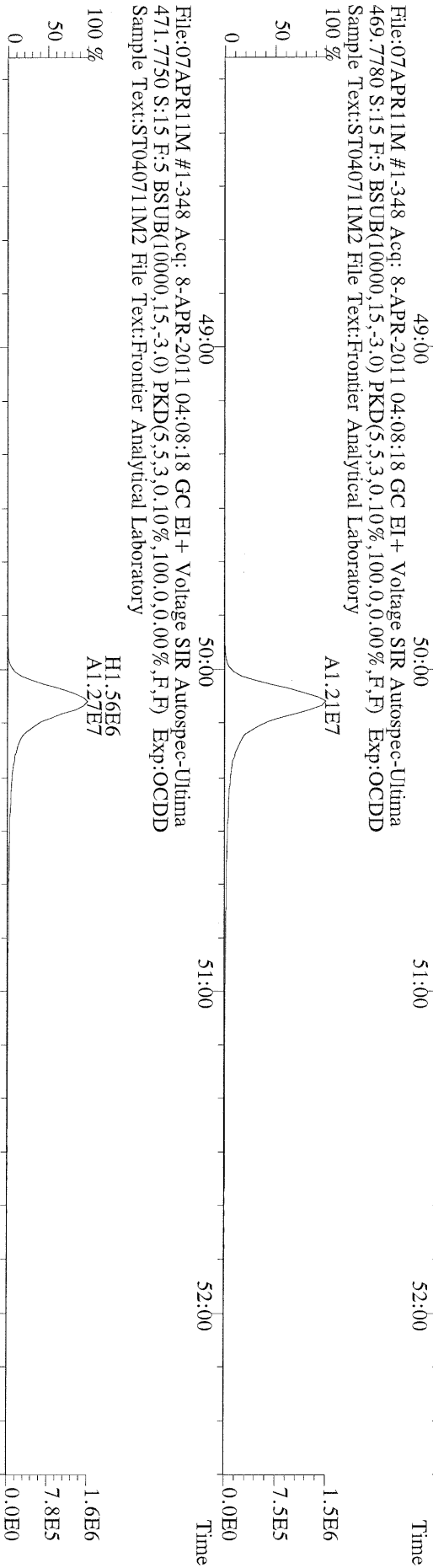
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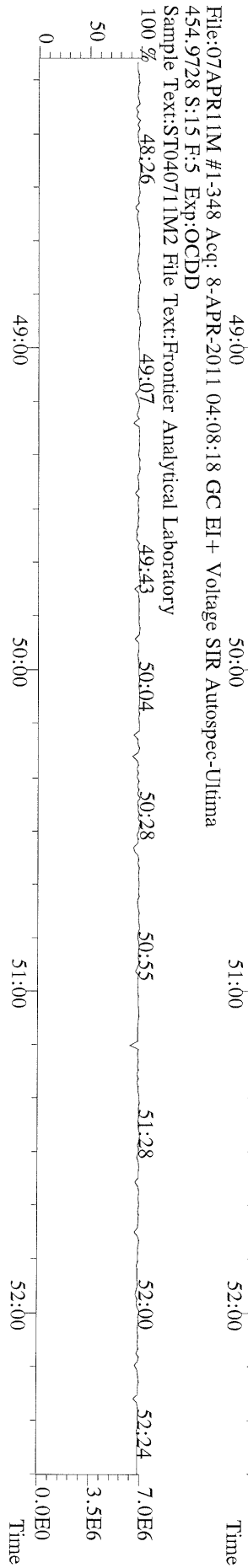
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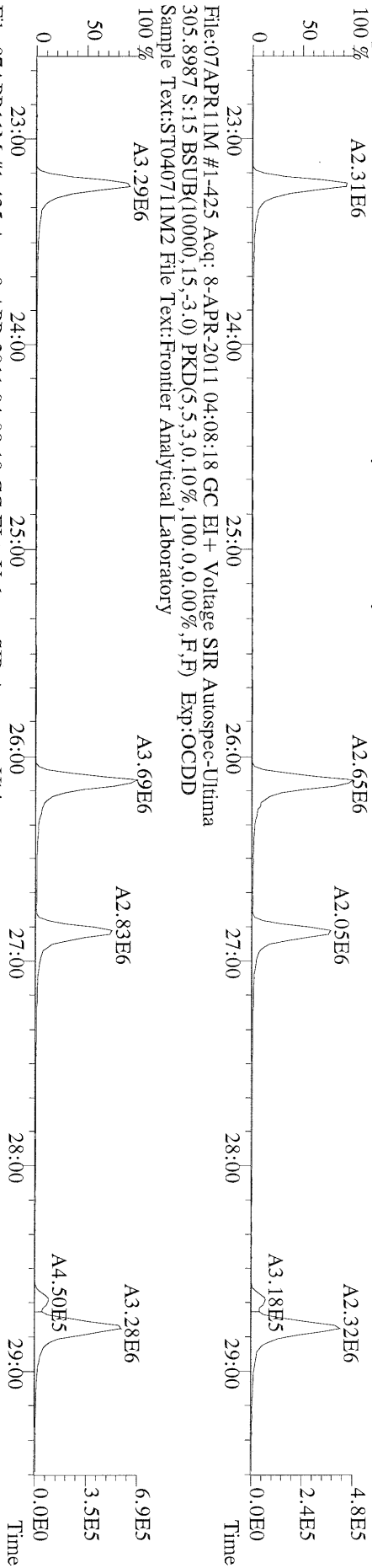
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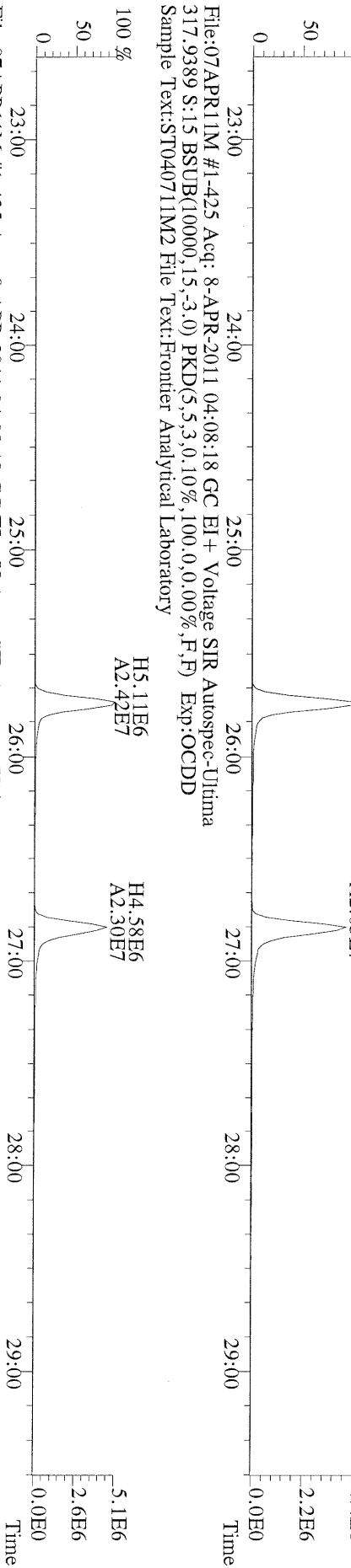
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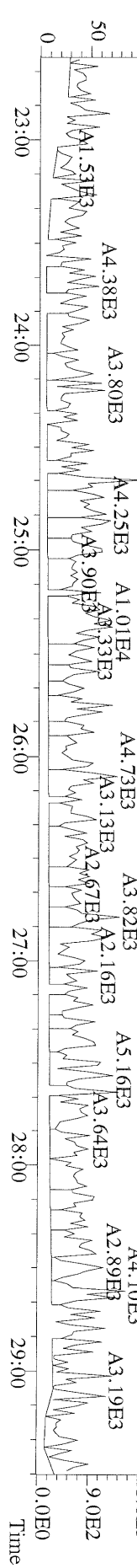
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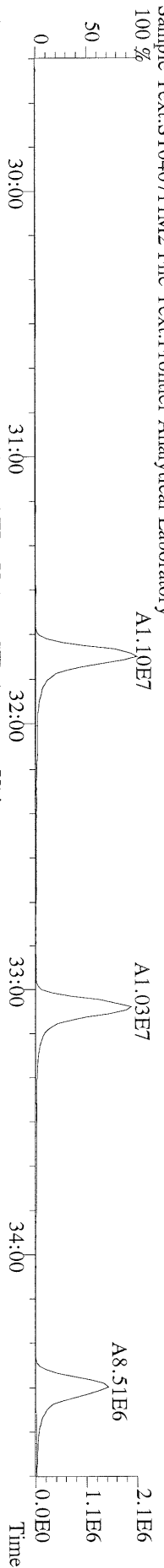
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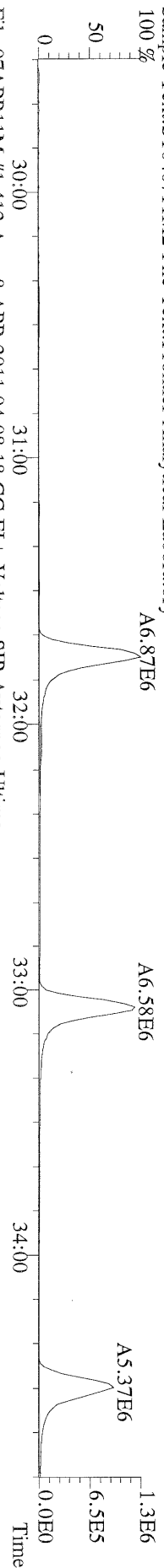
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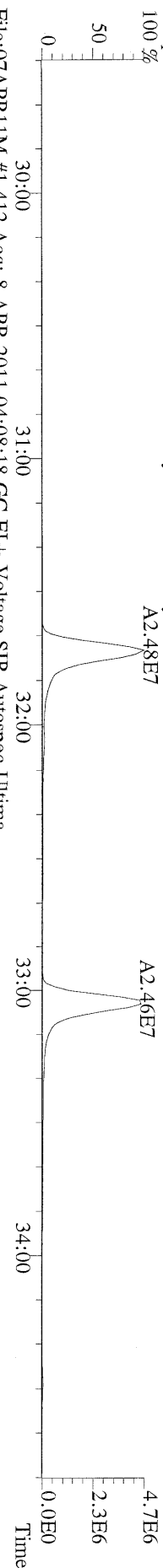
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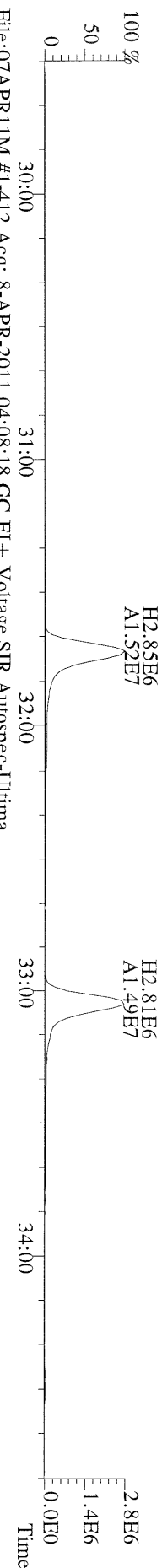
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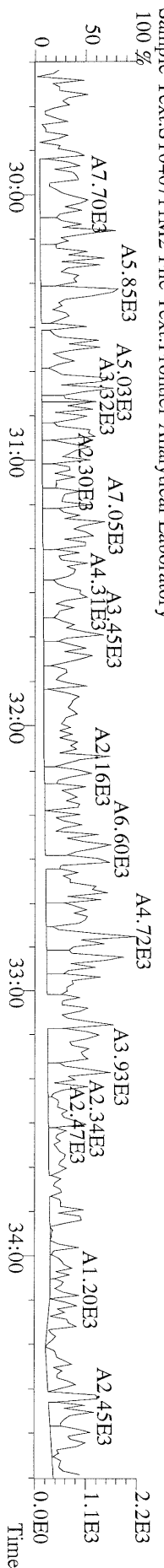
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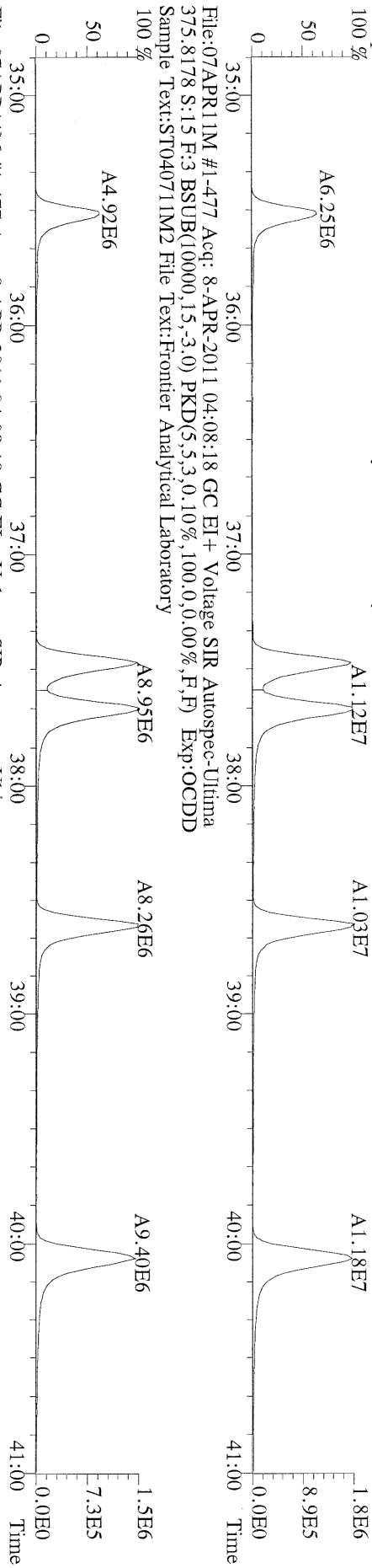
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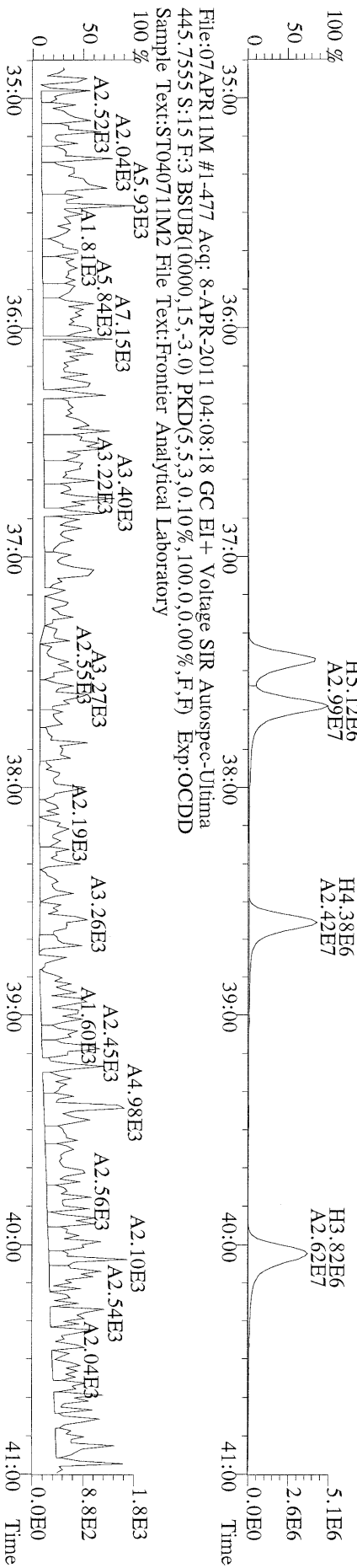
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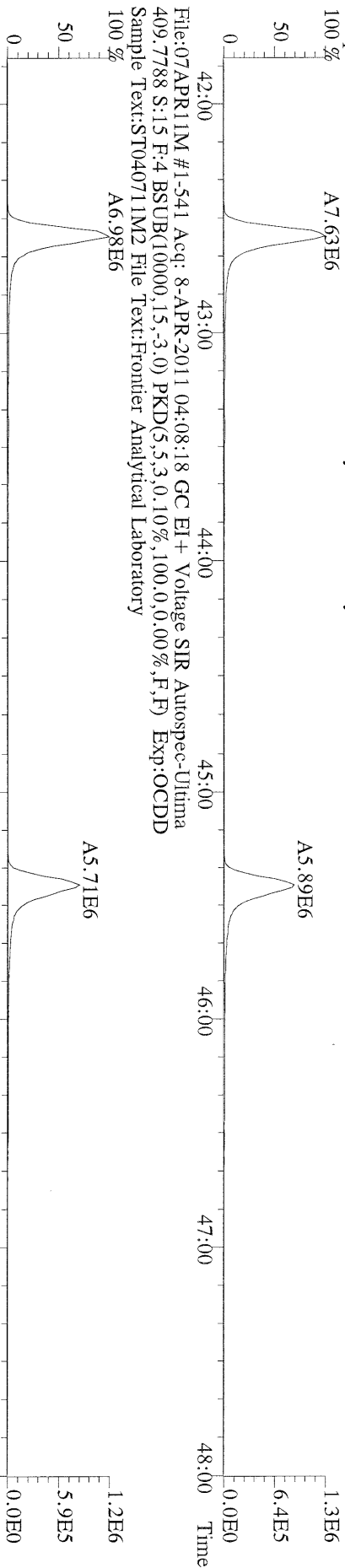
File:07APR11M #1-477 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Ultima
373.8207 S:15 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040711M2 File Text:Frontier Analytical Laboratory



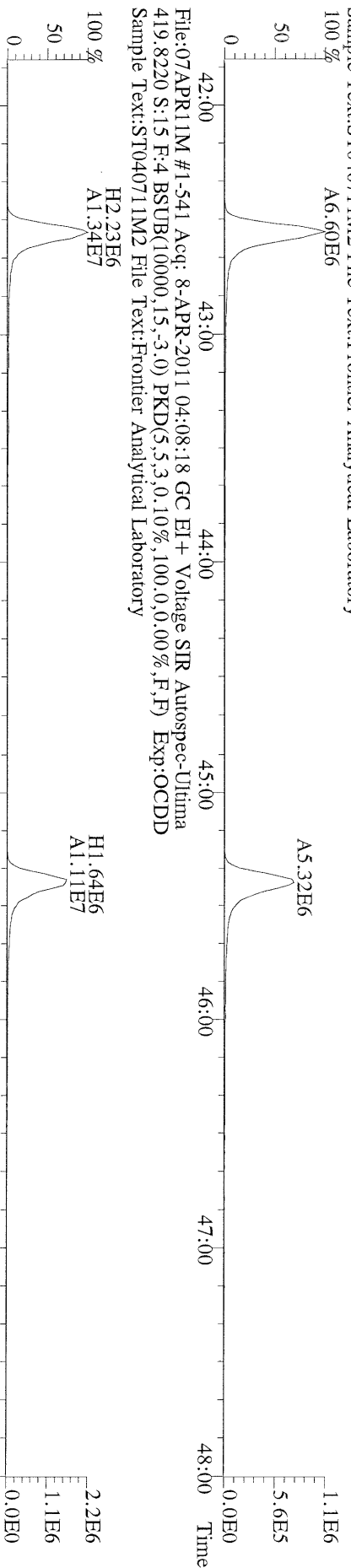
File:07APR11M #1-477 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Ultima
383.8639 S:15 F:3 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100.0,0.00%,F,F) Exp:OCDD
Sample Text:ST040711M2 File Text:Frontier Analytical Laboratory



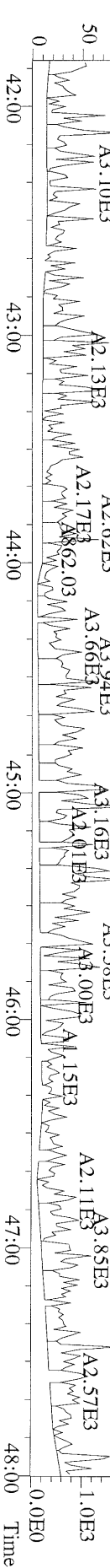
File:07APR11M #1-541 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Utima
407.7818 S:15 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory



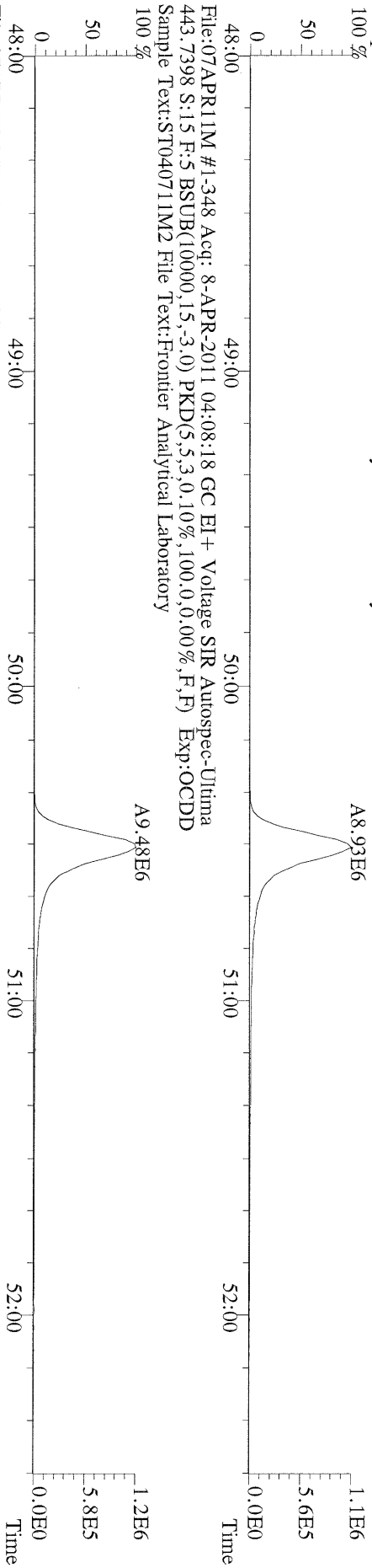
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417.8253 S:15 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory



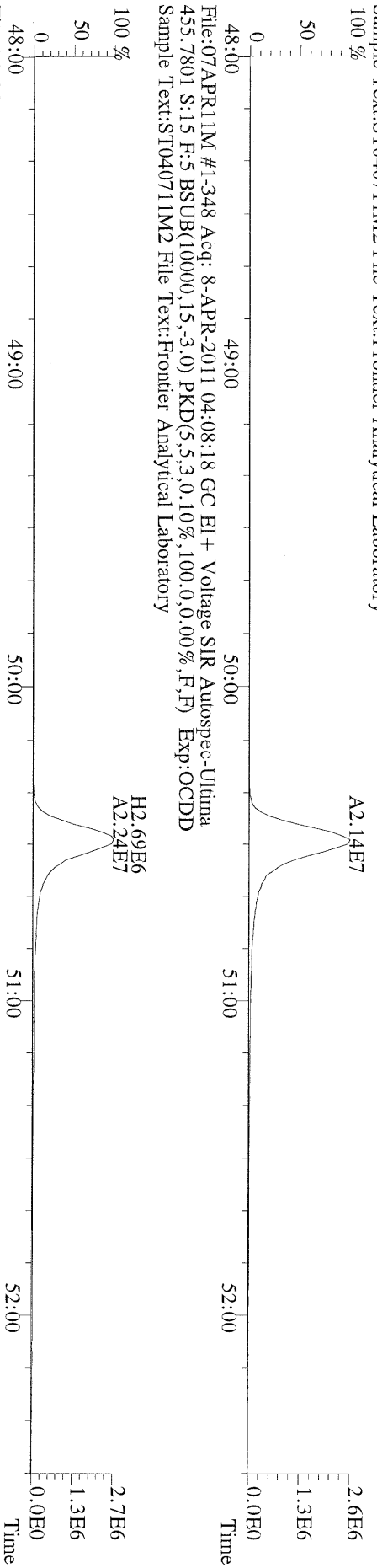
File:07APR11M #1-541 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Utima
419.8220 S:15 F:4 BSUB(10000,15,-3,0) PKD(5,5,3,0,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory



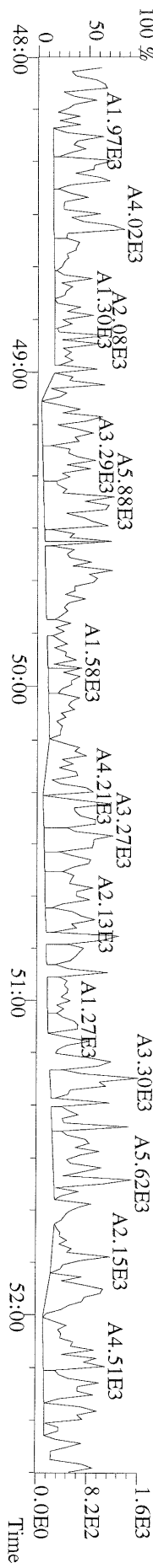
File:07APR11M #1-348 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Ultima
441.7428 S:15 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory
100 %



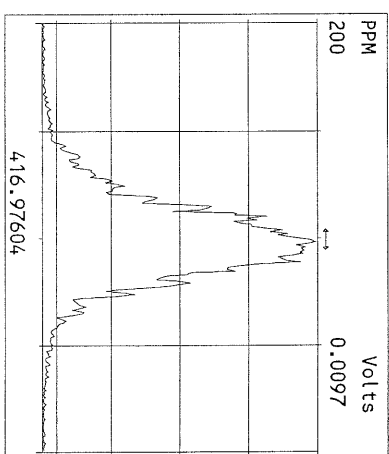
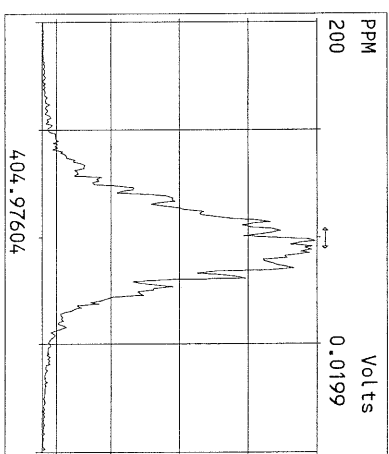
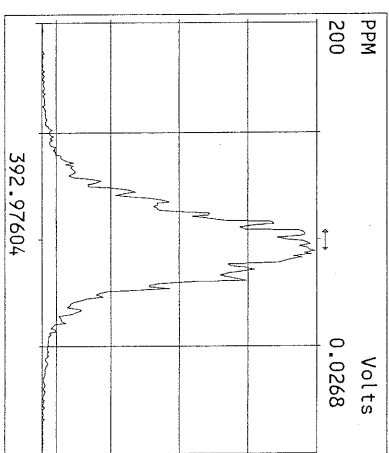
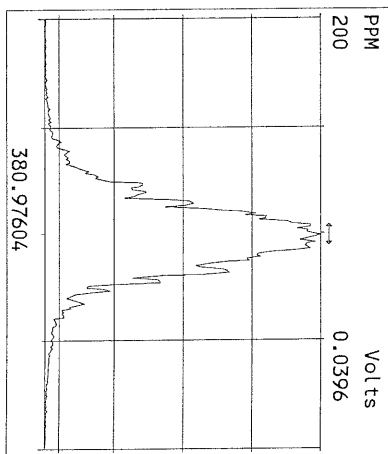
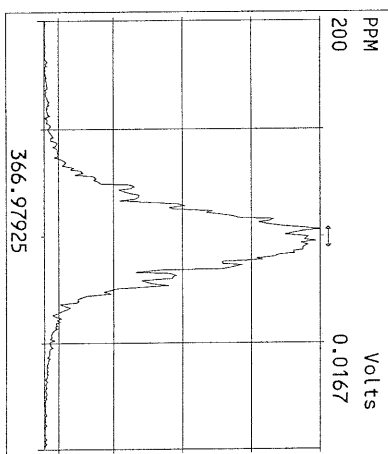
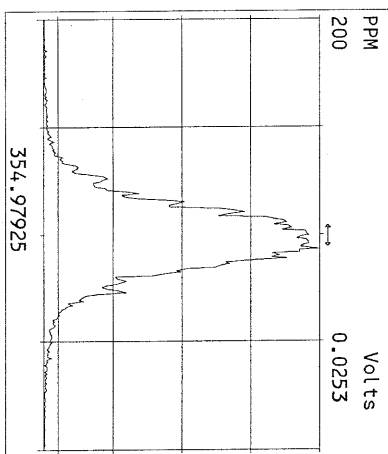
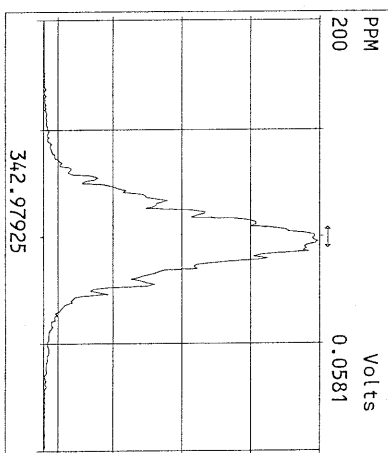
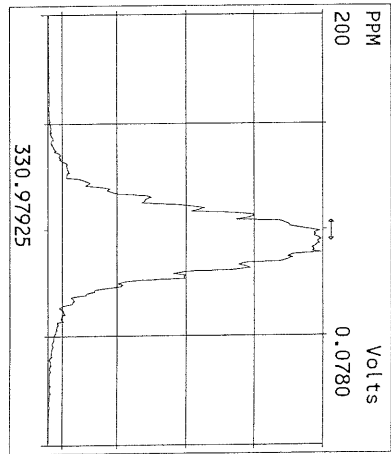
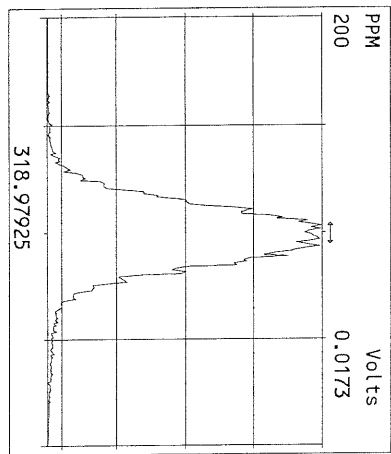
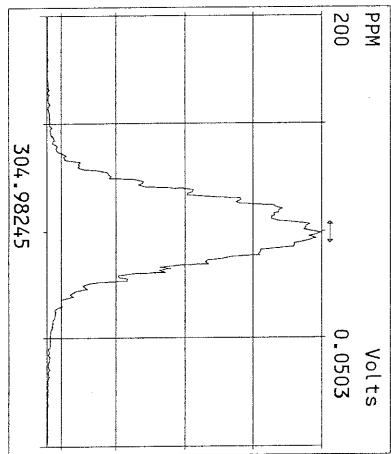
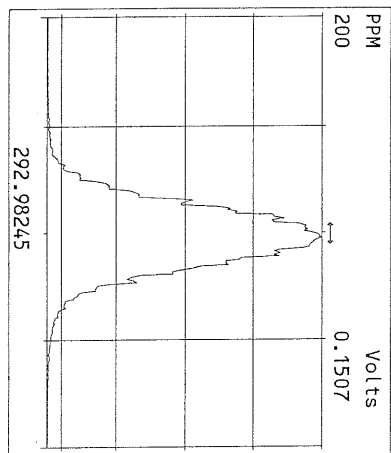
File:07APR11M #1-348 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Ultima
453.7831 S:15 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory
100 %

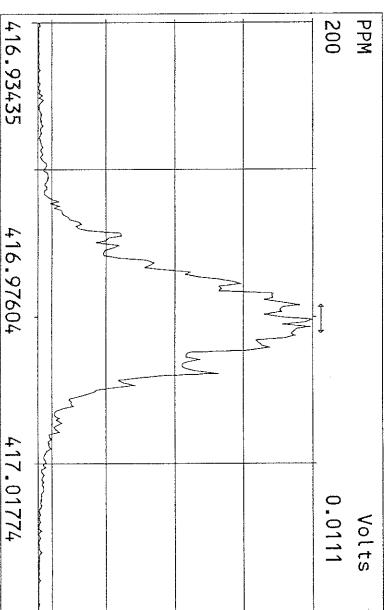
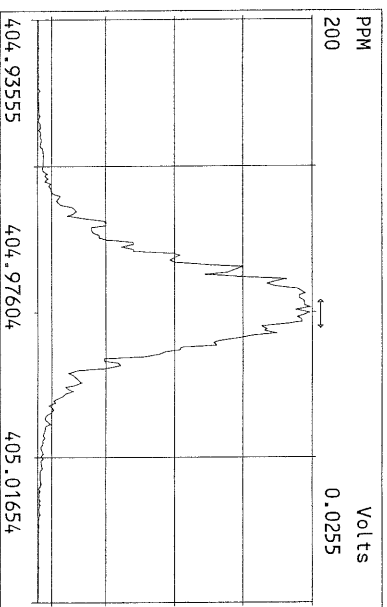
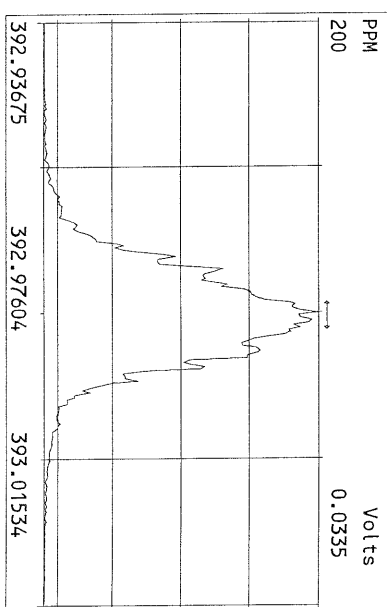
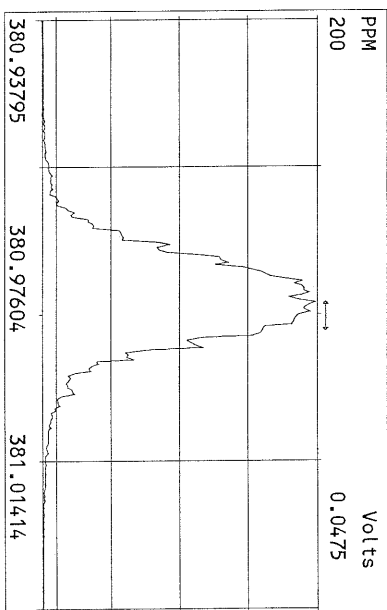
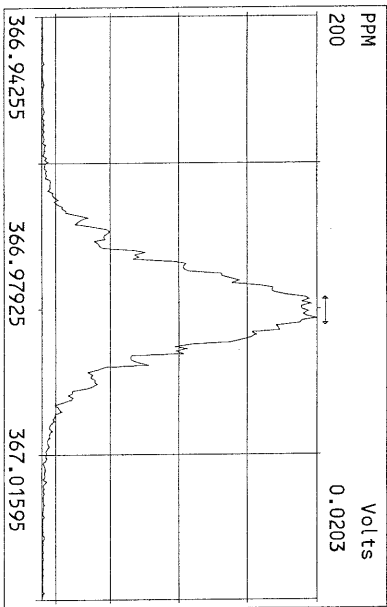
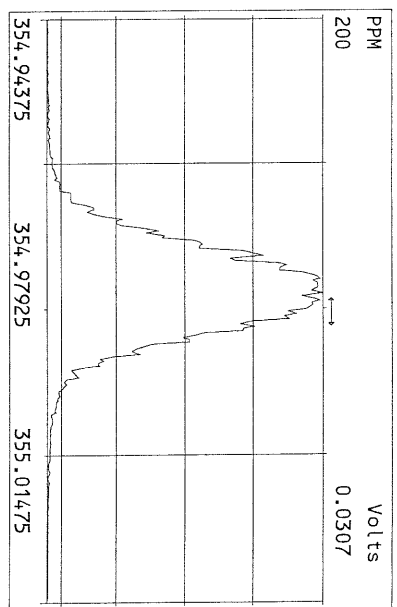
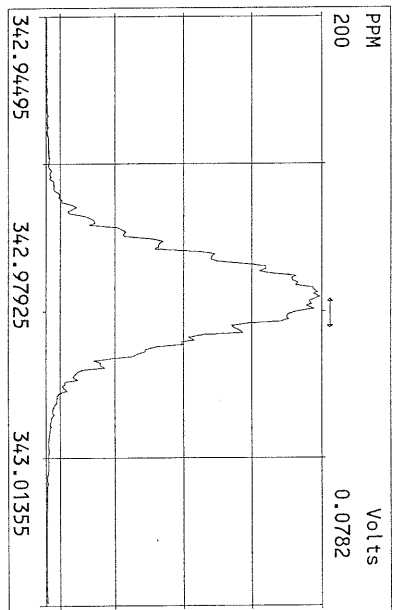
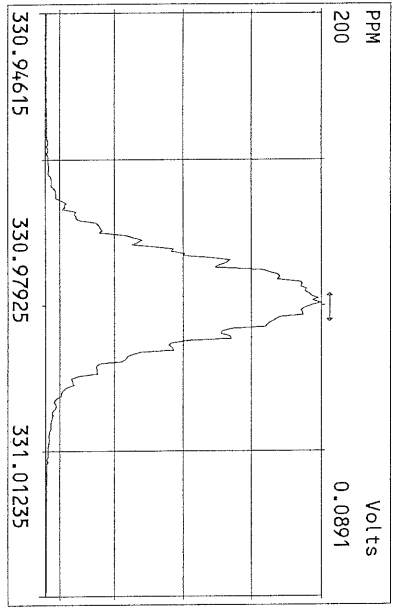


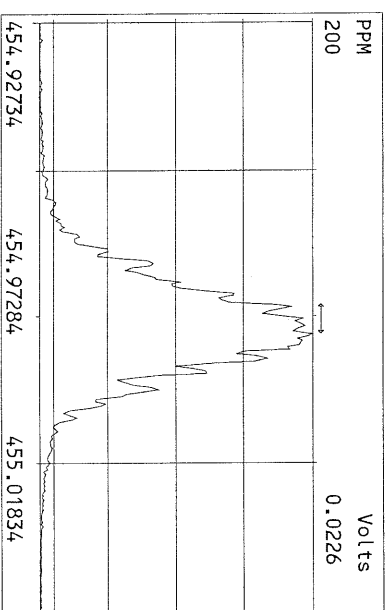
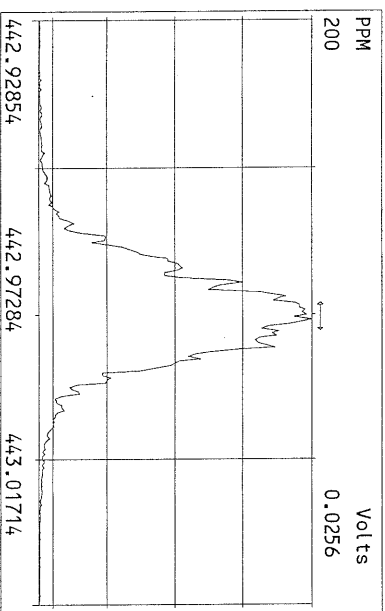
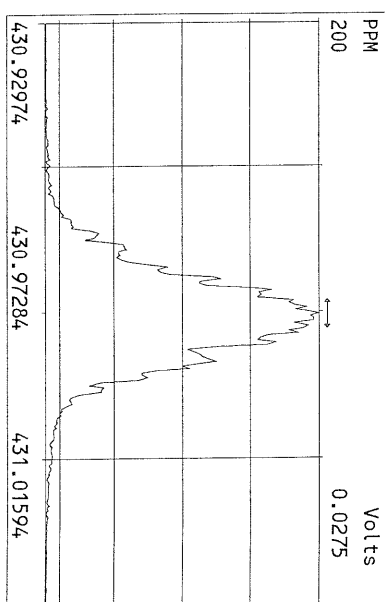
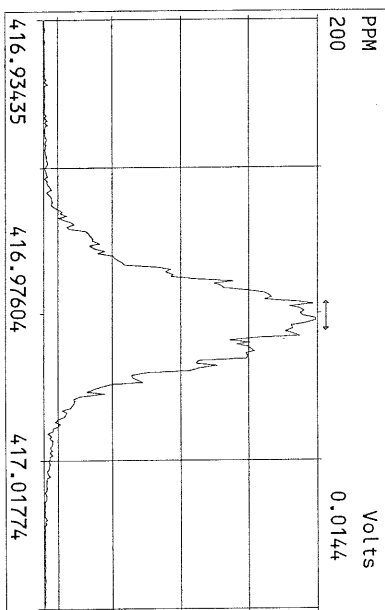
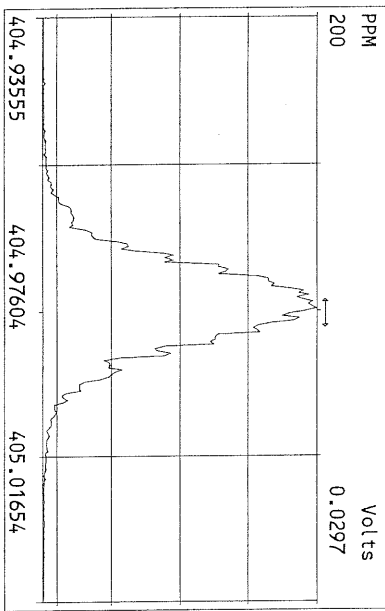
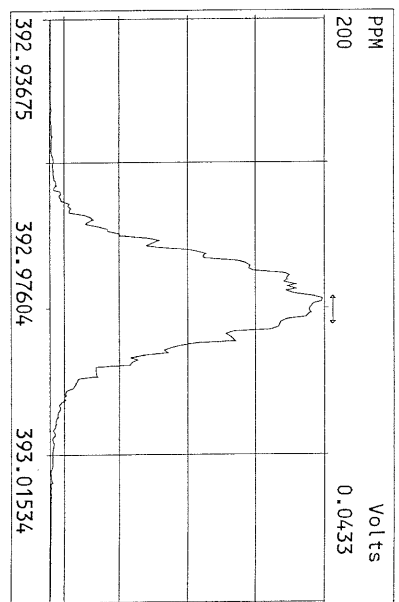
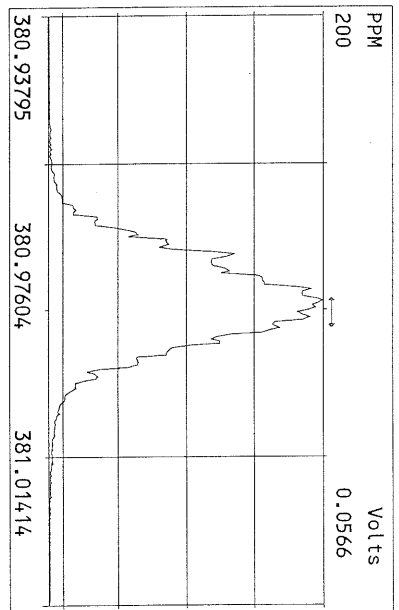
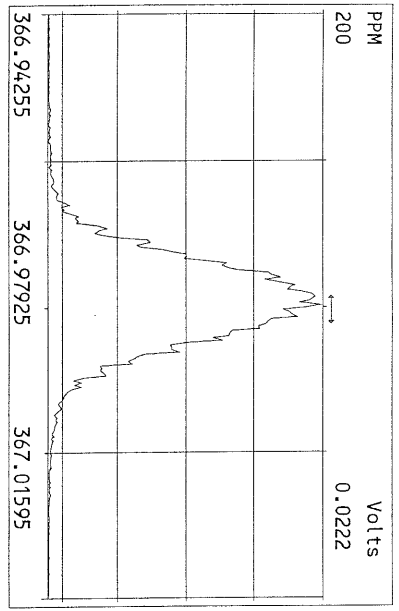
File:07APR11M #1-348 Acq: 8-APR-2011 04:08:18 GC EI+ Voltage SIR Autospec-Ultima
513.6775 S:15 F:5 BSUB(10000,15,-3,0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM2 File Text:Frontier Analytical Laboratory
100 %



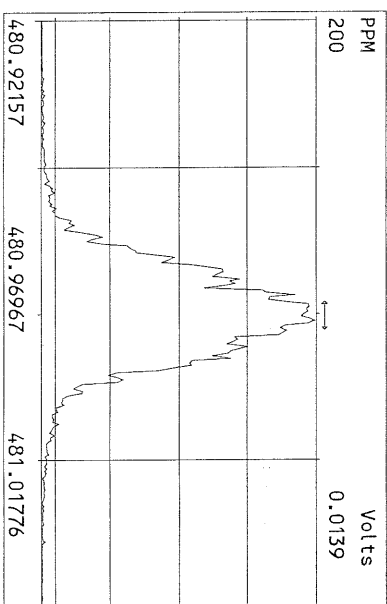
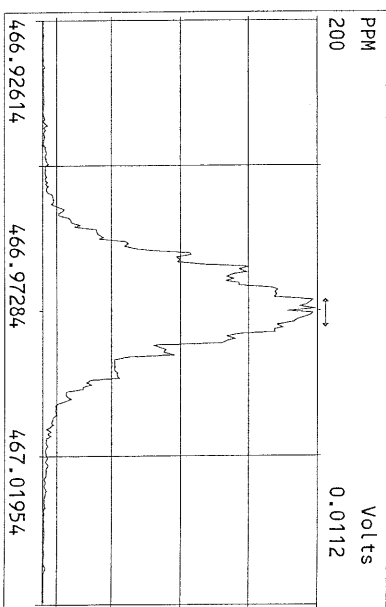
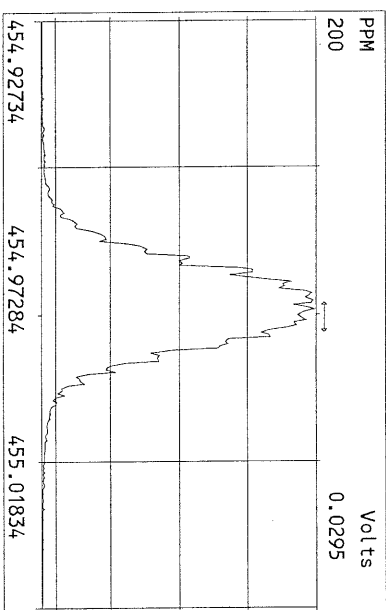
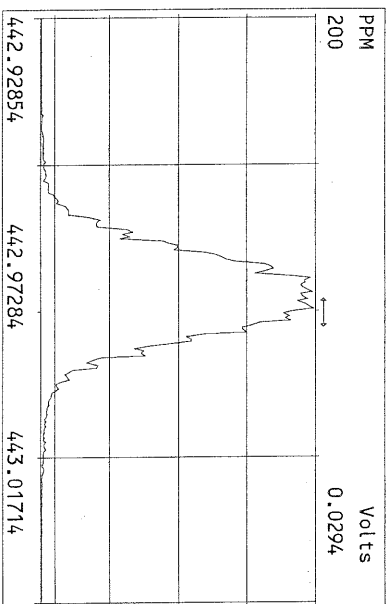
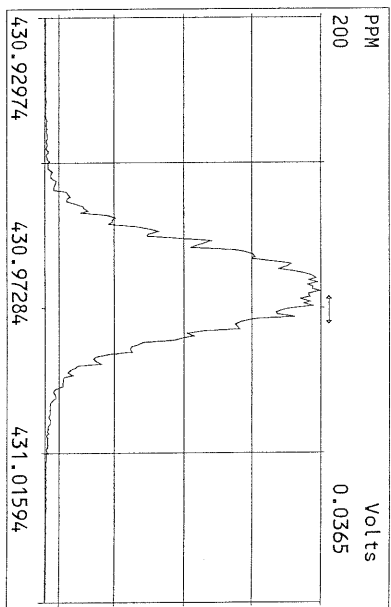
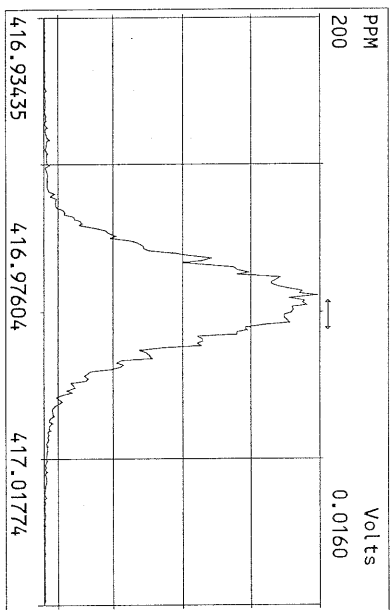
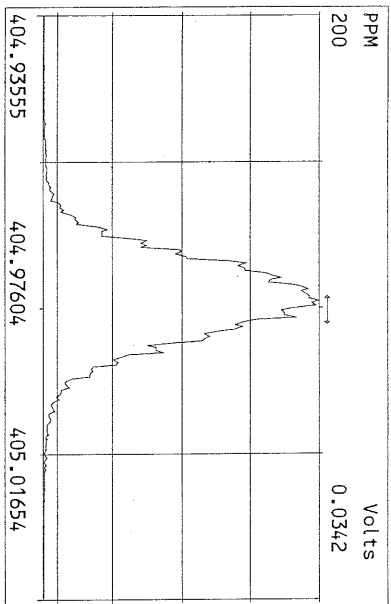
Peak Locate Examination: 8-APR-2011:08:48 File:07APR11M_RES_CHECK
Experiment:OCDD Function:1 Reference:PFK

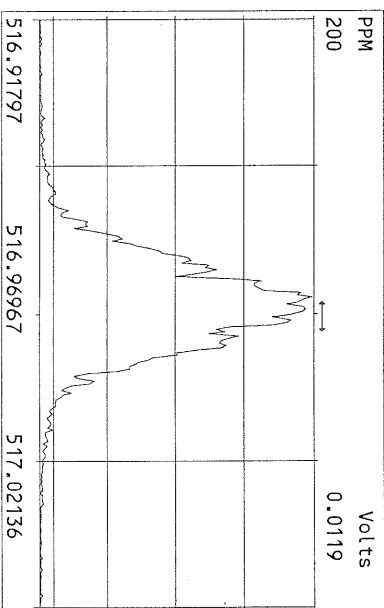
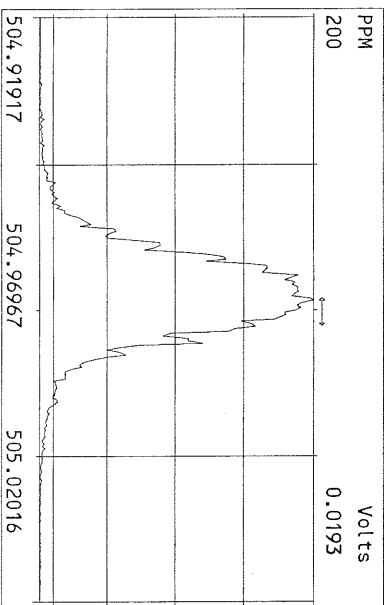
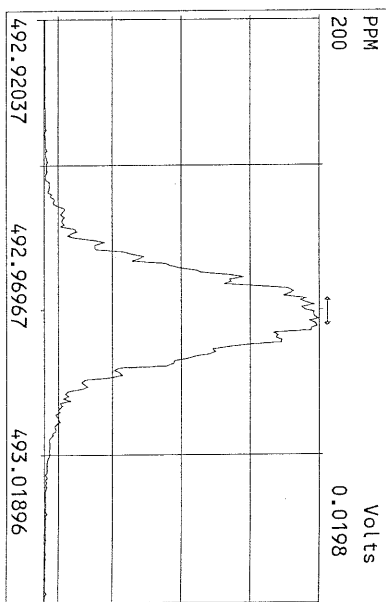
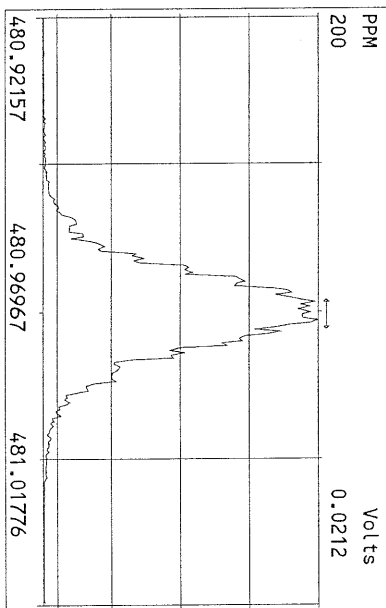
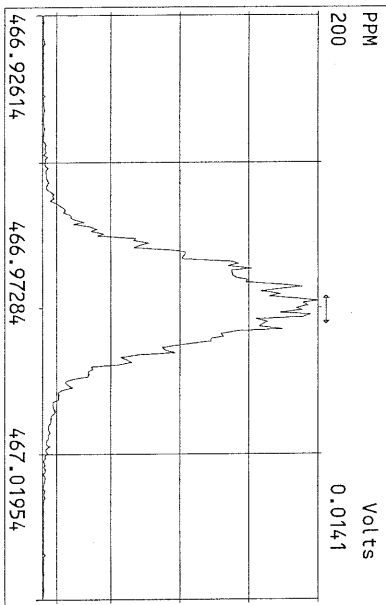
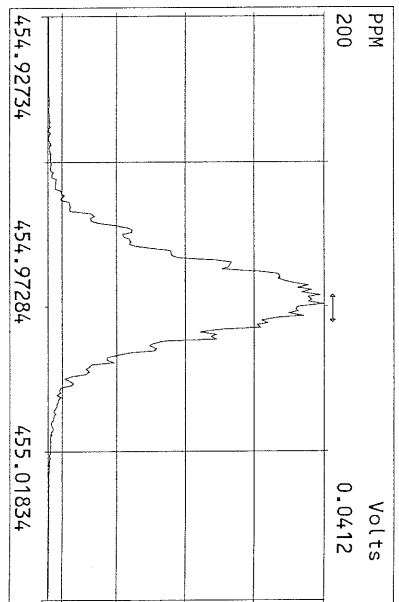
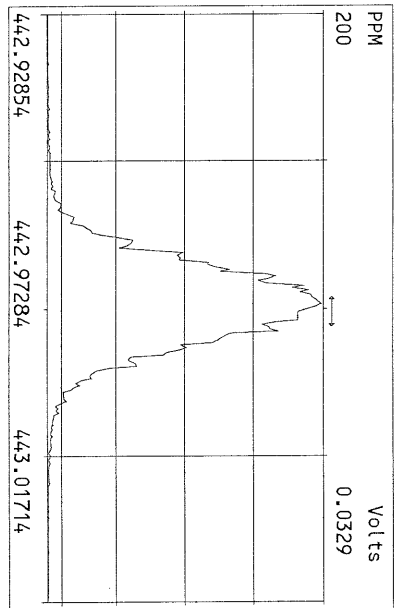
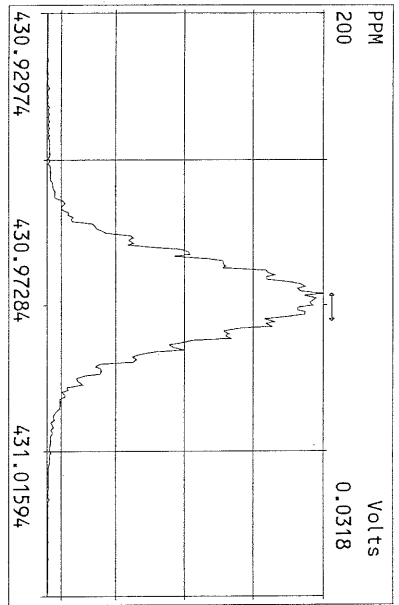






Peak Locate Examination: 8-APR-2011:08:50 File:07APR11M_RES_CHECK
Experiment:OCDD Function:4 Reference:PFK





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: 07APR11M Sam:19

Analysis Date: 8-APR-11 07:49:36

| 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 | 0.78 | 0.65-0.89 | y | 10.8 | 7.80 - 12.9 |
| 1,2,3,7,8-PeCDD | M+2/M+4 | 1.53 | 1.32-1.78 | y | 51.2 | 39.0 - 65.0 |
| 1,2,3,4,7,8-HxCDD | M+2/M+4 | 1.27 | 1.05-1.43 | y | 48.0 | 39.0 - 64.0 |
| 1,2,3,6,7,8-HxCDD | M+2/M+4 | 1.28 | 1.05-1.43 | y | 47.2 | 39.0 - 64.0 |
| 1,2,3,7,8,9-HxCDD | M+2/M+4 | 1.24 | 1.05-1.43 | y | 49.4 | 41.0 - 61.0 |
| 1,2,3,4,6,7,8-HpCDD | M+2/M+4 | 0.92 | 0.88-1.20 | y | 51.1 | 43.0 - 58.0 |
| OCDD | M+2/M+4 | 0.94 | 0.76-1.02 | y | 96.4 | 79.0 - 126 |
| 2,3,7,8-TCDF | M/M+2 | 0.72 | 0.65-0.89 | y | 10.2 | 8.40 - 12.0 |
| 1,2,3,7,8-PeCDF | M+2/M+4 | 1.61 | 1.32-1.78 | y | 52.2 | 41.0 - 60.0 |
| 2,3,4,7,8-PeCDF | M+2/M+4 | 1.63 | 1.32-1.78 | y | 50.3 | 41.0 - 60.0 |
| 1,2,3,4,7,8-HxCDF | M+2/M+4 | 1.22 | 1.05-1.43 | y | 49.1 | 45.0 - 56.0 |
| 1,2,3,6,7,8-HxCDF | M+2/M+4 | 1.23 | 1.05-1.43 | y | 49.2 | 44.0 - 57.0 |
| 2,3,4,6,7,8-HxCDF | M+2/M+4 | 1.23 | 1.05-1.43 | y | 48.6 | 44.0 - 57.0 |
| 1,2,3,7,8,9-HxCDF | M+2/M+4 | 1.25 | 1.05-1.43 | y | 49.8 | 45.0 - 56.0 |
| 1,2,3,4,6,7,8-HpCDF | M+2/M+4 | 1.08 | 0.88-1.20 | y | 49.4 | 45.0 - 55.0 |
| 1,2,3,4,7,8,9-HpCDF | M+2/M+4 | 1.10 | 0.88-1.20 | y | 50.5 | 43.0 - 58.0 |
| OCDF | M+2/M+4 | 0.92 | 0.76-1.02 | y | 99.3 | 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: 4/8/11

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: 07APR11M Sam:19

Analysis Date: 8-APR-11 07:49:36

| 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 | 0.73 | 0.65-0.89 | y | 98.4 | 82.0 - 121 |
| 13C-1,2,3,7,8-PeCDD | M+2/M+4 | 1.78 | 1.32-1.78 | y | 96.0 | 62.0 - 160 |
| 13C-1,2,3,4,7,8-HxCDD | M+2/M+4 | 1.31 | 1.05-1.43 | y | 86.6 | 85.0 - 117 |
| 13C-1,2,3,6,7,8-HxCDD | M+2/M+4 | 1.31 | 1.05-1.43 | y | 105 | 85.0 - 118 |
| 13C-1,2,3,4,6,7,8-HpCDD | M+2/M+4 | 1.03 | 0.88-1.20 | y | 106 | 72.0 - 138 |
| 13C-OCDD | M+2/M+4 | 0.96 | 0.76-1.02 | y | 212 | 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 | 1.69 | 1.32-1.78 | y | 107 | 76.0 - 130 |
| 13C-2,3,4,7,8-PeCDF | M+2/M+4 | 1.61 | 1.32-1.78 | y | 107 | 77.0 - 130 |
| 13C-1,2,3,4,7,8-HxCDF | M/M+2 | 0.48 | 0.43-0.59 | y | 89.2 | 76.0 - 131 |
| 13C-1,2,3,6,7,8-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 97.3 | 70.0 - 143 |
| 13C-2,3,4,6,7,8-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 96.3 | 73.0 - 137 |
| 13C-1,2,3,7,8,9-HxCDF | M/M+2 | 0.47 | 0.43-0.59 | y | 96.8 | 74.0 - 135 |
| 13C-1,2,3,4,6,7,8-HpCDF | M/M+2 | 0.49 | 0.37-0.51 | y | 101 | 78.0 - 129 |
| 13C-1,2,3,4,7,8,9-HpCDF | M/M+2 | 0.49 | 0.37-0.51 | y | 101 | 77.0 - 129 |
| 13C-OCDF | M+2/M+4 | 0.94 | 0.76-1.02 | y | 200 | 96.0 - 415 |
| CLEANUP STANDARD (4) | | | | | | |
| 37Cl-2,3,7,8-TCDD | | | | | 9.28 | 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: 07APR11M Sam:19 Analysis Date: 8-APR-11 Time: 07:49:36
DB-5 IS Data Filename: 07APR11M Sam:19 Analysis Date: 8-APR-11 Time: 07:49:36
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) | 28:47 |
| 1,2,4,7,9-PeCDD (F) | 30:29 | 1,3,4,6,8-PeCDF (F) | 28:38 |
| 1,2,3,8,9-PeCDD (L) | 34:03 | 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) | 39:28 | 1,2,3,7,8,9-HxCDF (L) | 40:03 |
| 1,2,3,4,6,7,9-HpCDD (F) | 43:05 | 1,2,3,4,6,7,8-HpCDF (F) | 42:34 |
| 1,2,3,4,6,7,8-HpCDD (L) | 44:28 | 1,2,3,4,7,8,9-HpCDF (L) | 45:24 |

(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: 4/8/11

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: 8-APR-11 07:49:36

CS3 or VER Data Filename: 07APR11M

Sam:19

| 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 | 1.000 | 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 | 1.023 | 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 | | 1.240 | 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: 

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: 8-APR-11 07:49:36

CS3 or VER Data Filename: 07APR11M


Sam:19

| NATIVE ANALYTES | RETENTION TIME | | RRT | RRT |
|-------------------------|-------------------------|--|-------|---------------|
| | REFERENCE | | | 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 | | 1.001 | 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 | | 1.000 | 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 | | 1.000 | 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 | | 1.001 | 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 | | | 0.955 | 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 | | | 1.151 | 1.057-1.154 |
| 13C-OCDD | | | 1.270 | 1.032-1.311 |
| 13C-OCDF | | | 1.280 | 1.000-1.311 |

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

Analyst: Date: 4/8/11

| Name | Resp | RA | RT | RRF | Conc | Qual | Fac Noise-1 | Noise-2 | DL | #Hom |
|--------------------------|----------|--------|-------|------|------|------|-------------|---------|----|---------|
| 2,3,7,8-TCDD | 3.01e+06 | 0.78 y | 27:36 | 1.13 | 10.8 | | 2.50 | - | * | |
| 1,2,3,7,8-PeCDD | 1.23e+07 | 1.53 y | 33:29 | 1.02 | 51.2 | | 2.50 | - | * | |
| 1,2,3,4,7,8-HxCDD | 1.29e+07 | 1.27 y | 38:52 | 1.45 | 48.0 | | 2.50 | - | * | |
| 1,2,3,6,7,8-HxCDD | 1.21e+07 | 1.28 y | 39:01 | 1.45 | 47.2 | | 2.50 | - | * | |
| 1,2,3,7,8,9-HxCDD | 1.31e+07 | 1.24 y | 39:28 | 1.47 | 49.4 | | 2.50 | - | * | |
| 1,2,3,4,6,7,8-HpCDD | 1.05e+07 | 0.92 y | 44:28 | 1.30 | 51.1 | | 2.50 | - | * | |
| OCDD | 1.62e+07 | 0.94 y | 50:06 | 1.45 | 96.4 | | 2.50 | - | * | |
| 2,3,7,8-TCDF | 4.86e+06 | 0.72 y | 26:51 | 1.15 | 10.2 | | 2.50 | - | * | |
| 1,2,3,7,8-PeCDF | 1.71e+07 | 1.61 y | 31:44 | 0.89 | 52.2 | | 2.50 | - | * | |
| 2,3,4,7,8-PeCDF | 1.61e+07 | 1.63 y | 33:04 | 0.89 | 50.3 | | 2.50 | - | * | |
| 1,2,3,4,7,8-HxCDF | 1.46e+07 | 1.22 y | 37:28 | 1.01 | 49.1 | | 2.50 | - | * | |
| 1,2,3,6,7,8-HxCDF | 1.76e+07 | 1.23 y | 37:40 | 0.89 | 49.2 | | 2.50 | - | * | |
| 2,3,4,6,7,8-HxCDF | 1.59e+07 | 1.23 y | 38:36 | 1.02 | 48.6 | | 2.50 | - | * | |
| 1,2,3,7,8,9-HxCDF | 1.89e+07 | 1.25 y | 40:03 | 1.10 | 49.8 | | 2.50 | - | * | |
| 1,2,3,4,6,7,8-HpCDF | 1.32e+07 | 1.08 y | 42:34 | 1.48 | 49.4 | | 2.50 | - | * | |
| 1,2,3,4,7,8,9-HpCDF | 1.00e+07 | 1.10 y | 45:24 | 1.43 | 50.5 | | 2.50 | - | * | |
| OCDF | 1.75e+07 | 0.92 y | 50:30 | 0.84 | 99.3 | | 2.50 | - | * | |
| | | | | | | | | | | Rec |
| 13C-2,3,7,8-TCDD | 2.47e+07 | 0.73 y | 27:35 | 1.03 | 98.4 | | | | | 98.4 |
| 13C-1,2,3,7,8-PeCDD | 2.37e+07 | 1.78 y | 33:28 | 1.01 | 96.0 | | | | | 96.0 |
| 13C-1,2,3,4,7,8-HxCDD | 1.85e+07 | 1.31 y | 38:50 | 1.19 | 86.6 | | | | | 86.6 |
| 13C-1,2,3,6,7,8-HxCDD | 1.76e+07 | 1.31 y | 39:00 | 0.94 | 105 | | | | | 105 |
| 13C-1,2,3,4,6,7,8-HpCDD | 1.57e+07 | 1.03 y | 44:27 | 0.83 | 106 | | | | | 106 |
| 13C-OCDD | 2.32e+07 | 0.96 y | 50:06 | 0.61 | 212 | | | | | 106 |
| 13C-2,3,7,8-TCDF | 4.15e+07 | 0.88 y | 26:50 | 0.98 | 101 | | | | | 101 |
| 13C-1,2,3,7,8-PeCDF | 3.71e+07 | 1.69 y | 31:43 | 0.83 | 107 | | | | | 107 |
| 13C-2,3,4,7,8-PeCDF | 3.59e+07 | 1.61 y | 33:02 | 0.80 | 107 | | | | | 107 |
| 13C-1,2,3,4,7,8-HxCDF | 2.95e+07 | 0.48 y | 37:26 | 1.84 | 89.2 | | | | | 89.2 |
| 13C-1,2,3,6,7,8-HxCDF | 4.01e+07 | 0.47 y | 37:39 | 2.29 | 97.3 | | | | | 97.3 |
| 13C-2,3,4,6,7,8-HxCDF | 3.22e+07 | 0.47 y | 38:35 | 1.86 | 96.3 | | | | | 96.3 |
| 13C-1,2,3,7,8,9-HxCDF | 3.45e+07 | 0.47 y | 40:02 | 1.98 | 96.8 | | | | | 96.8 |
| 13C-1,2,3,4,6,7,8-HpCDF | 1.80e+07 | 0.49 y | 42:33 | 0.99 | 101 | | | | | 101 |
| 13C-1,2,3,4,7,8,9-HpCDF | 1.39e+07 | 0.49 y | 45:23 | 0.77 | 101 | | | | | 101 |
| 13C-OCDF | 4.18e+07 | 0.94 y | 50:28 | 1.17 | 200 | | | | | 99.9 |
| 37Cl-2,3,7,8-TCDD | 1.65e+06 | | 27:36 | 0.73 | 9.28 | | | | | 92.8 |
| 13C-1,2,3,4-TCDD | 2.44e+07 | 0.73 y | 26:60 | - | 64.2 | | | | | |
| 13C-1,2,3,4-TCDF | 4.18e+07 | 0.88 y | 25:44 | - | 58.1 | | | | | |
| 13C-1,2,3,7,8,9-HxCDD | 1.80e+07 | 1.29 y | 39:26 | - | 72.4 | | | | | |
| Total Tetra-Dioxins | 1.50e+07 | | 24:35 | 1.13 | 53.6 | | 2.50 | - | * | 20 |
| Total Penta-Dioxins | 2.62e+07 | | 30:29 | 1.02 | 109 | | 2.50 | - | * | 5 |
| Total Hexa-Dioxins | 4.28e+07 | | 36:23 | 1.46 | 163 | | 2.50 | - | * | 9 |
| Total Hepta-Dioxins | 2.18e+07 | | 43:05 | 1.30 | 107 | | 2.50 | - | * | 16 |
| Total Tetra-Furans | 2.27e+07 | | 23:13 | 1.15 | 47.6 | | 2.50 | - | * | 21 |
| 1st Fn. Tot Penta-Furans | 1.78e+07 | | 28:38 | 0.89 | 54.8 | | 2.50 | - | * | PeCDF 2 |
| Total Penta-Furans | 4.76e+07 | | 30:25 | 0.89 | 146 | | 2.50 | - | * | 201 12 |
| Total Hexa-Furans | 7.77e+07 | | 35:30 | 1.00 | 228 | | 2.50 | - | * | 19 |
| Total Hepta-Furans | 2.34e+07 | | 42:34 | 1.46 | 101 | | 2.50 | - | * | 13 |

Analyst: 

Date: 4/8/11

Frontier Analytical Laboratory - Acquisition Log

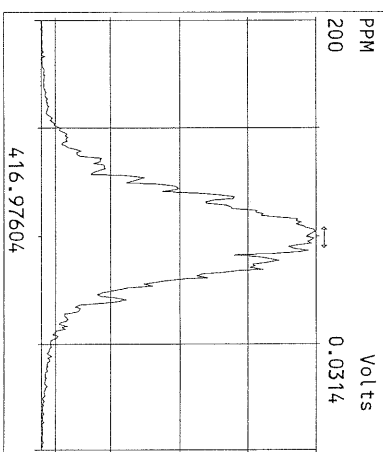
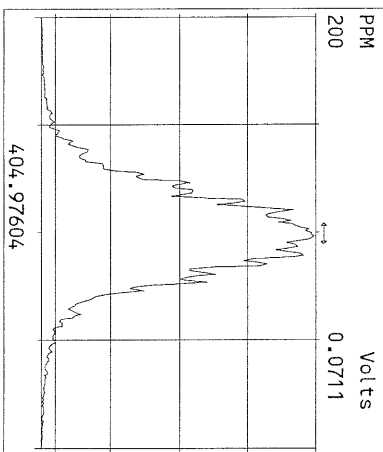
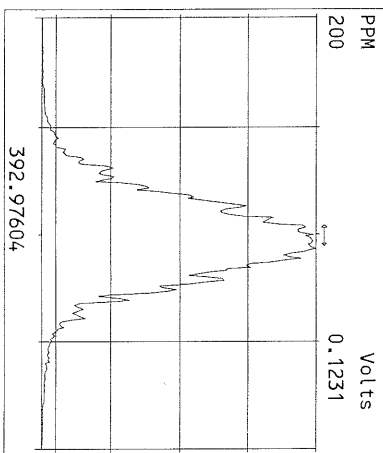
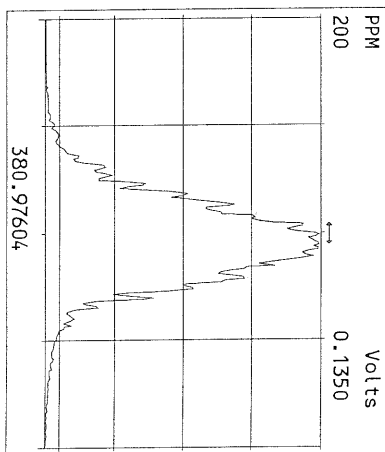
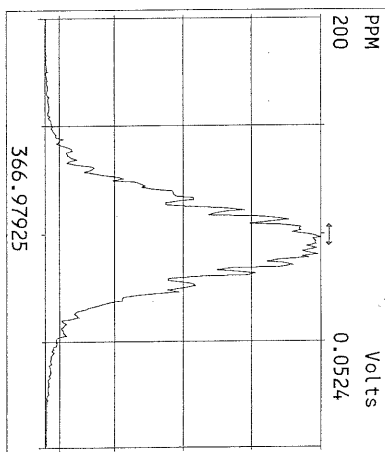
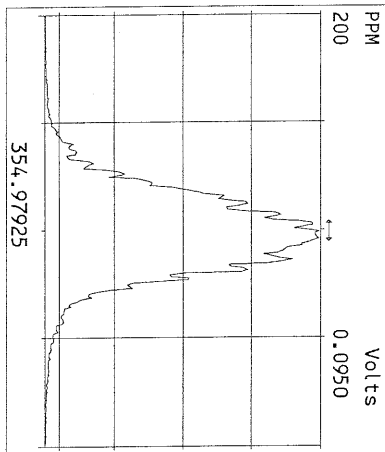
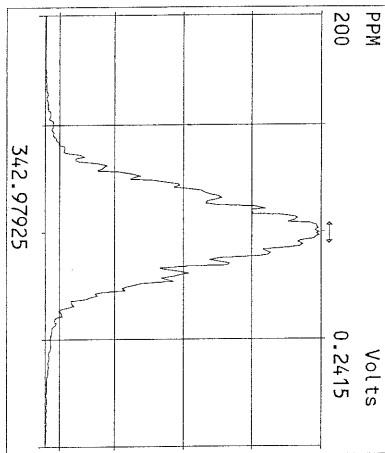
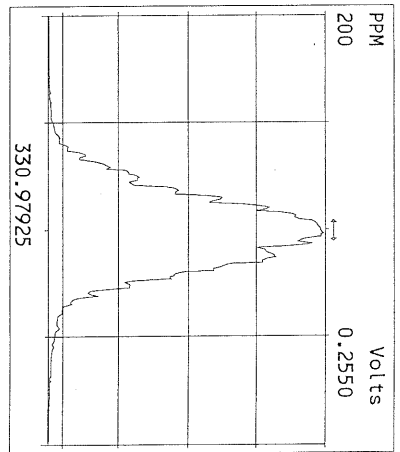
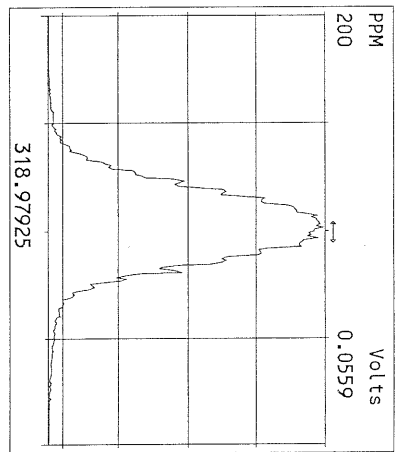
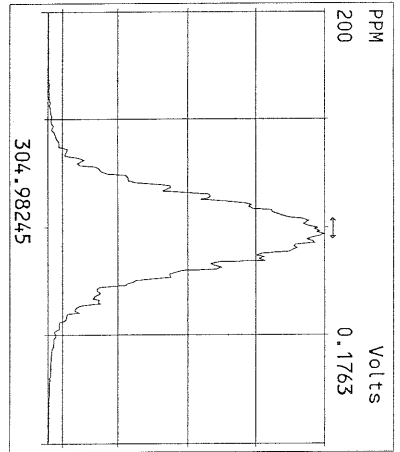
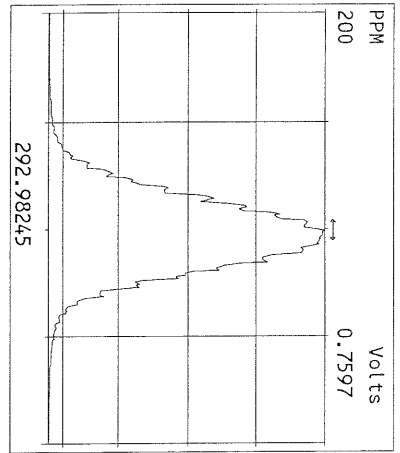
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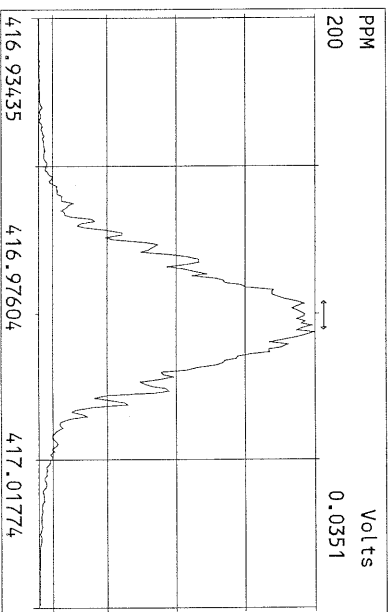
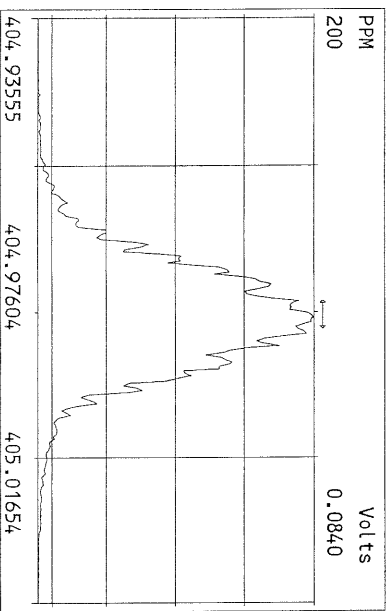
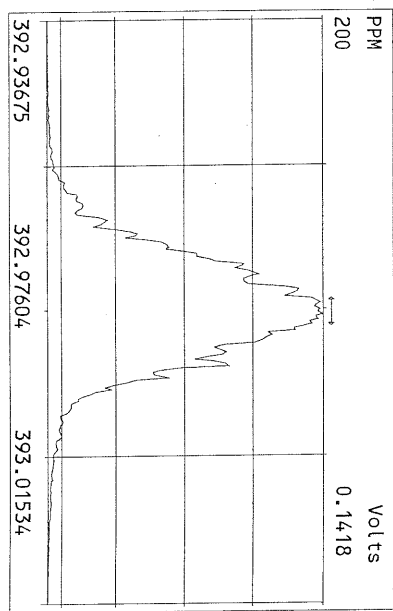
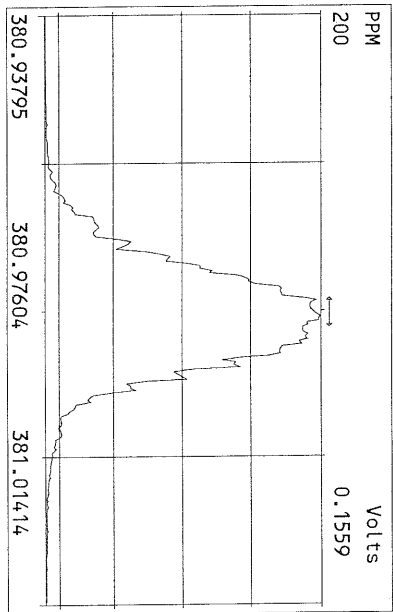
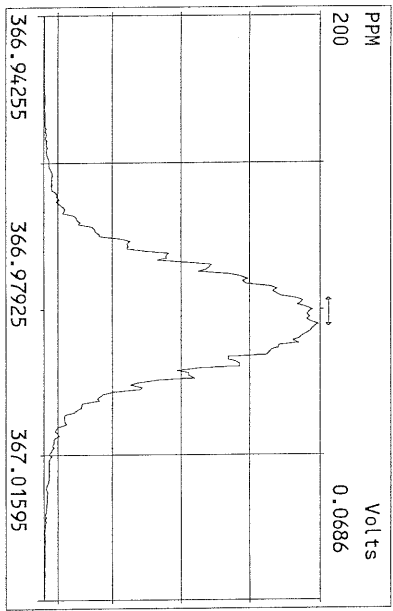
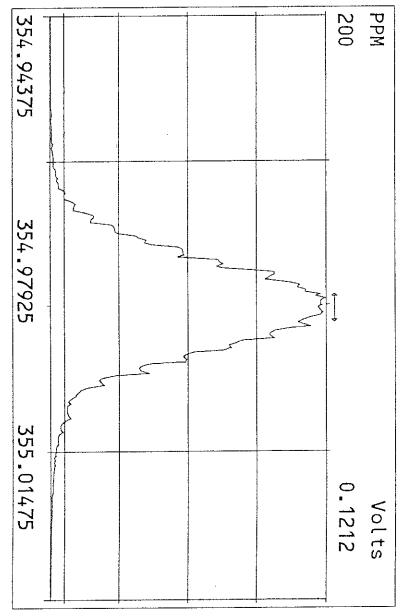
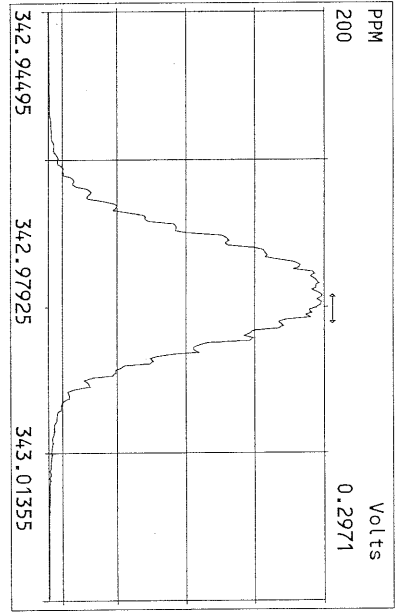
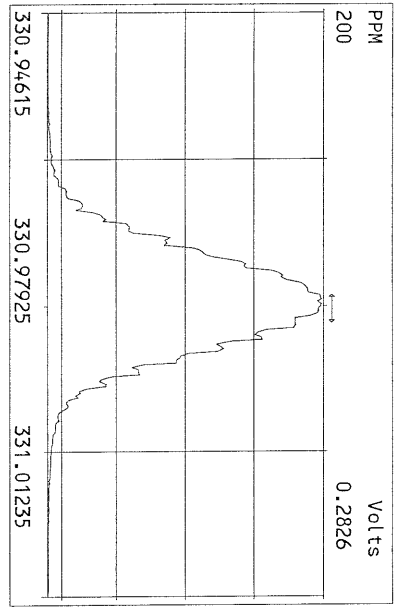
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| 07APR11M 3 | 2264-001-0001-OPR | OPR | 7-APR-11 17:03:57 | ST040711M1 | ST040711M2 | TC |
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| 07APR11M 5 | 6678-007-0001-SA | LL-SED2-0-56-031511-D | 7-APR-11 18:54:38 | ST040711M1 | ST040711M2 | TC |
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| 07APR11M 7 | 6678-005-0001-SA | LL-SED2-56-112-031511 | 7-APR-11 20:45:17 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 8 | 6678-004-0001-SA | LL-SED2-0-56-031511 | 7-APR-11 21:40:40 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 9 | 6678-003-0001-SA | LL-SED3-141-167-031511 | 7-APR-11 22:36:00 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 10 | 6678-002-0001-SA | LL-SED3-36-141-031511 | 7-APR-11 23:31:19 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 11 | 6678-001-0001-SA | LL-SED3-0-36-031511 | 8-APR-11 00:26:42 | ST040711M1 | ST040711M2 | TC |
| 07APR11M 12 | 6678-005-0002-MS | LL-SED2-56-112-031511 | 8-APR-11 01:22:09 | ST040711M1 | ST040711M2 | TC |
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| 07APR11M 14 | SB040711M1 | Solvent Blank | 8-APR-11 03:12:55 | ST040711M1 | ST040711M2 | TC |
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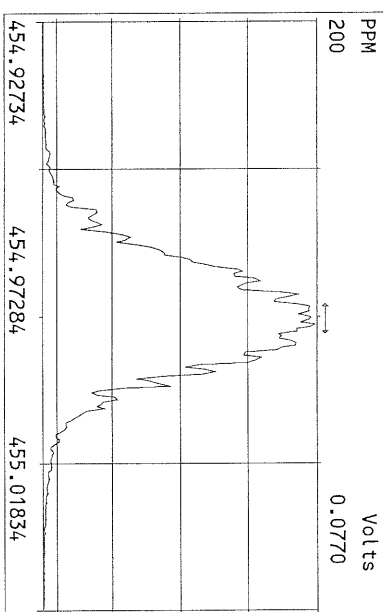
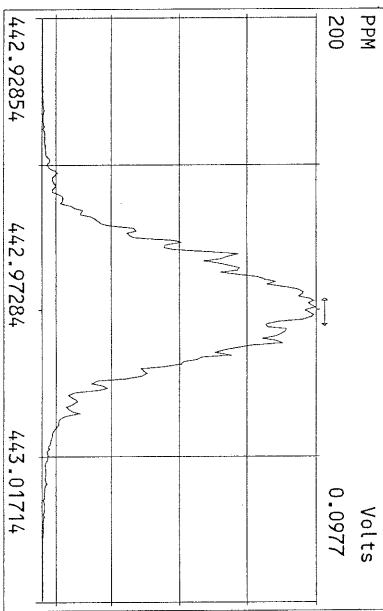
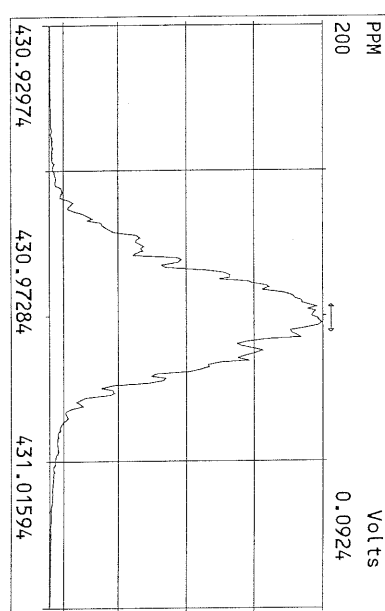
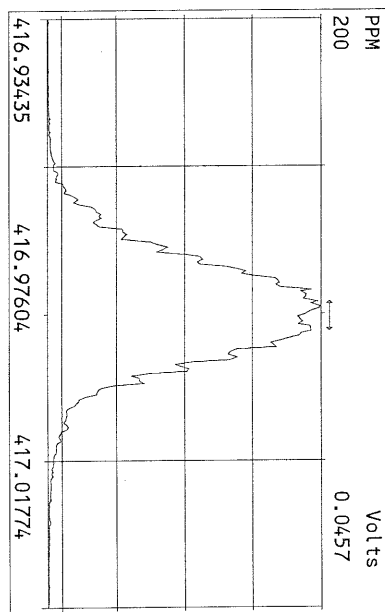
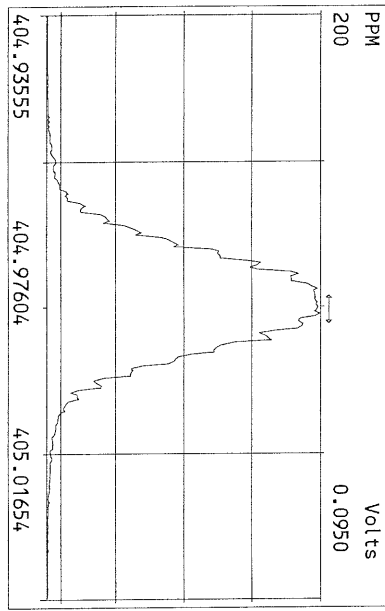
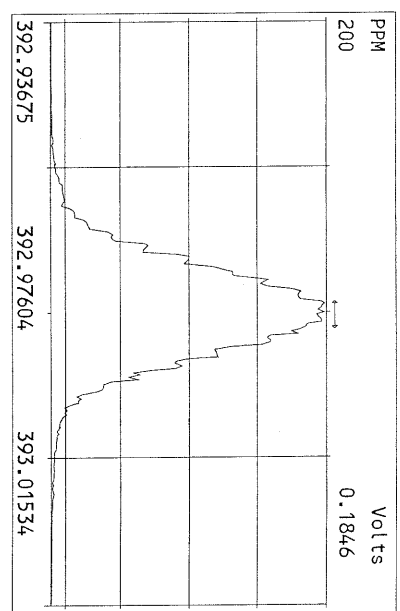
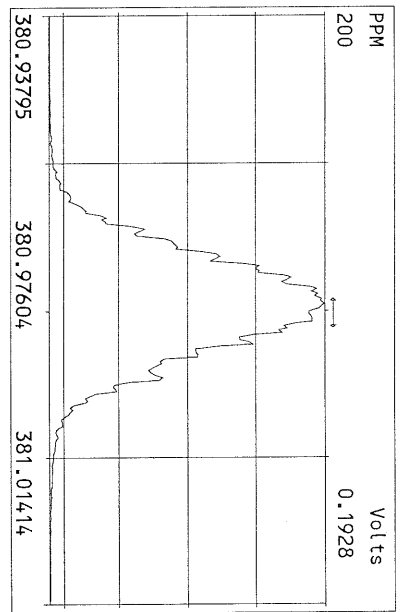
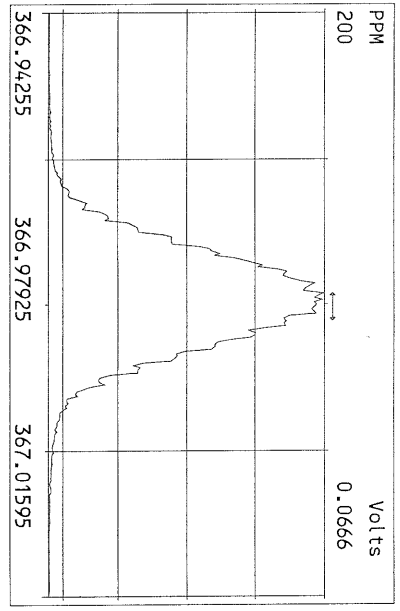


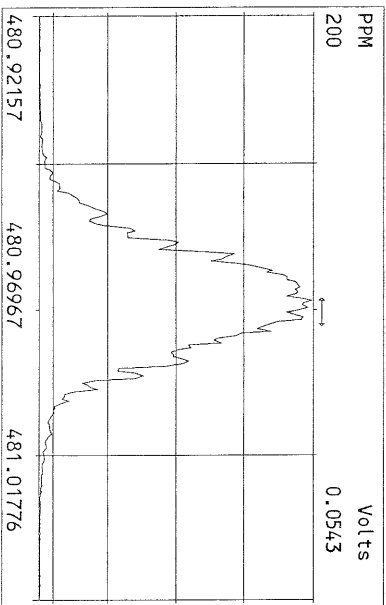
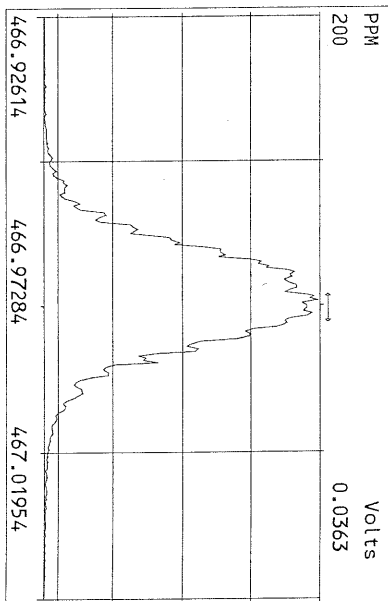
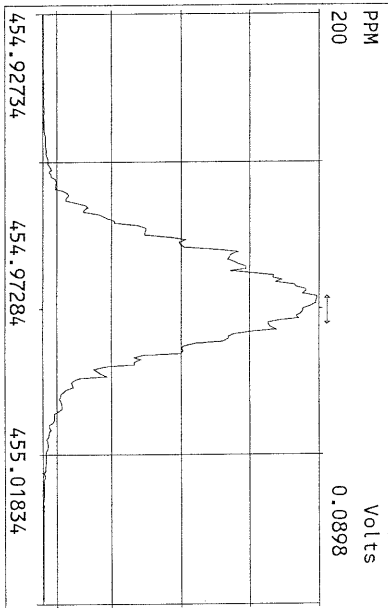
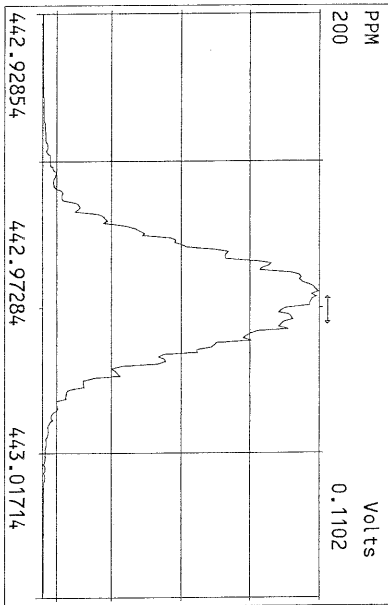
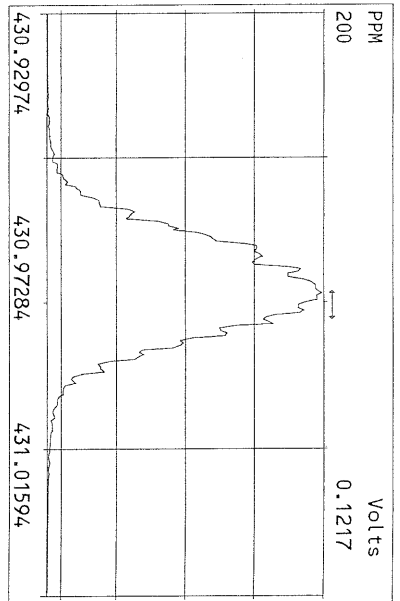
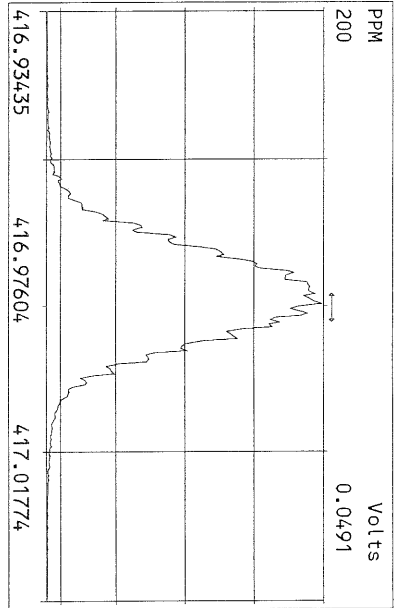
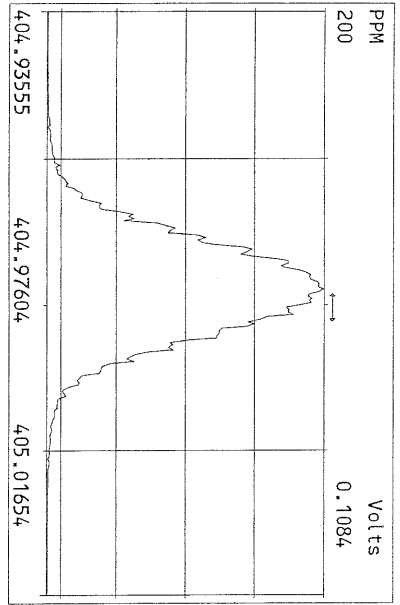

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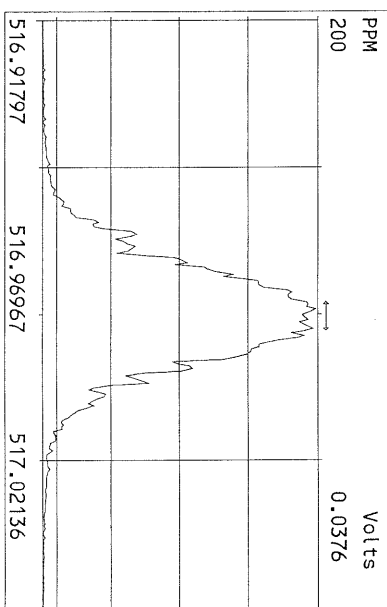
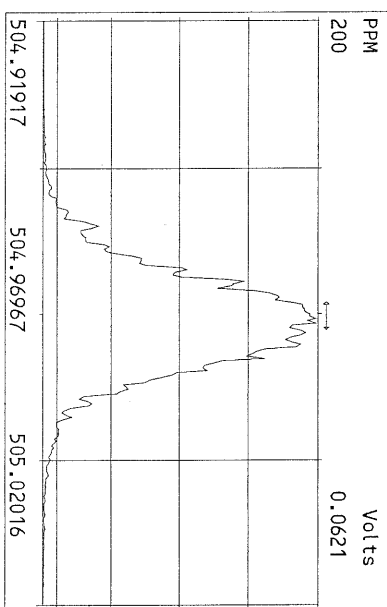
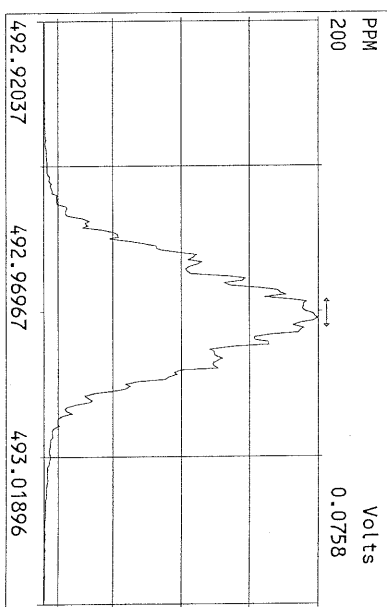
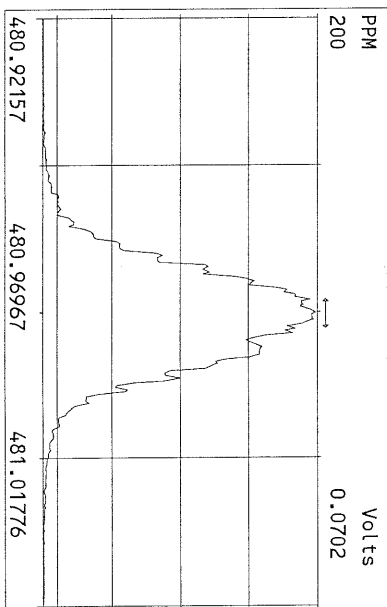
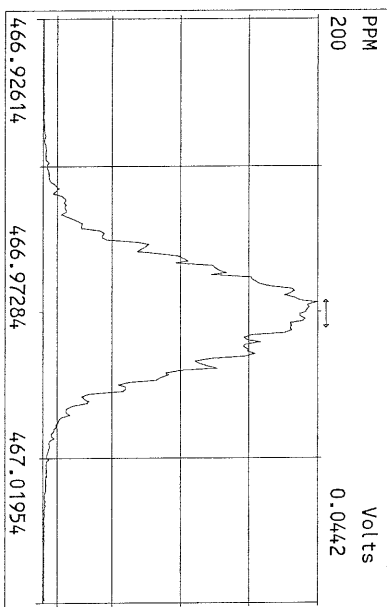
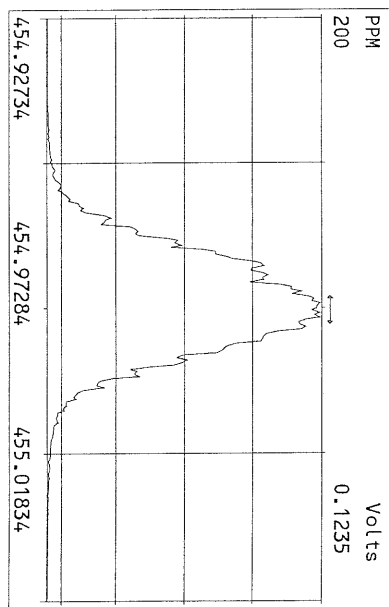
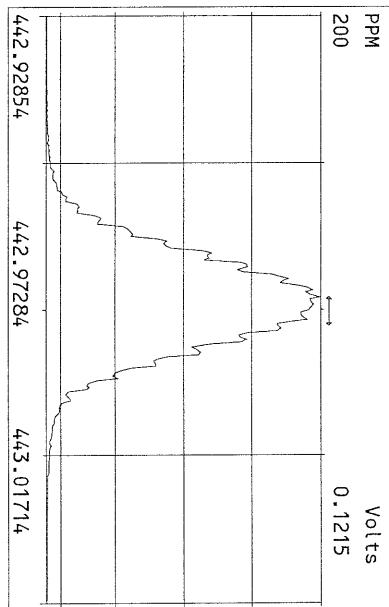
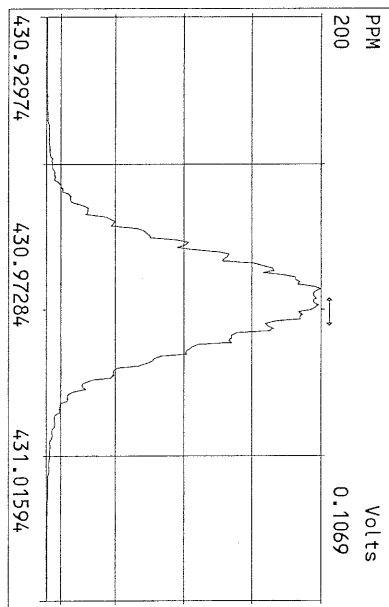
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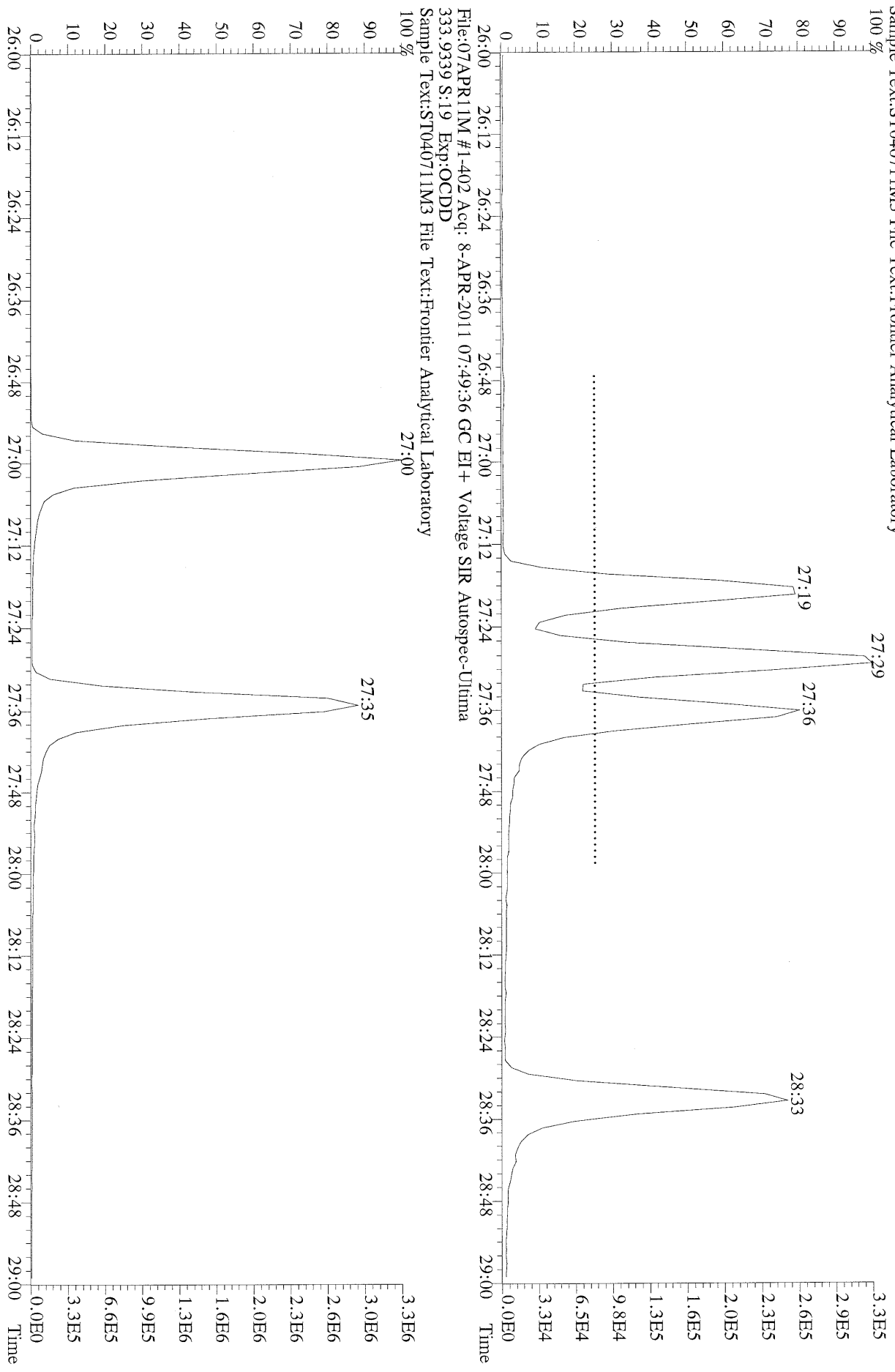




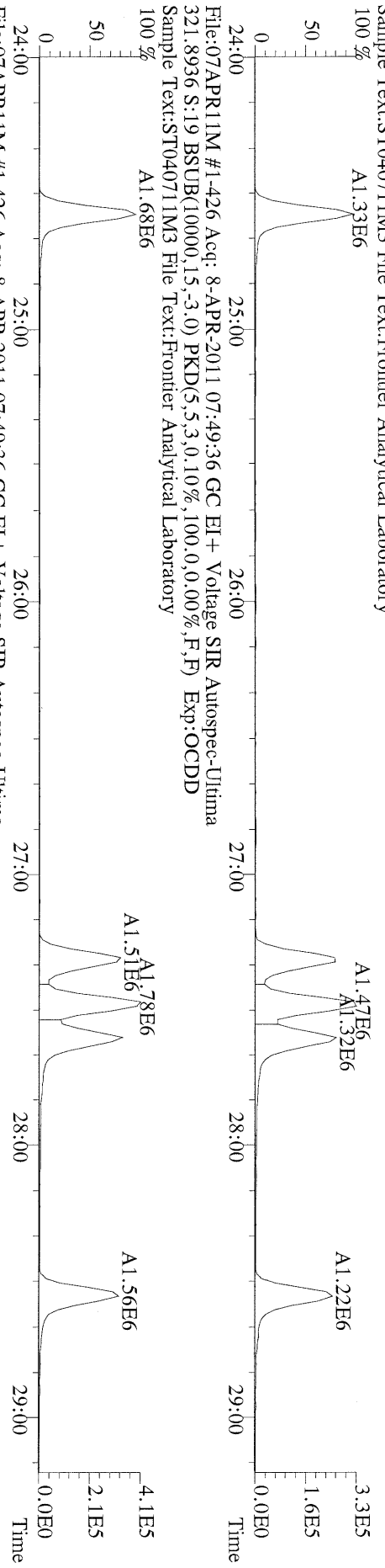




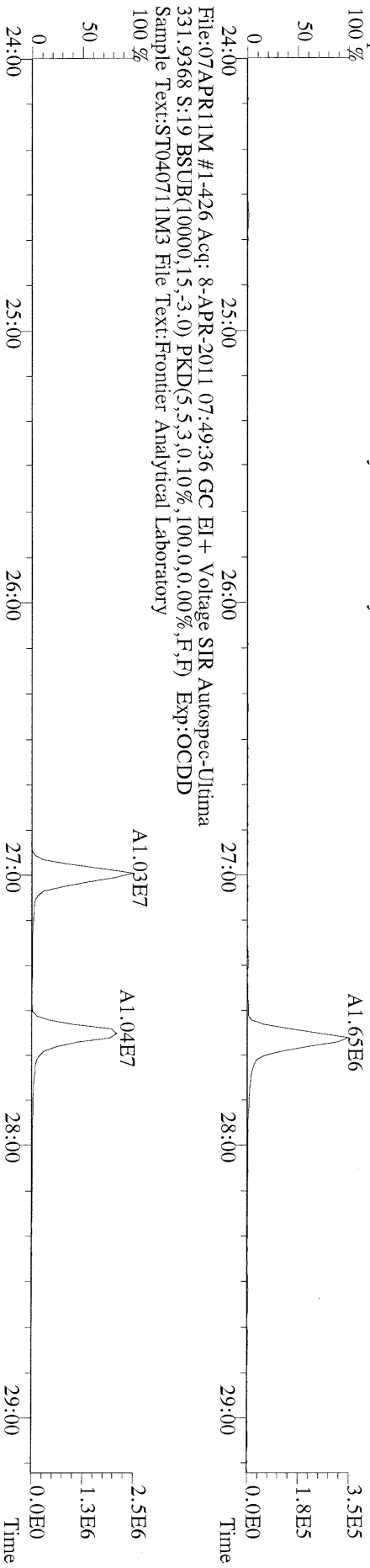
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Sample Text:ST040711M3 File Text:Frontier Analytical Laboratory



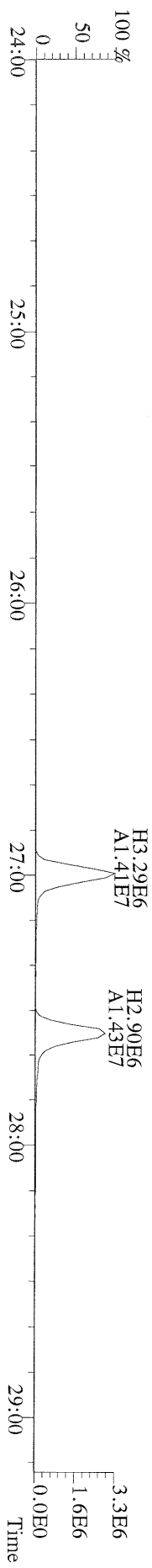
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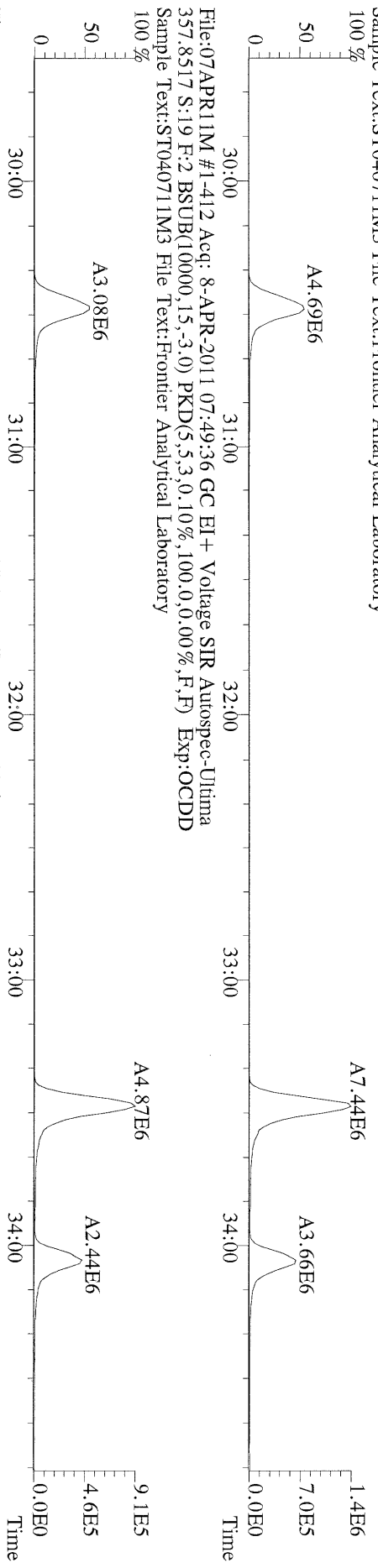
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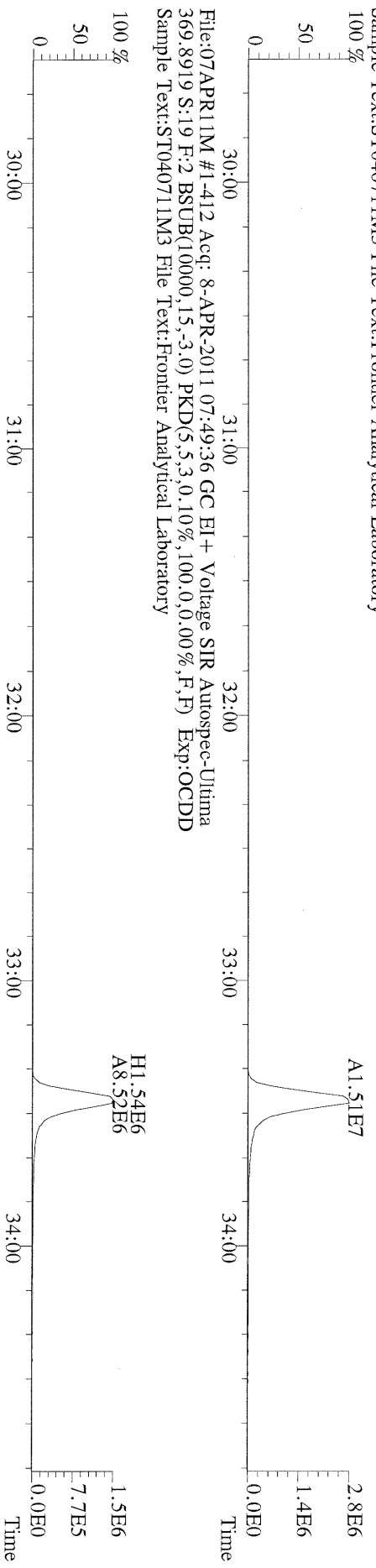
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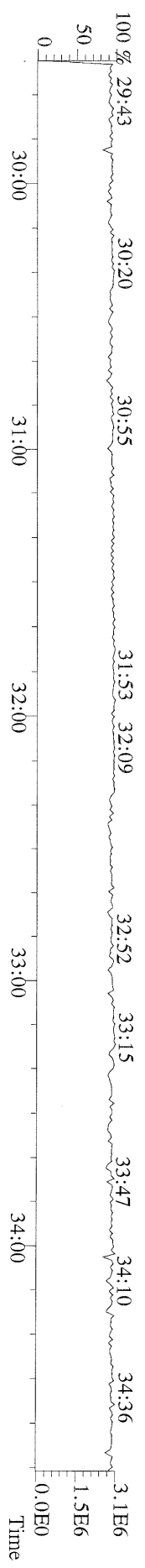
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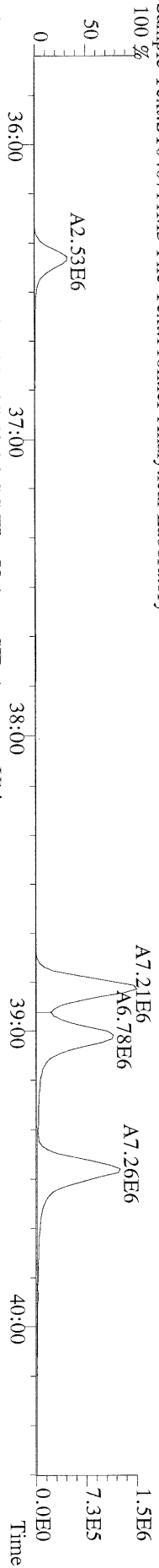
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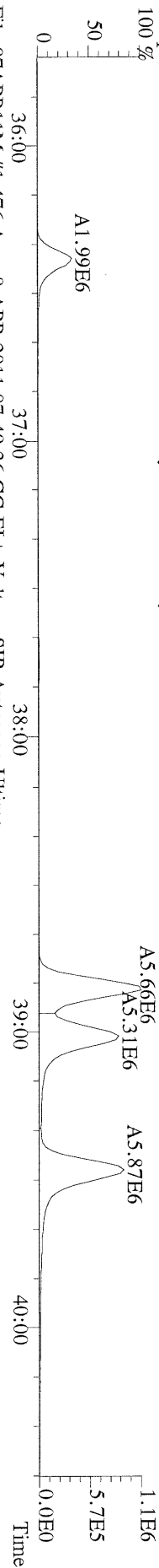
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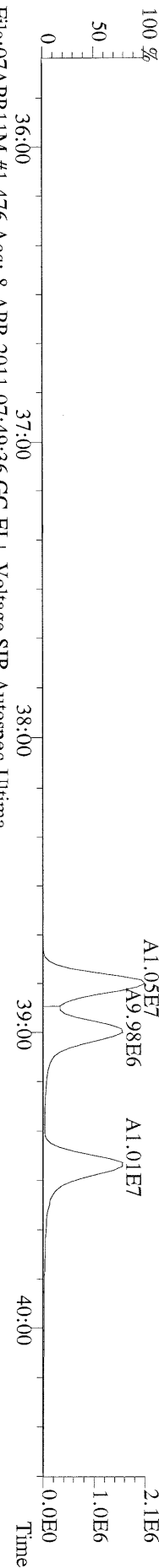
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Sample Text:ST04071IM3 File Text:Frontier Analytical Laboratory



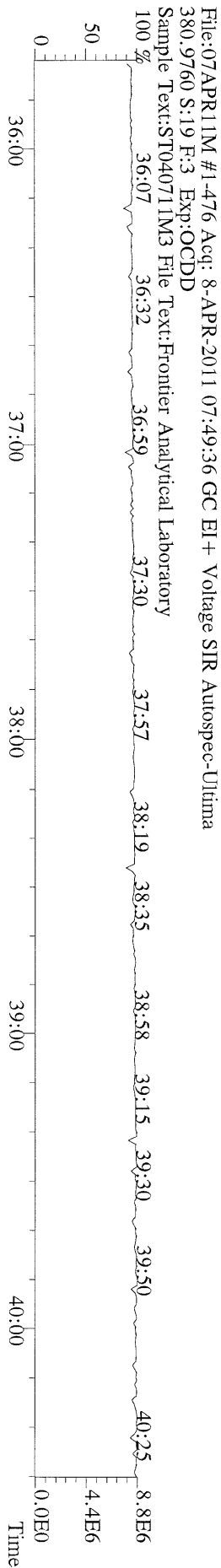
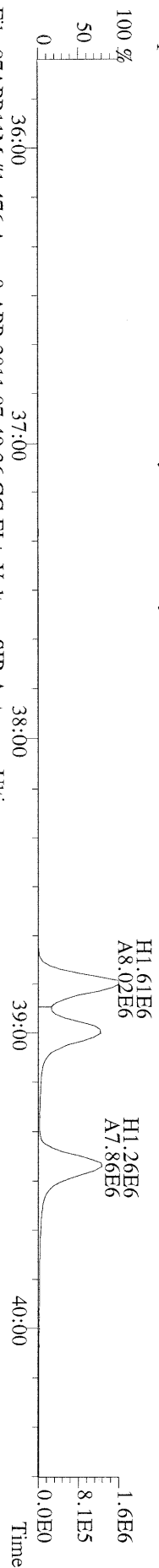
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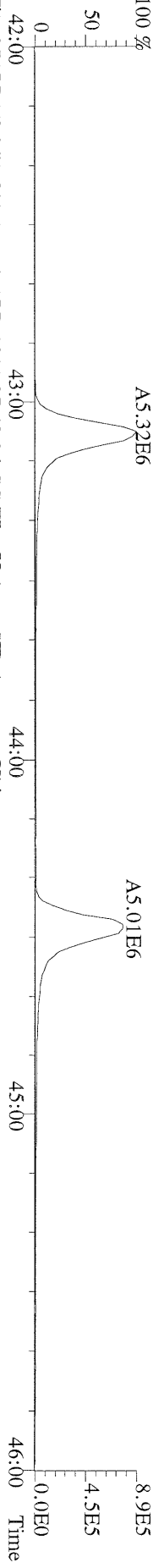
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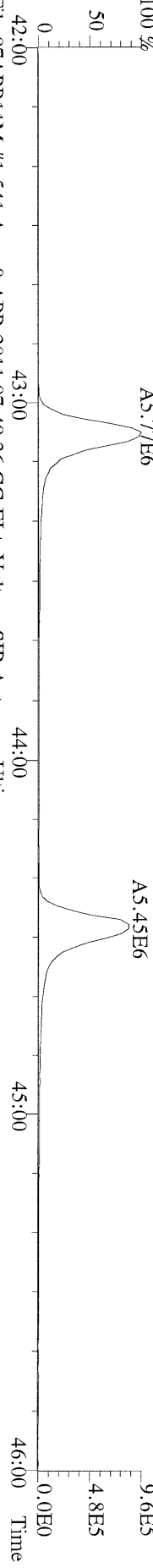
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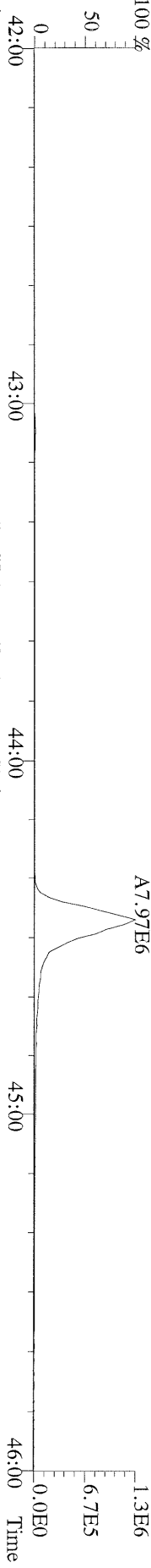
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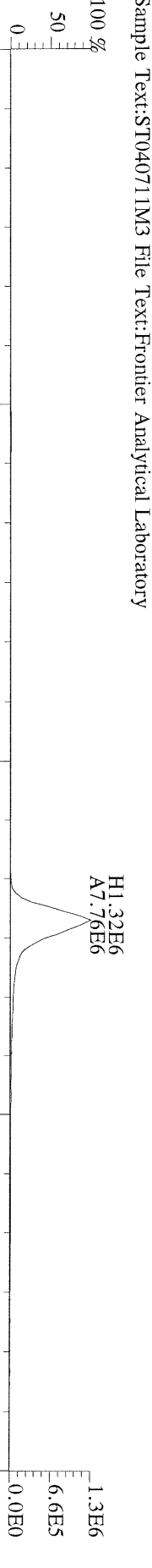
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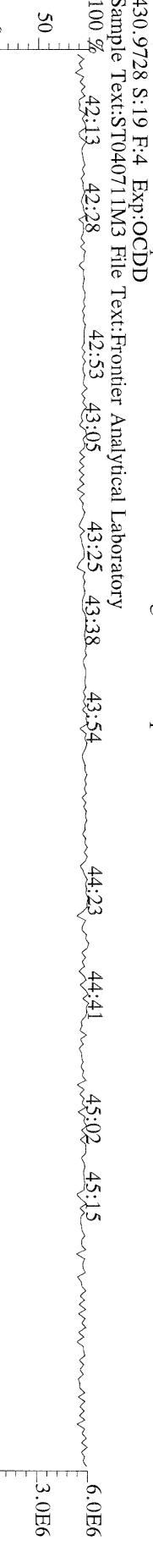
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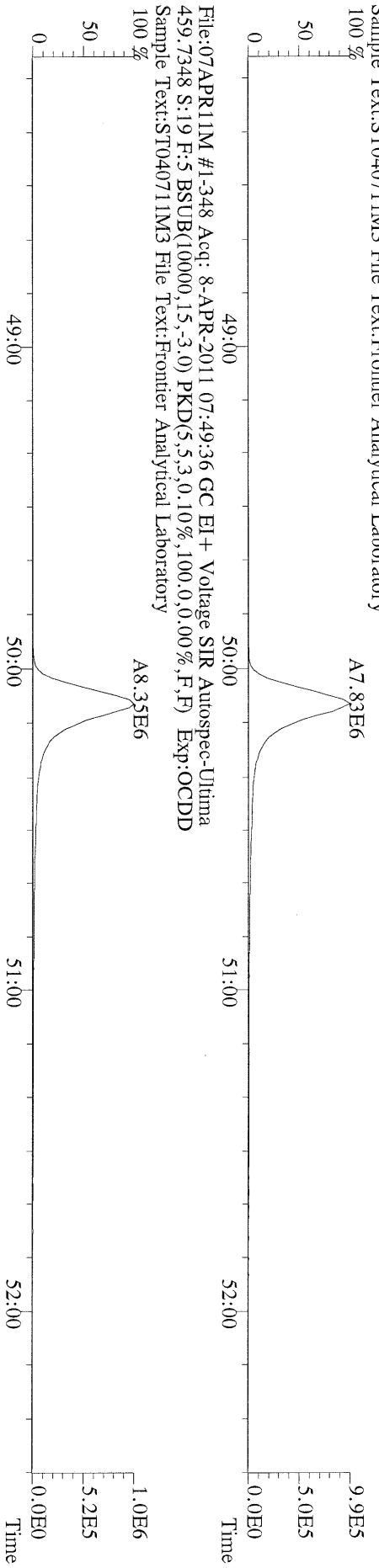
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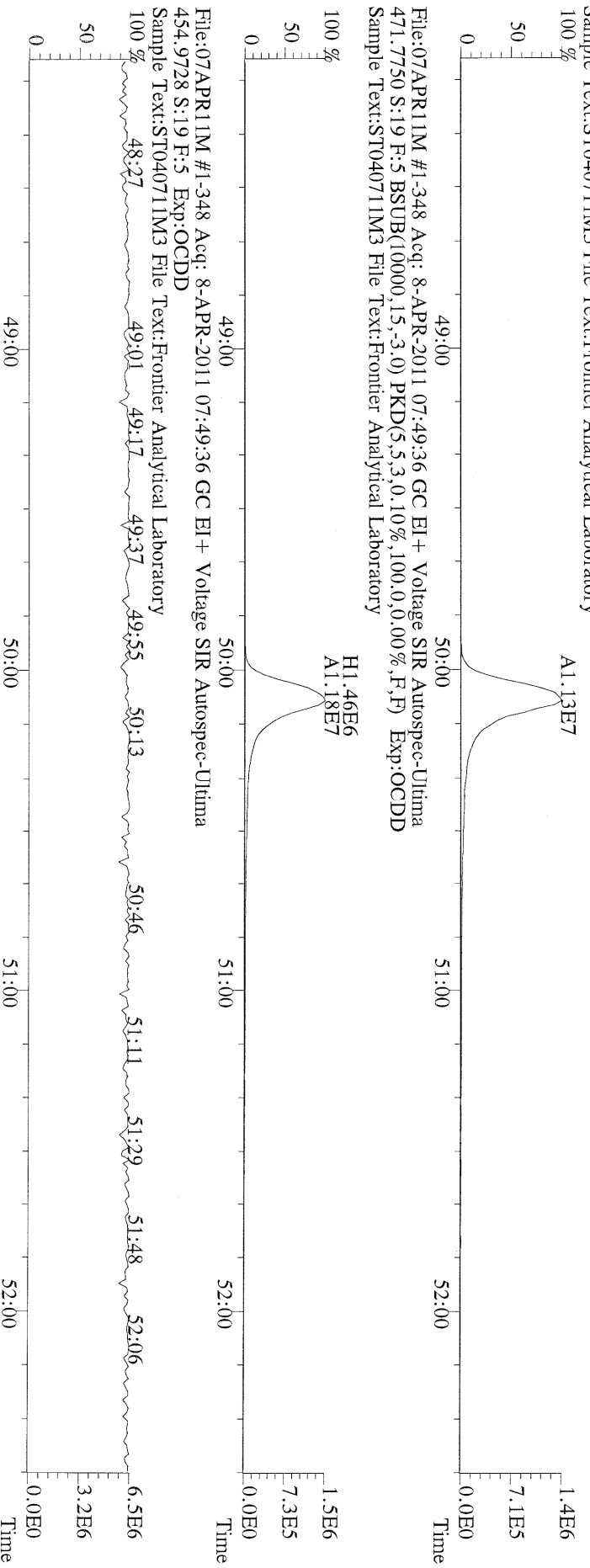
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Sample Text:ST04071IM3 File Text:Frontier Analytical Laboratory



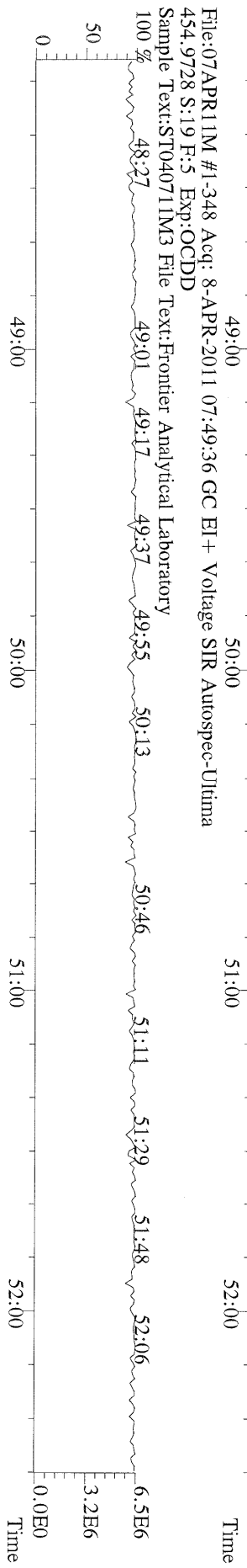
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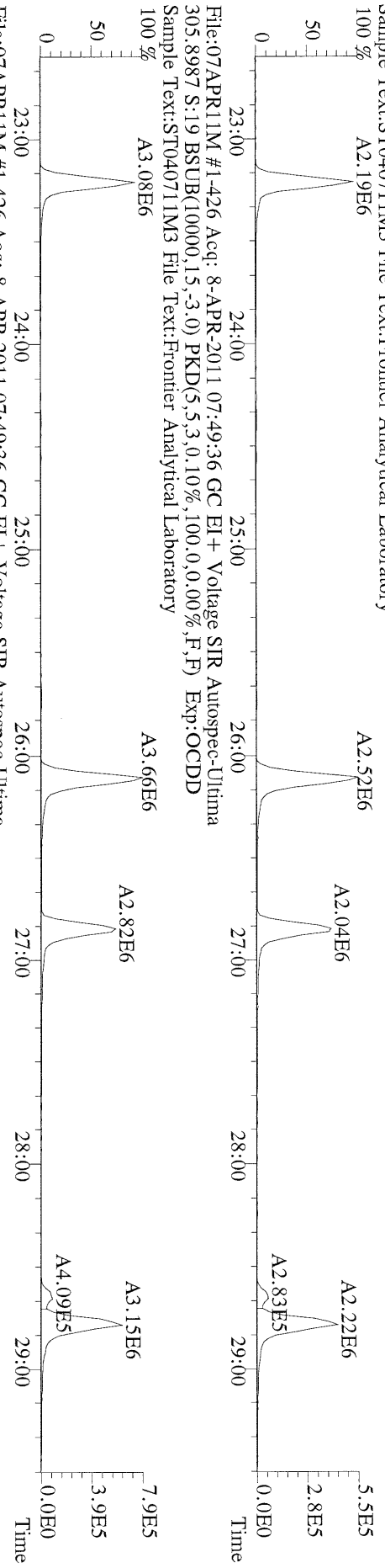
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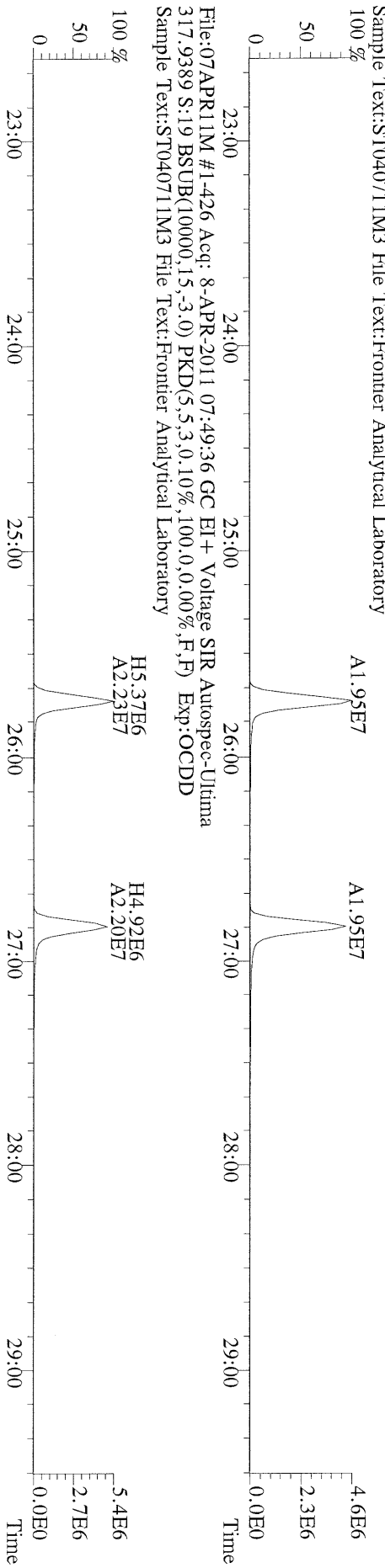
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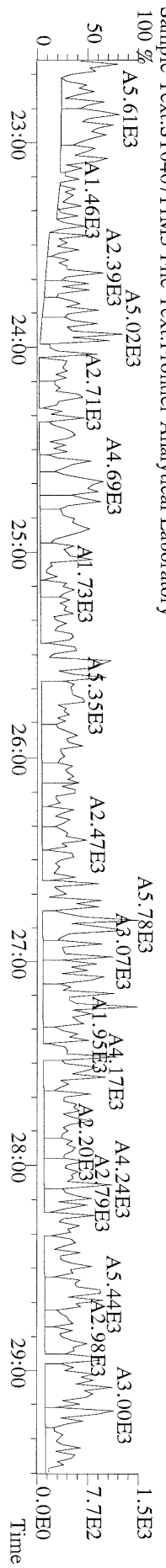
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 303.9016 S:19 BSUB(10000,15,-3.0) PKD(5,5,3,0.10%,100,0.0,0.00%,F,F) Exp:OCDD
 Sample Text:ST040711M3 File Text:Frontier Analytical Laboratory



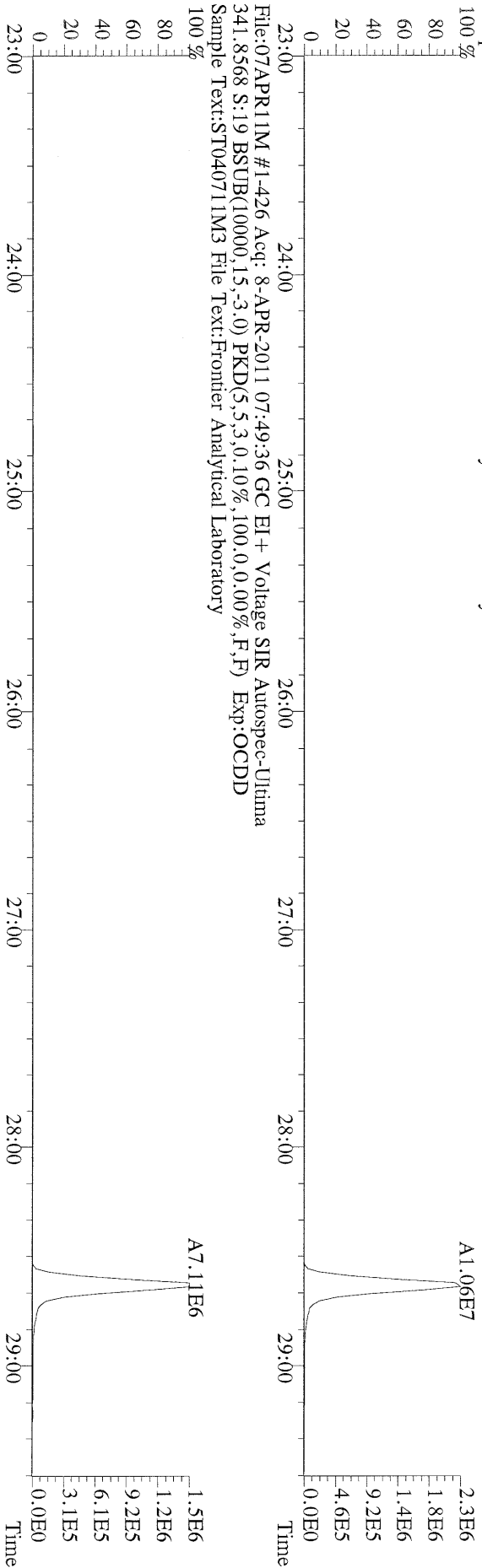
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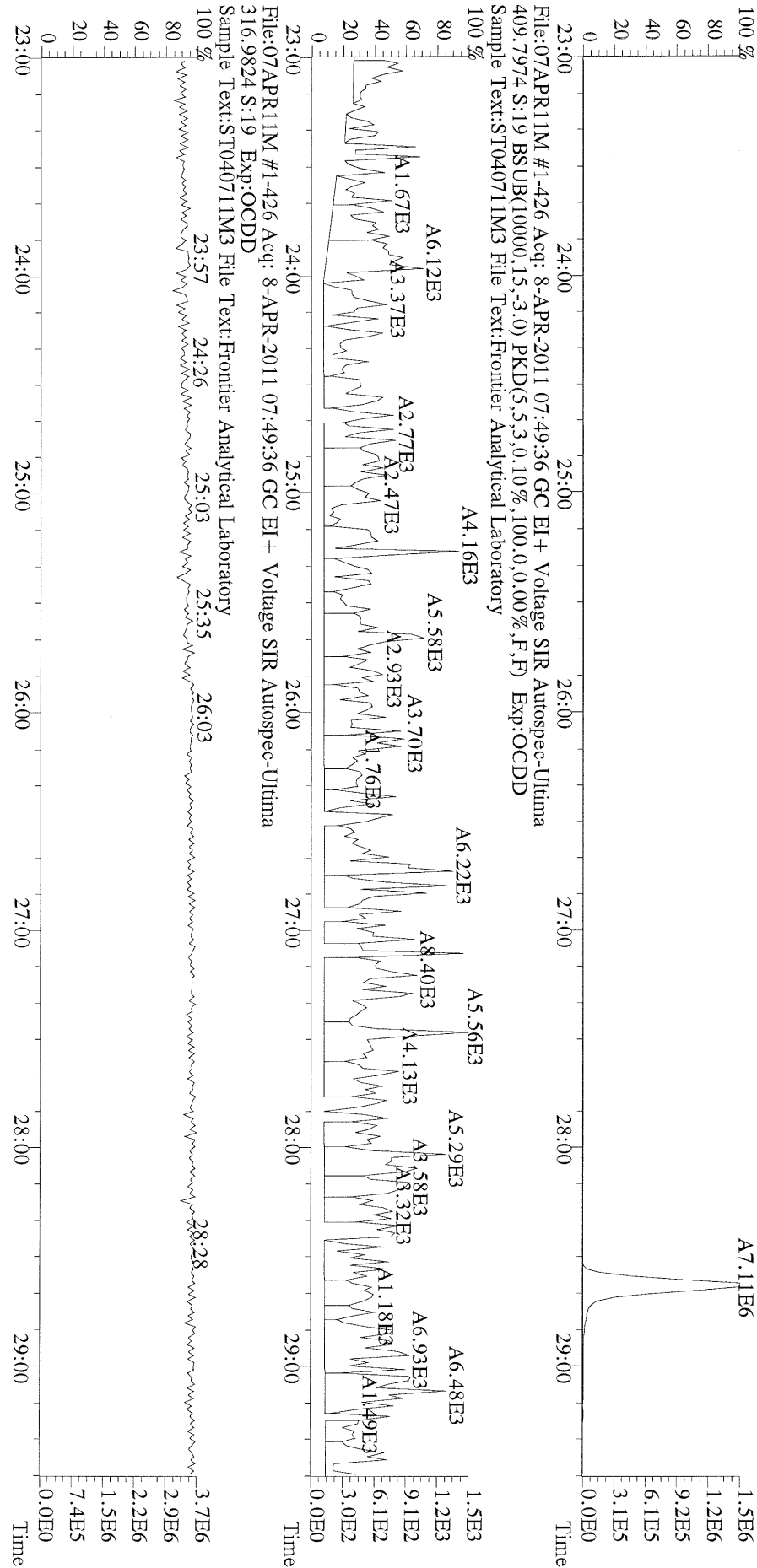
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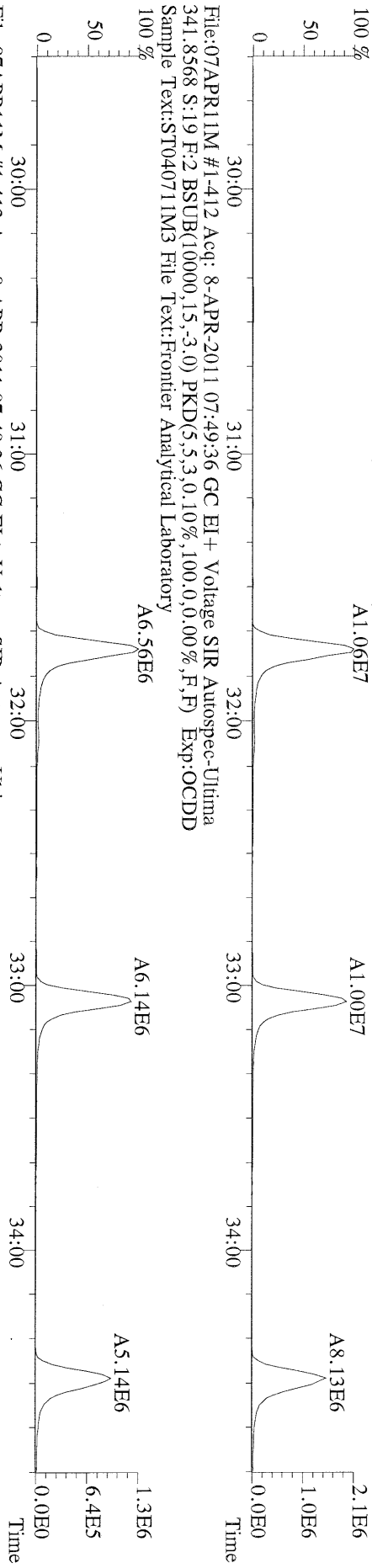
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 Sample Text:ST04071IM3 File Text:Frontier Analytical Laboratory



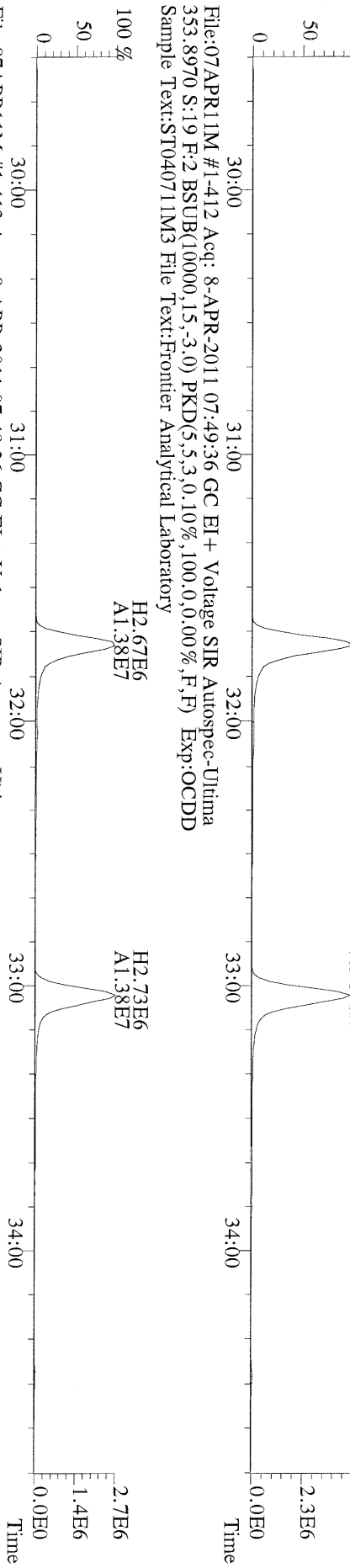
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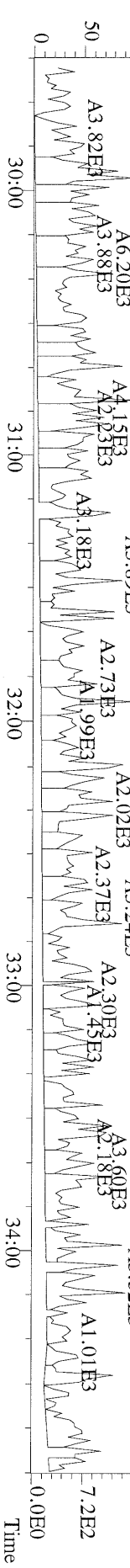
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Sample Text:ST040711M3 File Text:Frontier Analytical Laboratory



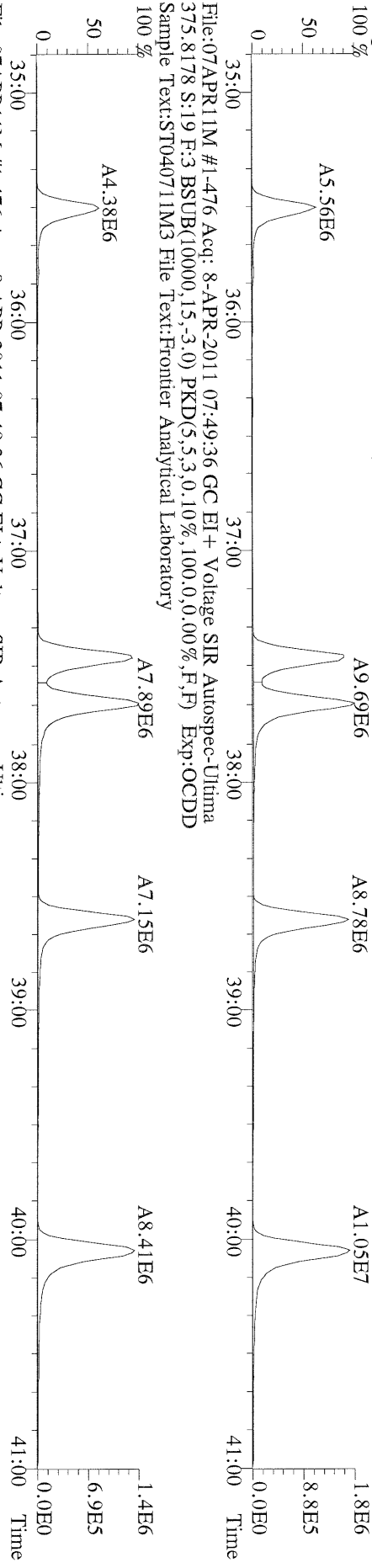
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Sample Text:ST040711M3 File Text:Frontier Analytical Laboratory



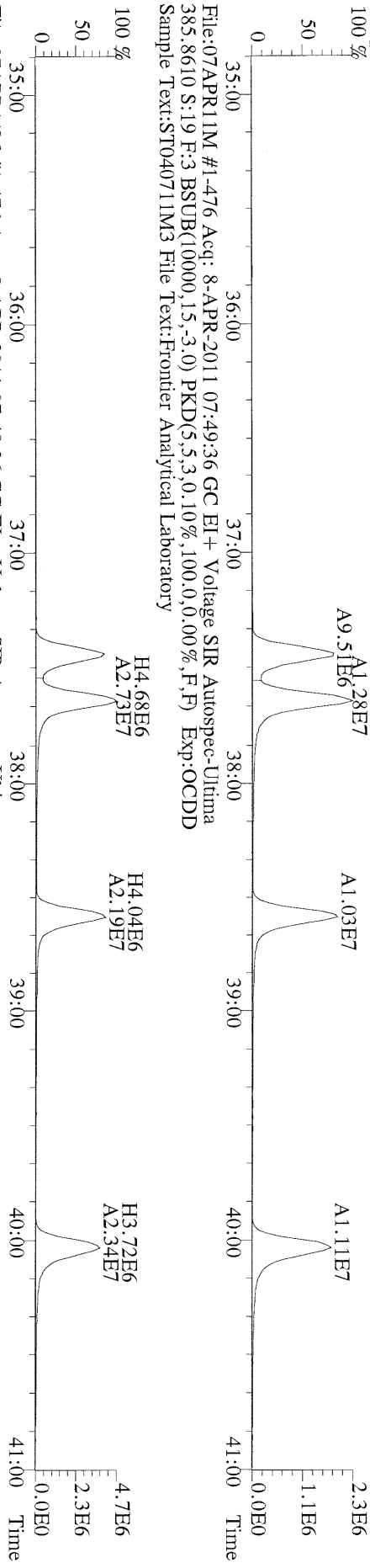
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Sample Text:ST040711M3 File Text:Frontier Analytical Laboratory



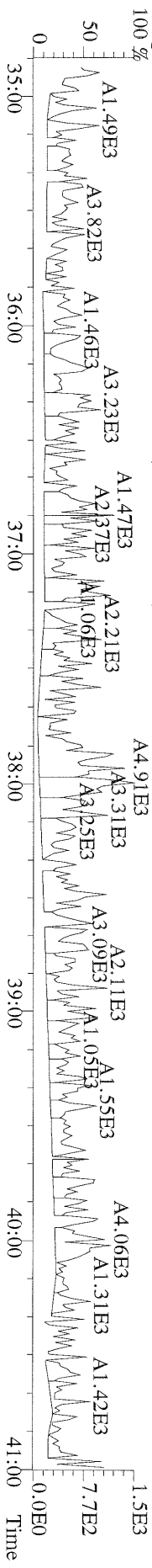
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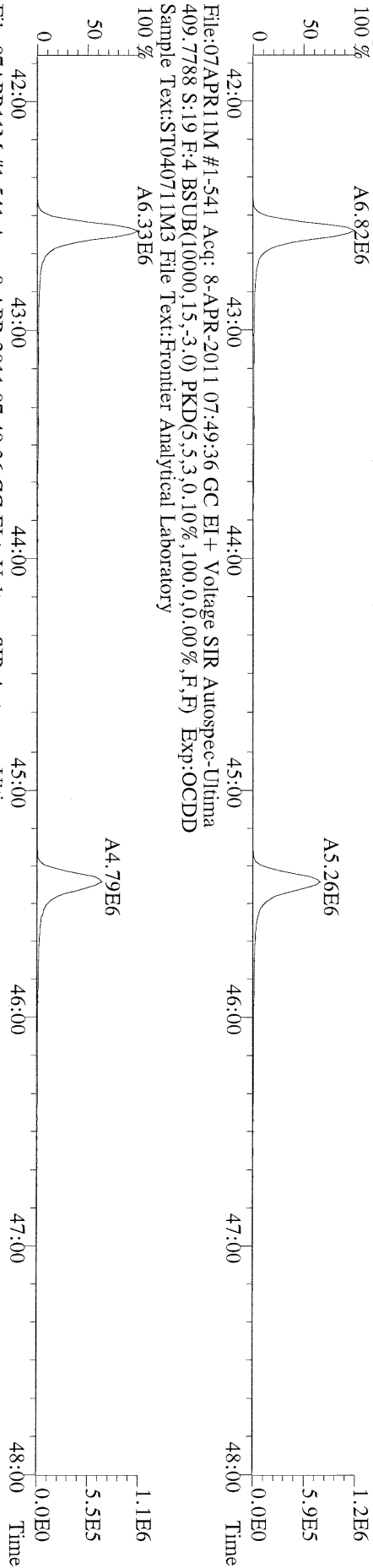
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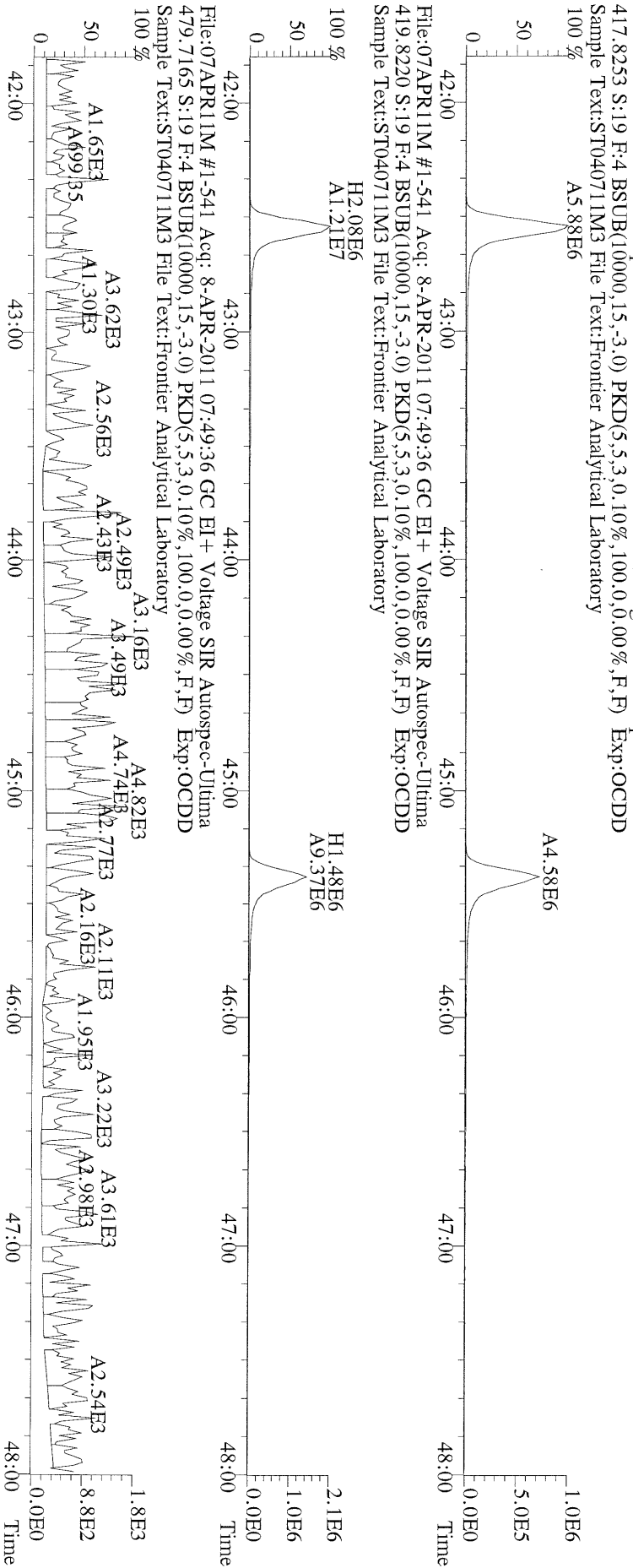
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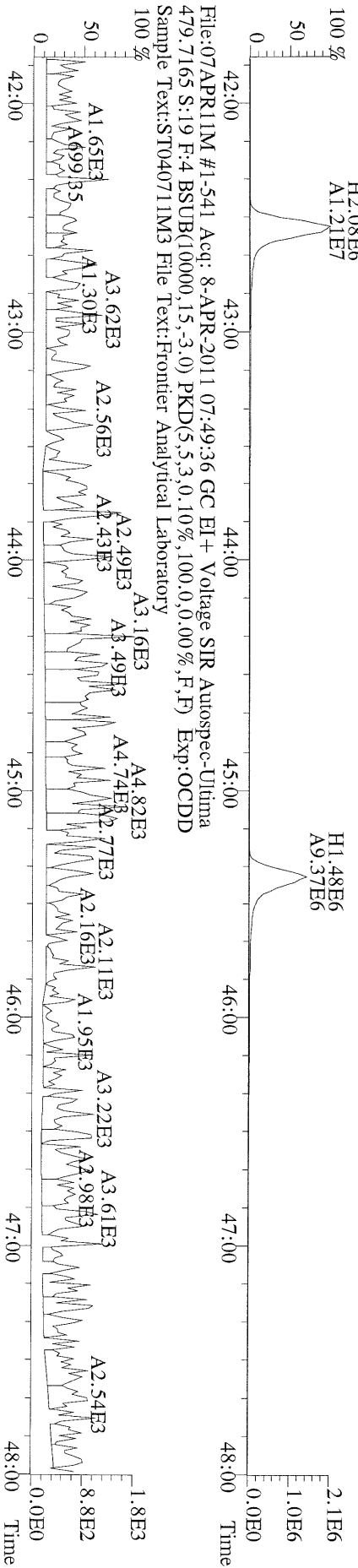
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Sample Text:ST040711M3 File Text:FronTier Analytical Laboratory
100 %



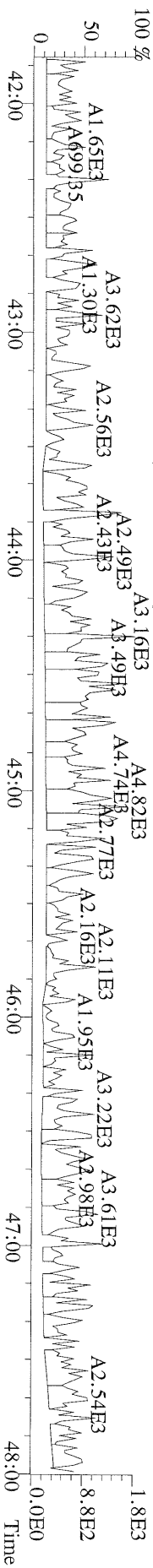
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Sample Text:ST040711M3 File Text:FronTier Analytical Laboratory
100 %



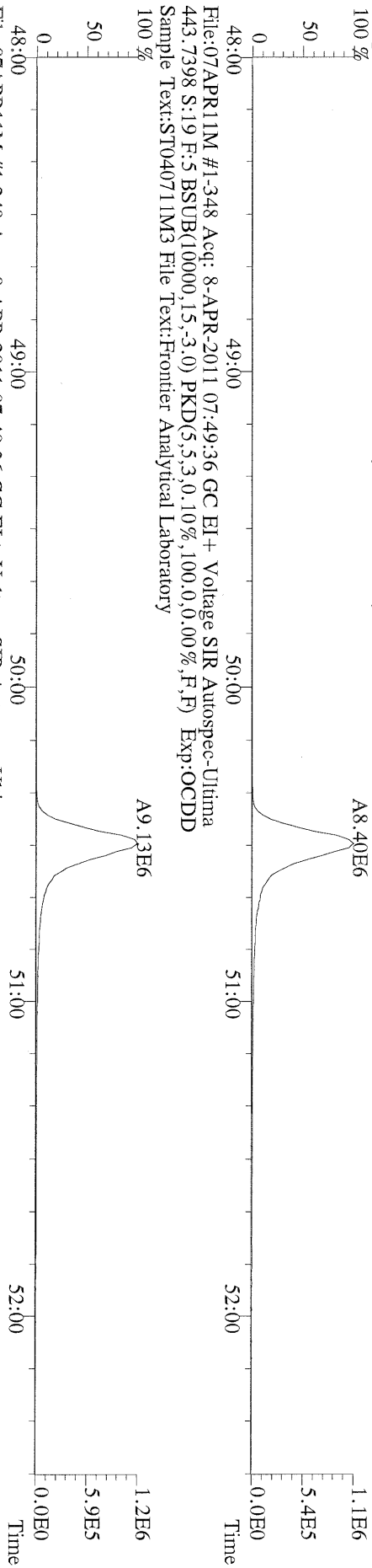
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100 %



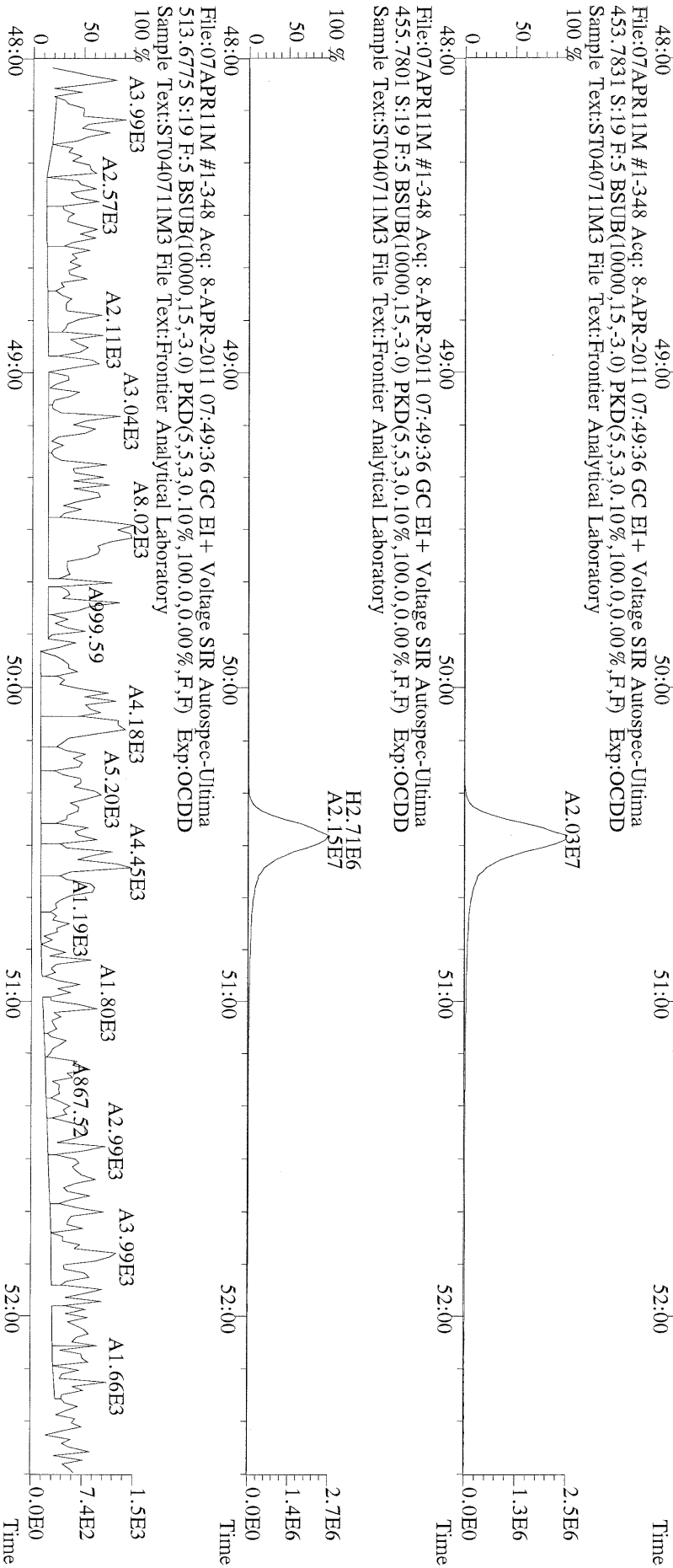
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100 %



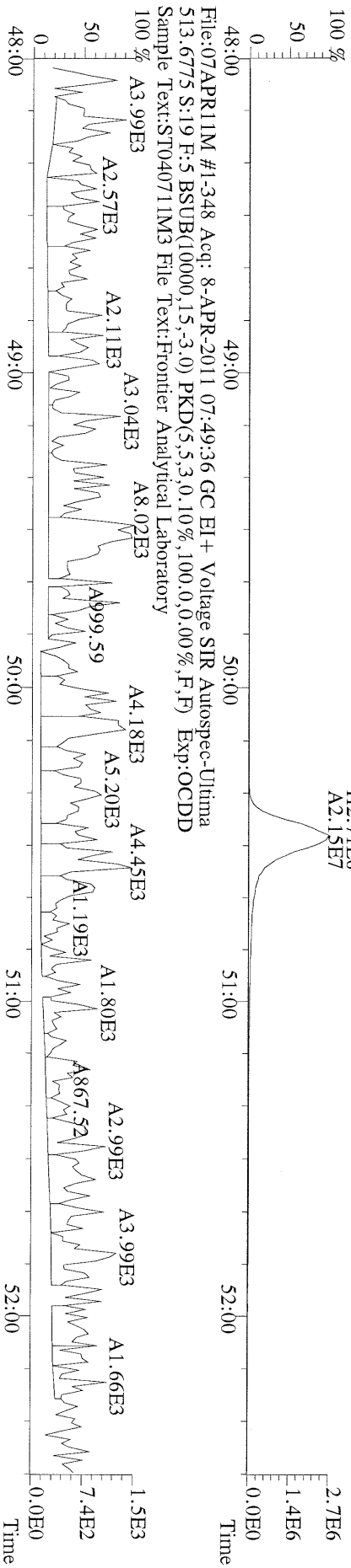
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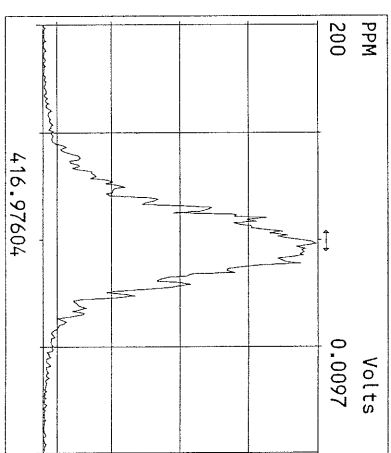
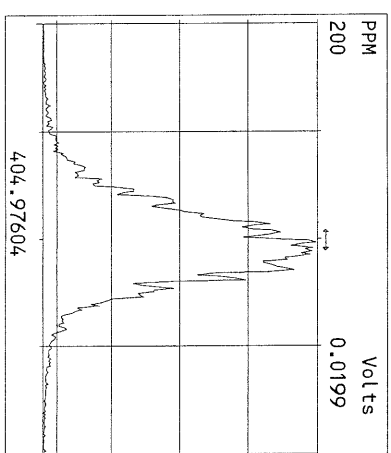
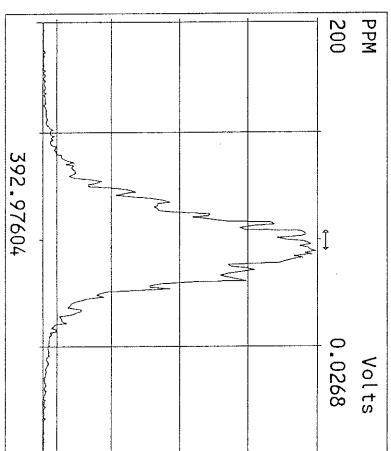
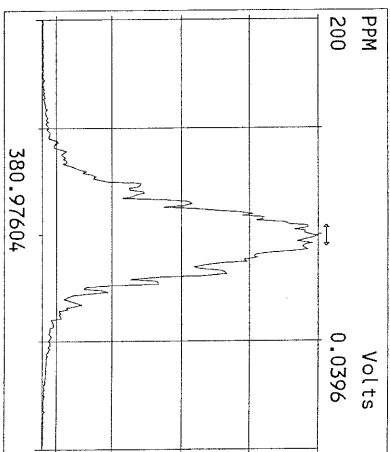
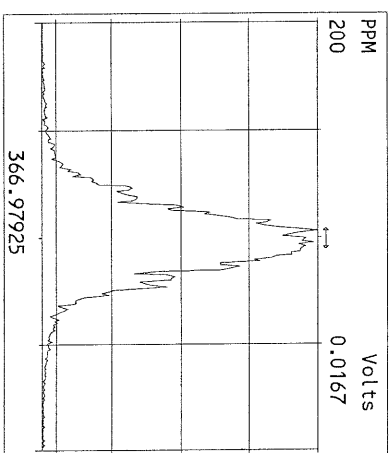
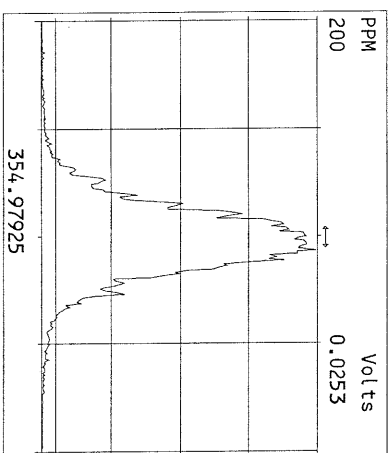
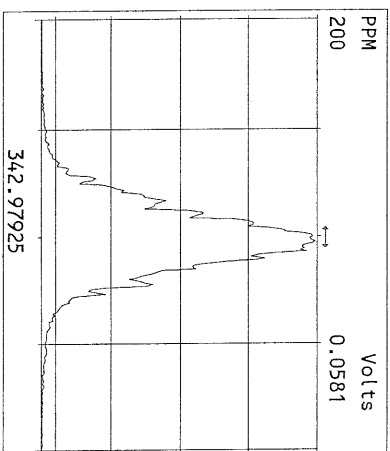
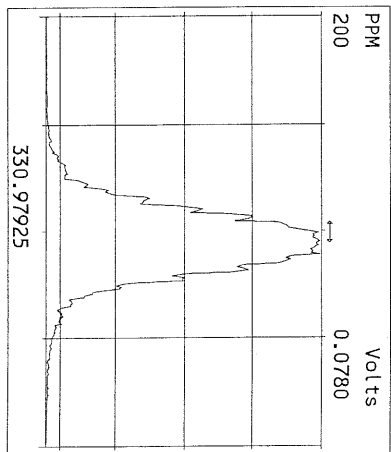
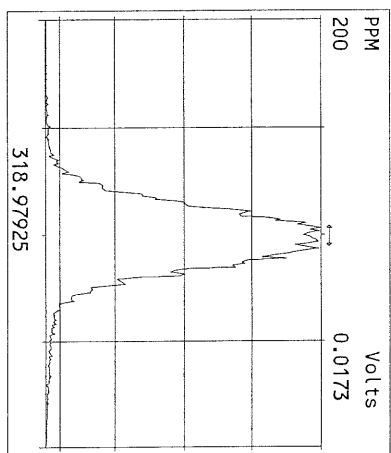
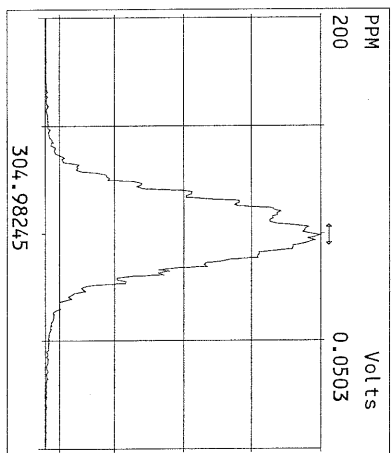
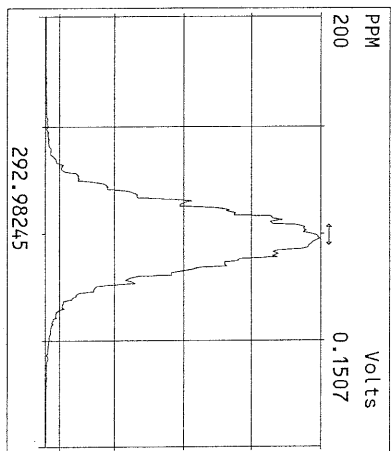


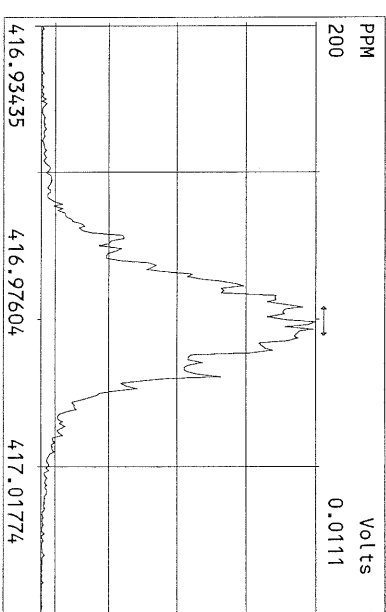
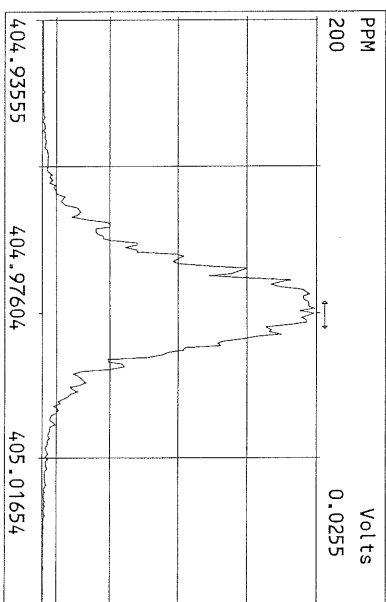
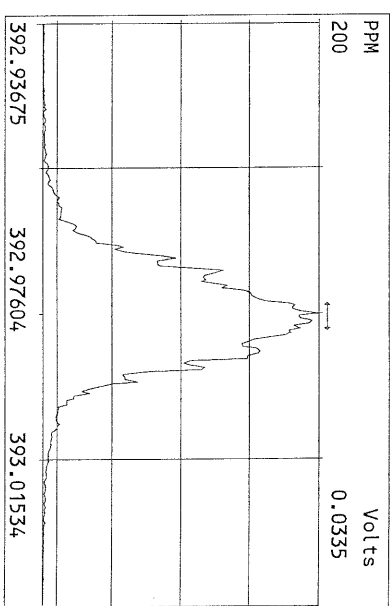
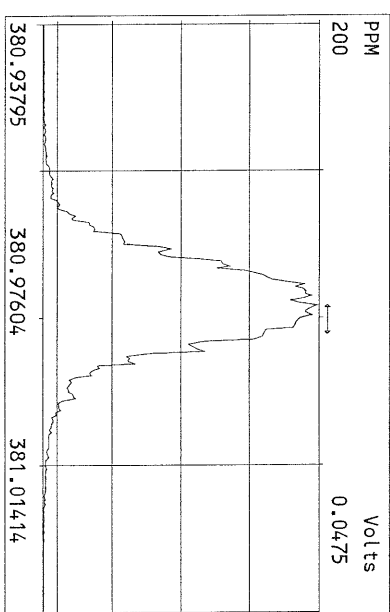
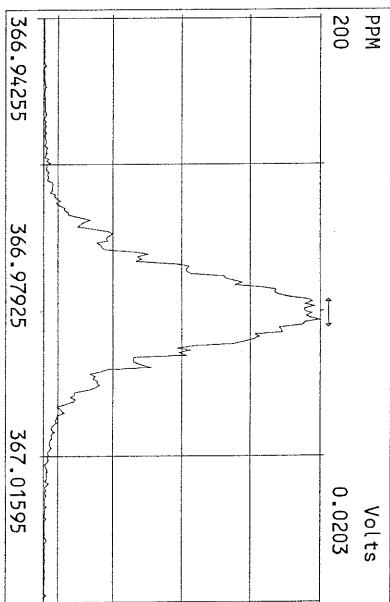
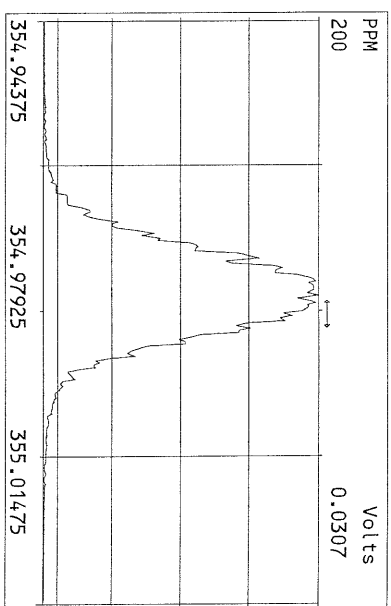
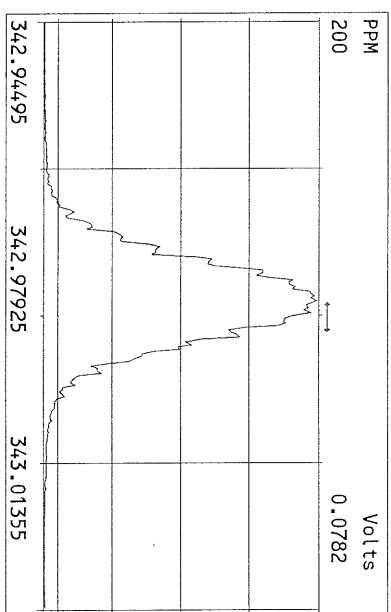
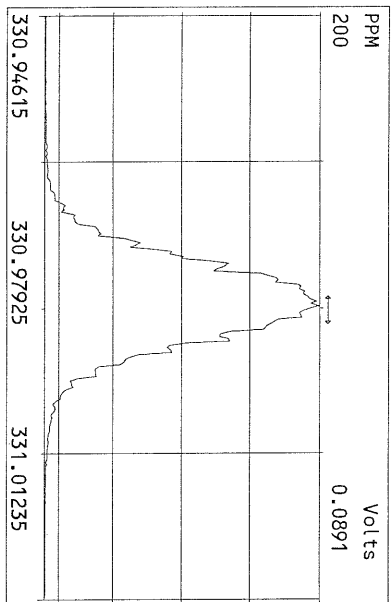
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453.7831 S:19 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM3 File Text:Frontier Analytical Laboratory

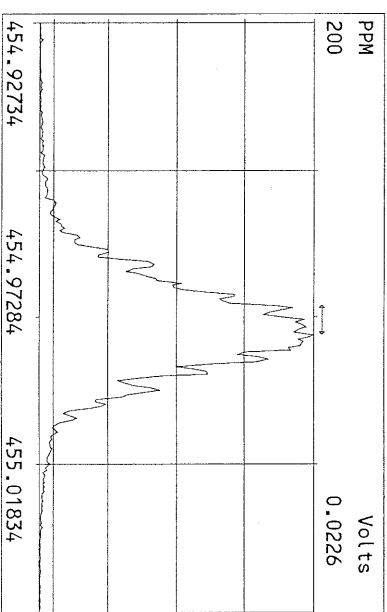
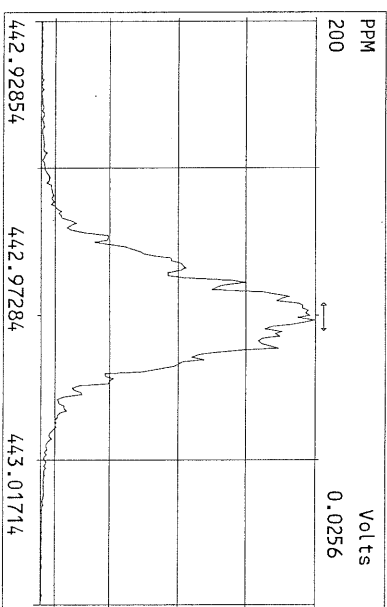
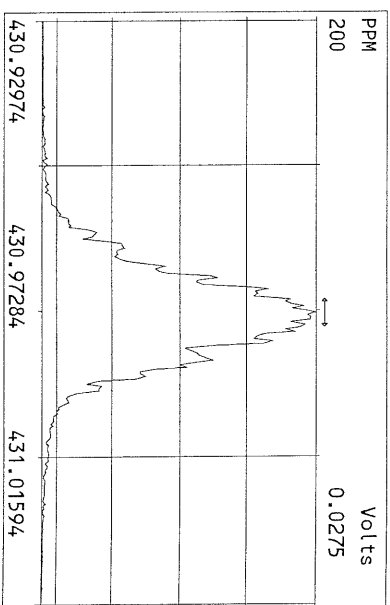
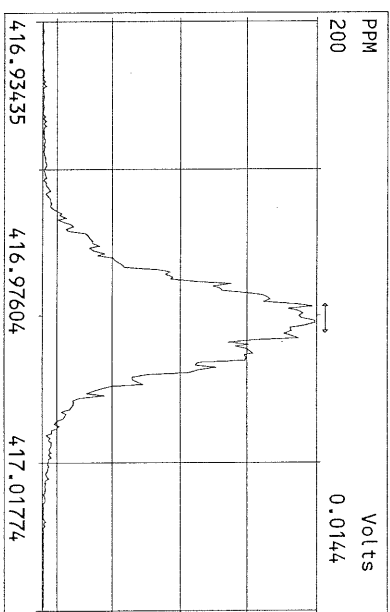
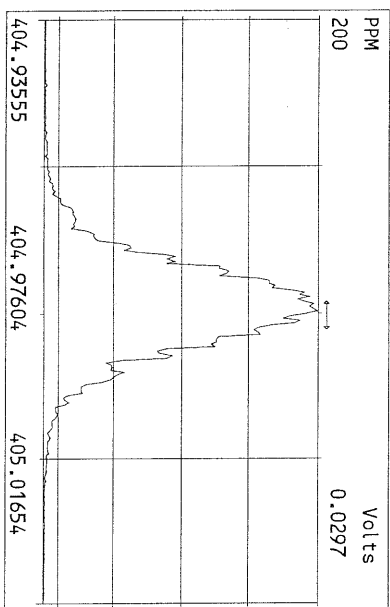
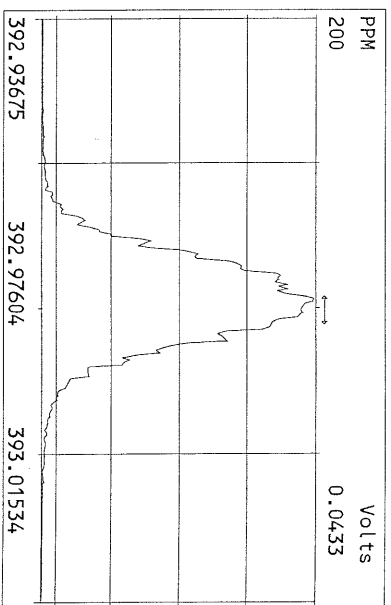
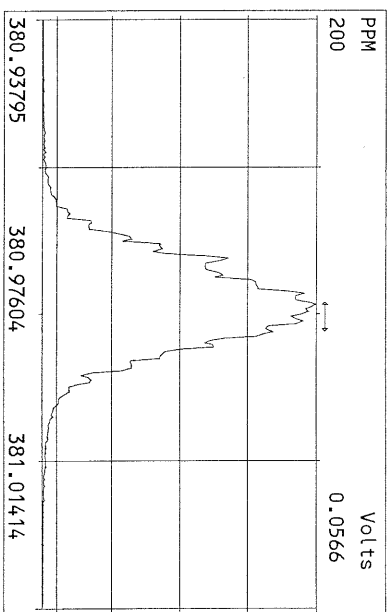
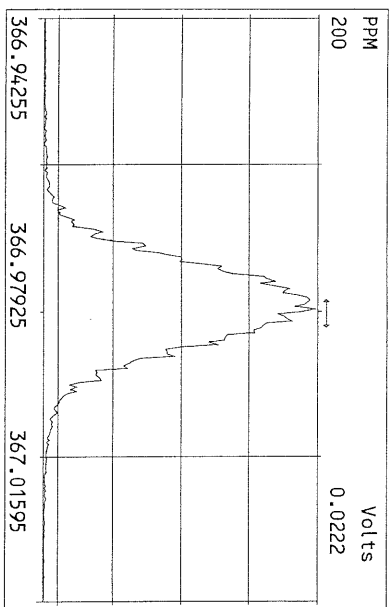


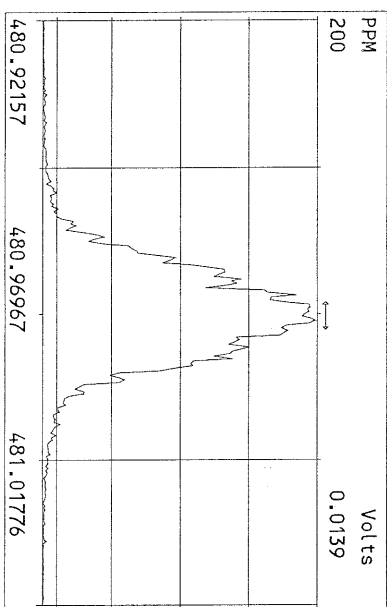
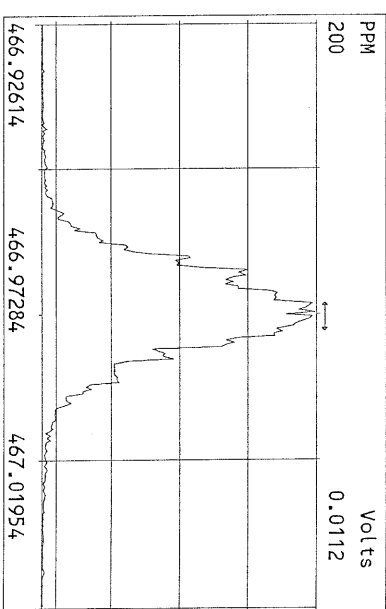
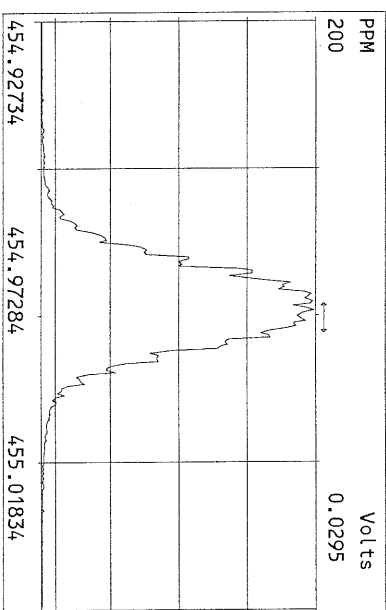
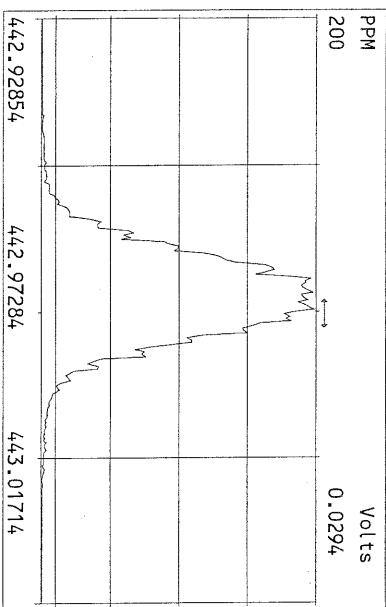
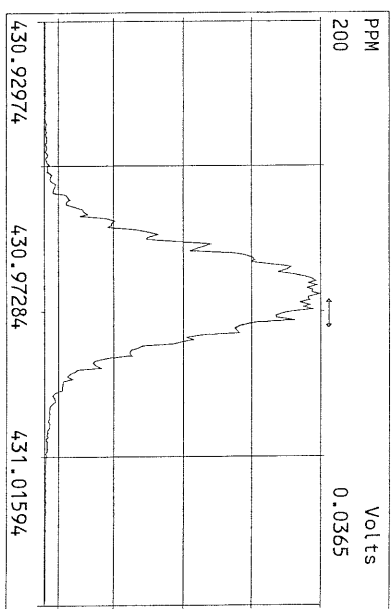
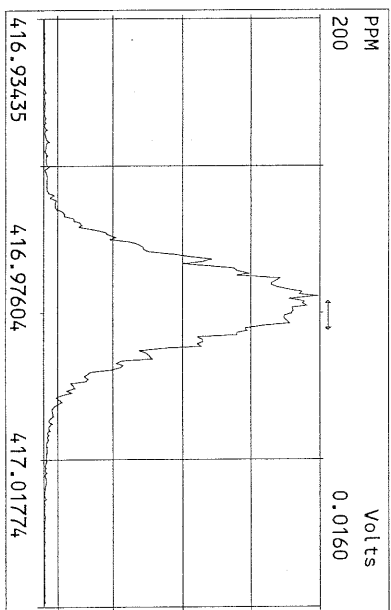
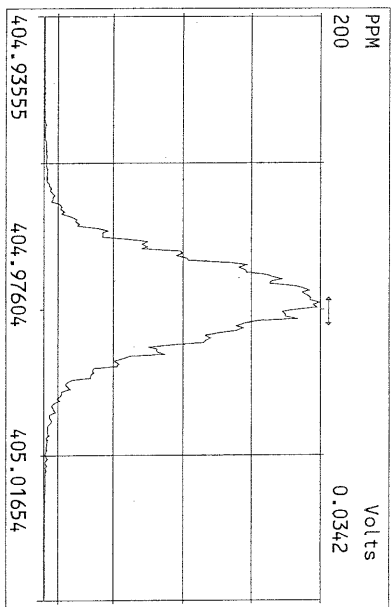
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513.6775 S:19 F:5 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,00%,F,F) Exp:OCDD
Sample Text:ST04071IM3 File Text:Frontier Analytical Laboratory

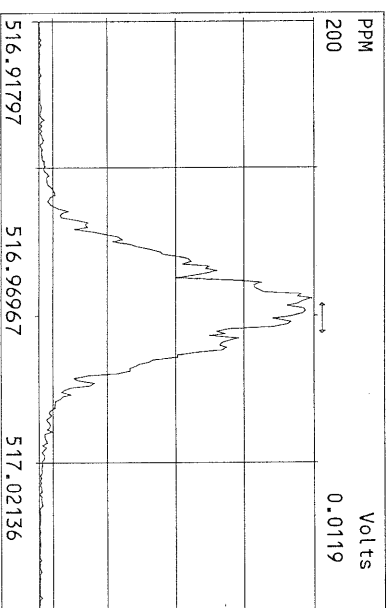
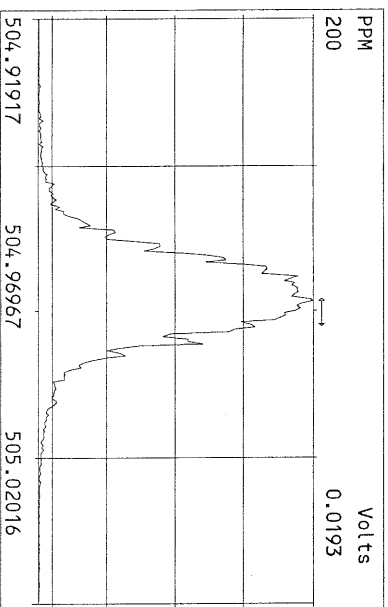
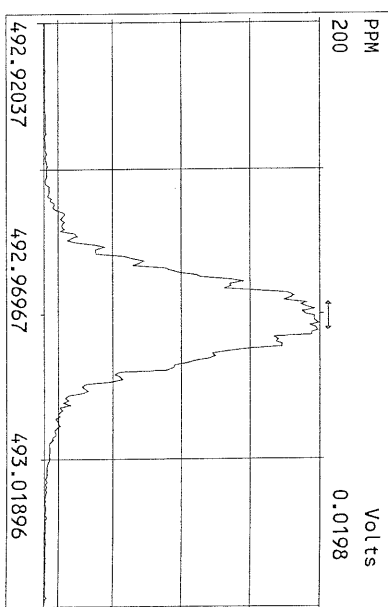
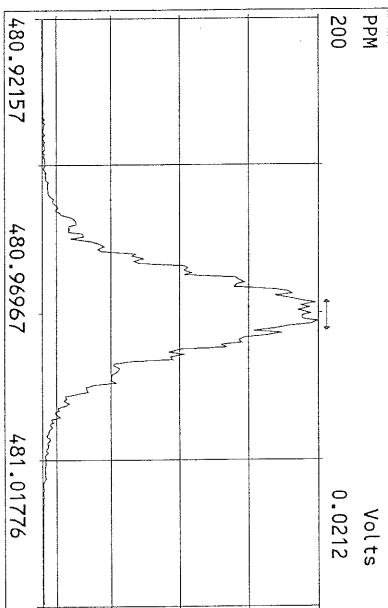
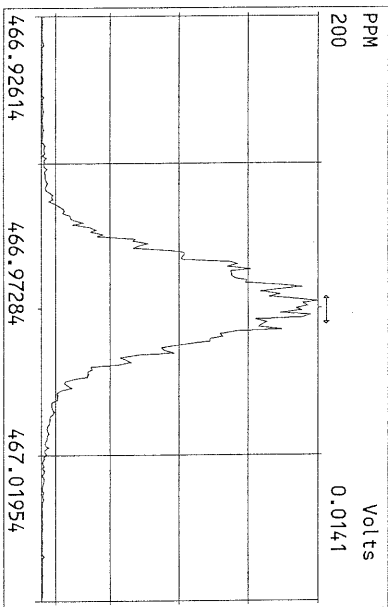
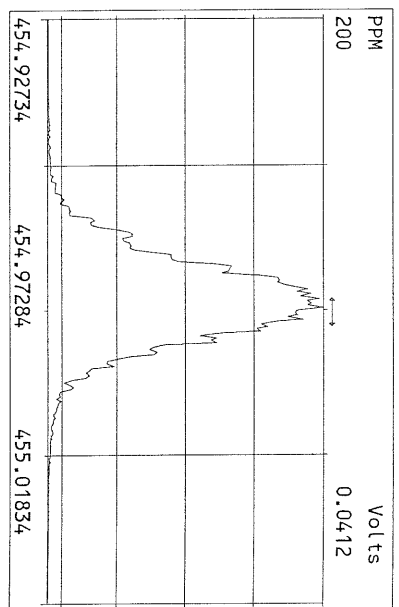
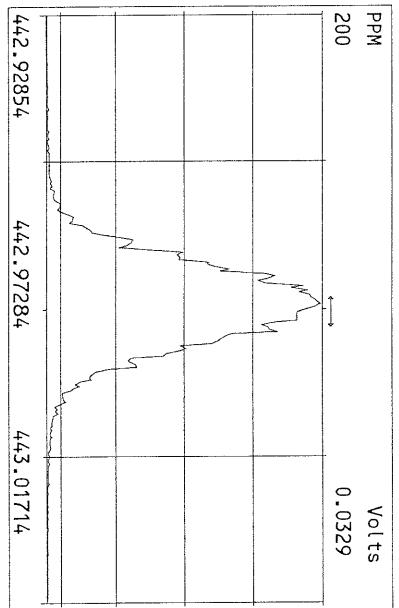
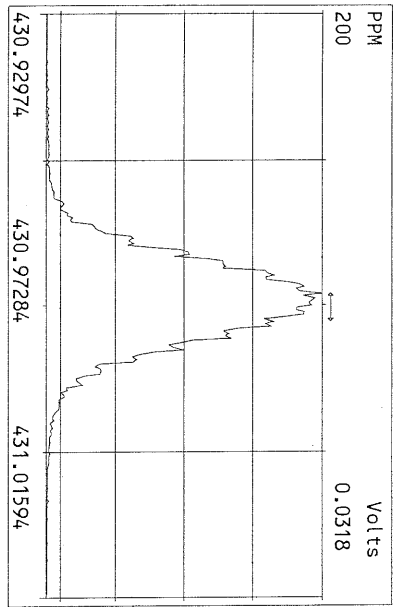












USEPA - ITD

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: 08APR11A Sam:1 Analysis Date: 8-APR-11 Time: 09:21:16


| | 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 | 0.78 | 0.65-0.89 | y | 9.02 | 8.40 - 12.0 |
| LABELED COMPOUNDS | | | | | | |
| 13C-2,3,7,8-TCDF | M/M+2 | 0.80 | 0.65-0.89 | y | 94.8 | 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: 4/8/11

FAL ID: ST040811A1 Filename: 08APR11A Sam:1 Acquired: 8-APR-11 09:21:16 ICal: tcdffa11-2-18-11
Client ID: 1613 CS3 100511J ConCal: ST040811A1 EndCal: ST040811A2
Results: 6624TCDF GC Column: DB225 Amount: 1.000

| Name | Resp | RA | RT | RRF | Conc | Qual | Fac | Noise | DL | #Hom | Rec |
|------------------|----------|--------|-------|------|------|------|------|-------|----|------|------|
| 2,3,7,8-TCDF | 2.41e+07 | 0.78 y | 19:22 | 1.16 | 9.02 | | 2.50 | - | - | 1 | |
| 13C-2,3,7,8-TCDF | 2.31e+08 | 0.80 y | 19:21 | 1.05 | 94.8 | | | | | | 94.8 |
| 13C-1,2,3,4-TCDF | 2.31e+08 | 0.80 y | 16:48 | - | 131 | | | | | | |

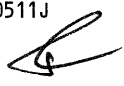
Analyst: 

Date: 4/8/11

Frontier Analytical Laboratory - Acquisition Log

Run Name:08APR11A Instrument: FAL1 GC: DB225 Experiment:TCDF

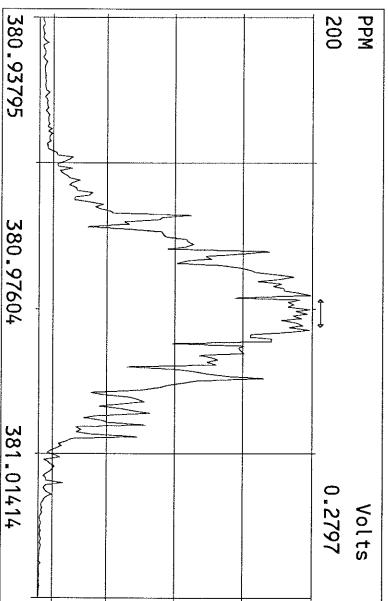
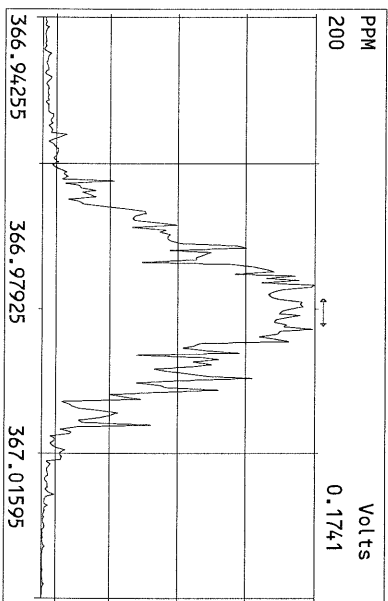
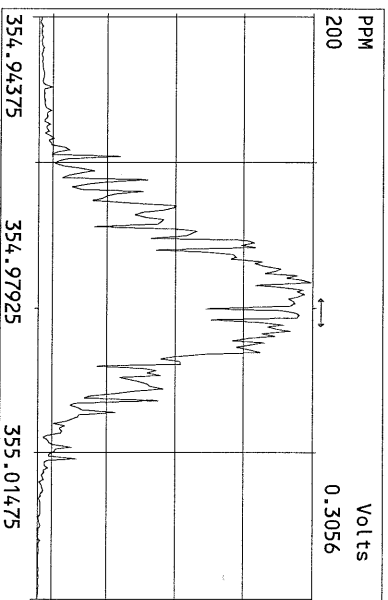
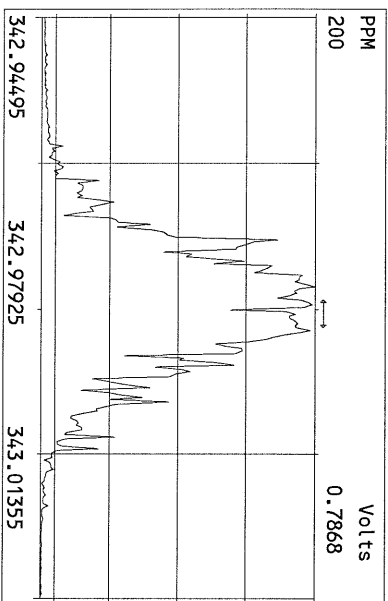
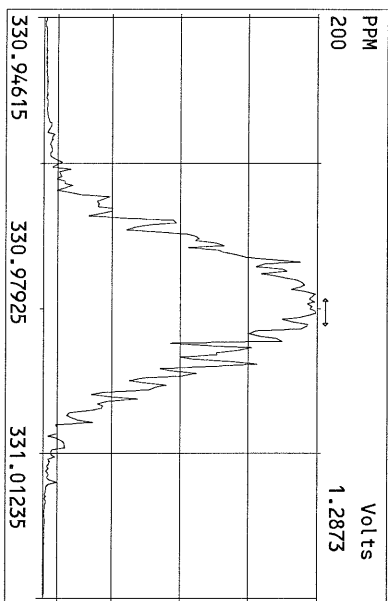
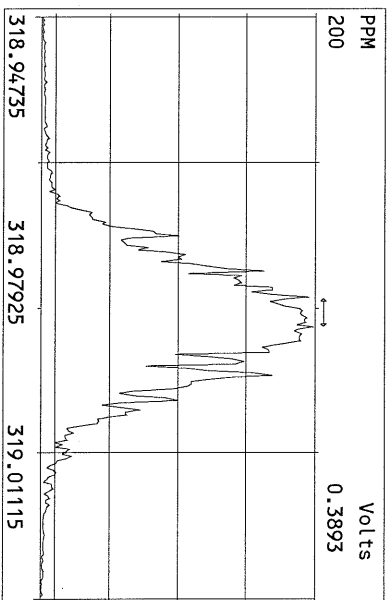
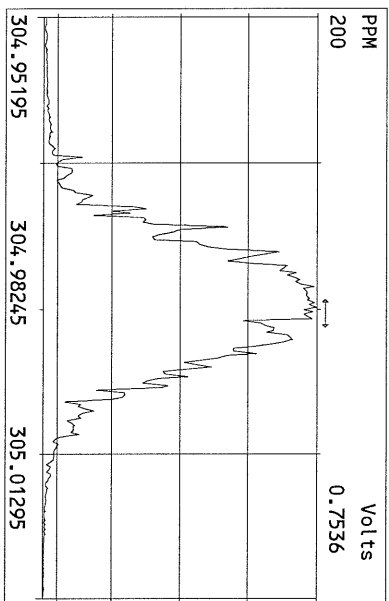
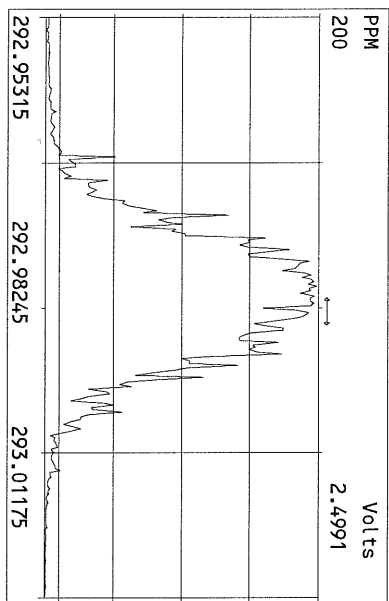
| Data File | S | FAL ID | Client ID | Acquired | ConCal | EndCal | Analyst |
|-----------|---|------------------|-----------------------|-------------------|------------|------------|---------|
| 08APR11A | 1 | ST040811A1 | 1613 CS3 100511J | 8-APR-11 09:21:16 | ST040811A1 | ST040811A2 | TC |
| 08APR11A | 2 | 6624-001-0001-SA | Lot #018690 | 8-APR-11 09:56:19 | ST040811A1 | ST040811A2 | TC |
| 08APR11A | 3 | 6678-004-0001-SA | LL-SED2-0-56-031511 | 8-APR-11 10:31:24 | ST040811A1 | ST040811A2 | TC |
| 08APR11A | 4 | 6678-007-0001-SA | LL-SED2-0-56-031511-D | 8-APR-11 11:06:27 | ST040811A1 | ST040811A2 | TC |
| 08APR11A | 5 | 6678-001-0001-SA | LL-SED3-0-36-031511 | 8-APR-11 11:41:32 | ST040811A1 | ST040811A2 | TC |
| 08APR11A | 6 | ST040811A2 | 1613 CS3 100511J | 8-APR-11 12:16:35 | ST040811A1 | ST040811A2 | TC |



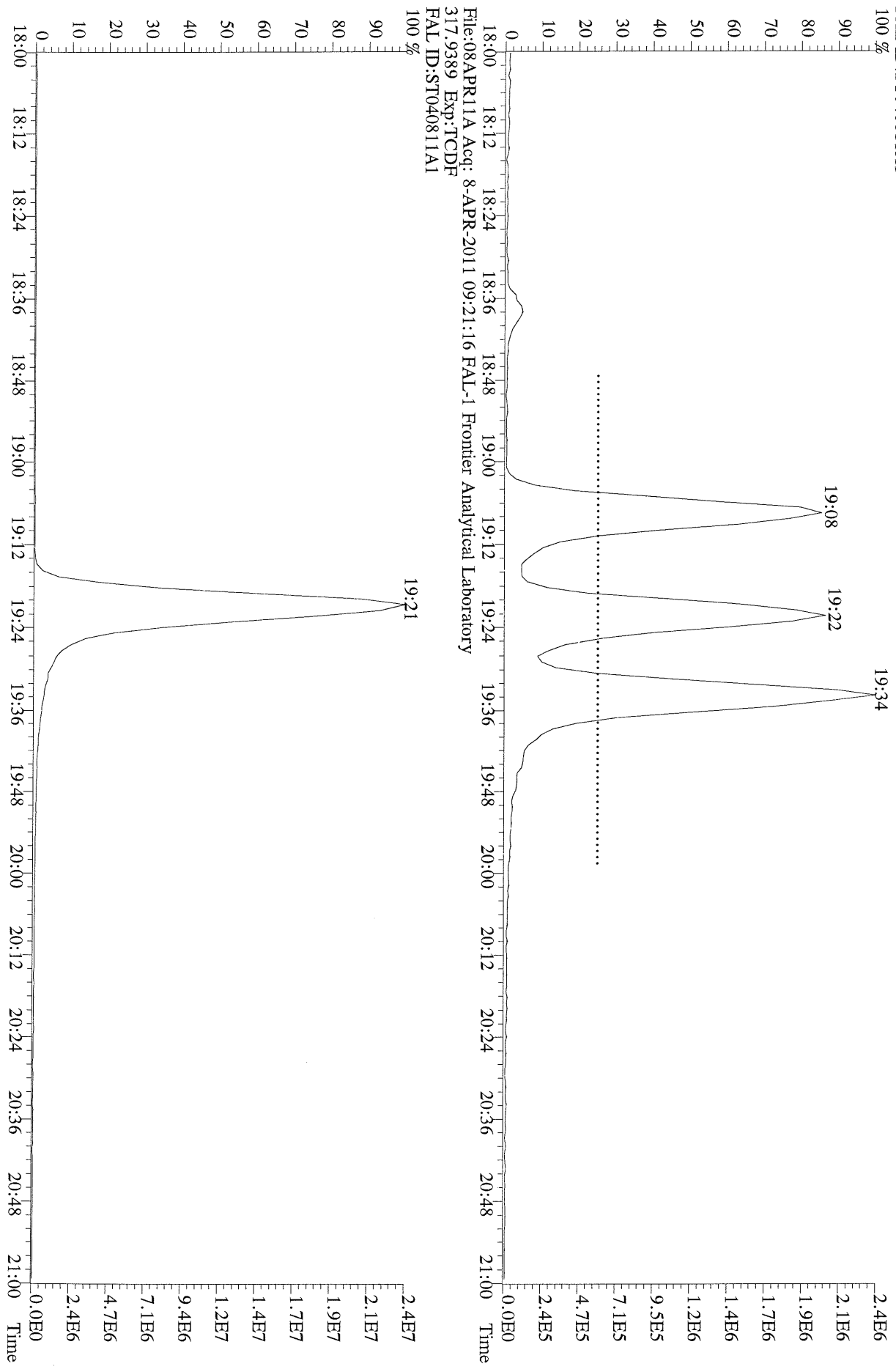
9/8/11

Data Backed Up: _____

Date: _____



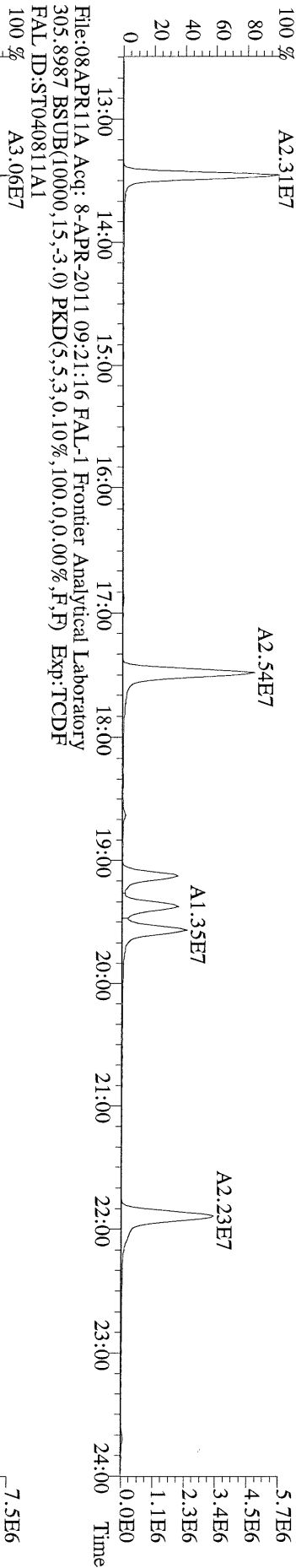
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303.9016 Exp:TCDF
FAL ID:ST040811A1



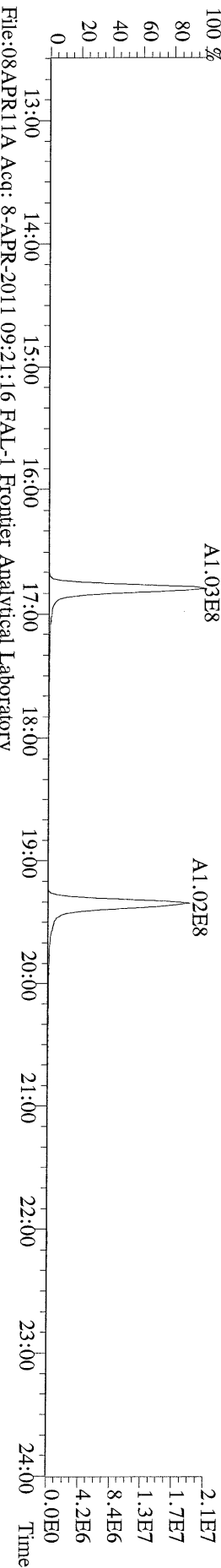
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317.9389 Exp:TCDF
FAL ID:ST040811A1

2.4E6
2.1E6
1.9E6
1.7E6
1.4E6
1.2E6
9.5E5
7.1E5
4.7E5
2.4E5
0.0E0
2.4E7
2.1E7
1.9E7
1.7E7
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1.2E7
9.4E6
7.1E6
4.7E6
2.4E6
0.0E0
Time

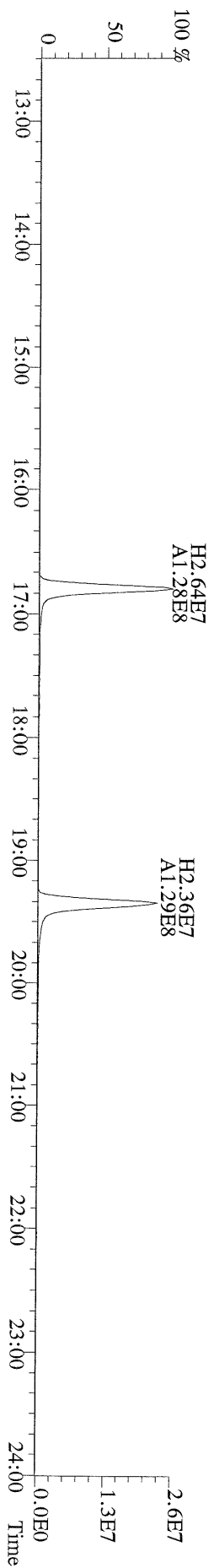
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FAL ID:ST040811A1

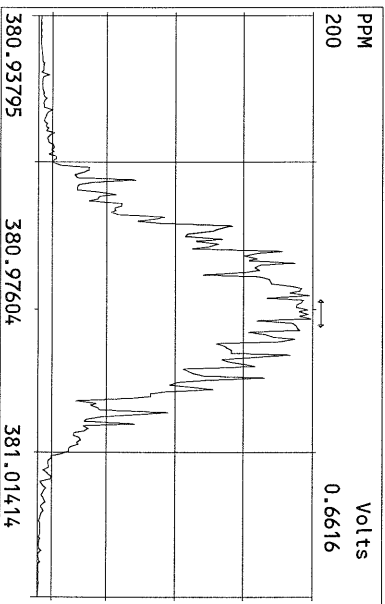
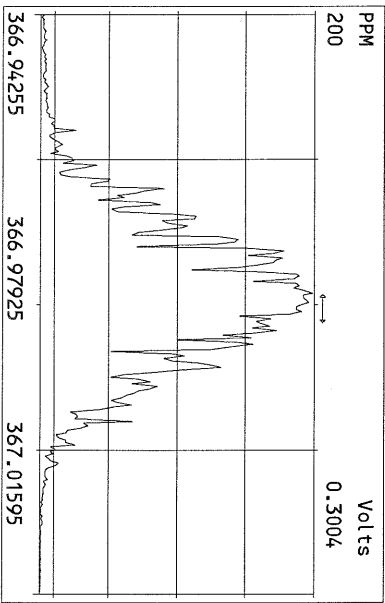
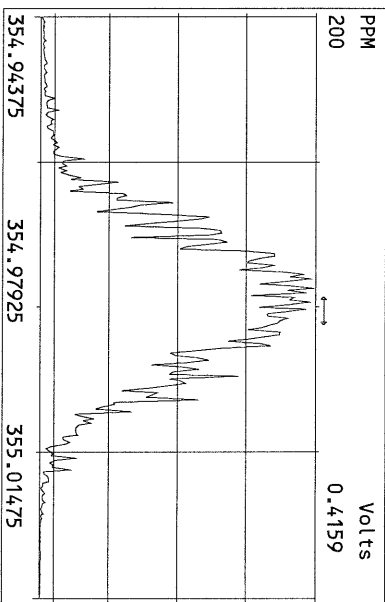
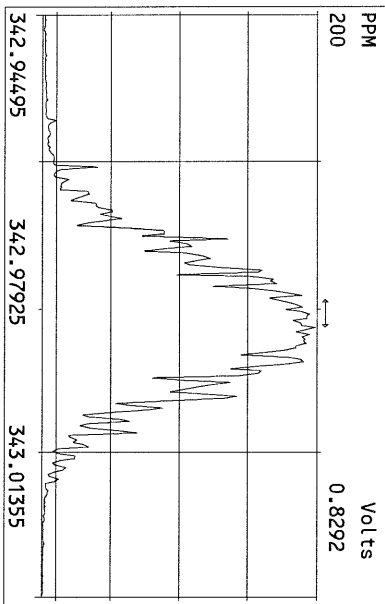
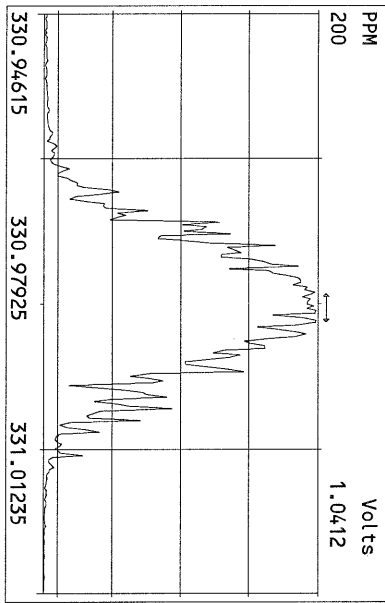
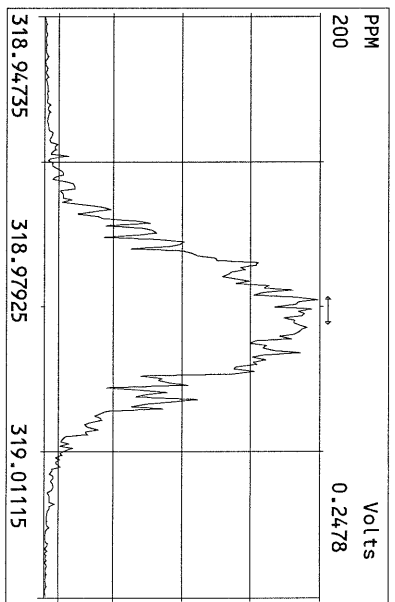
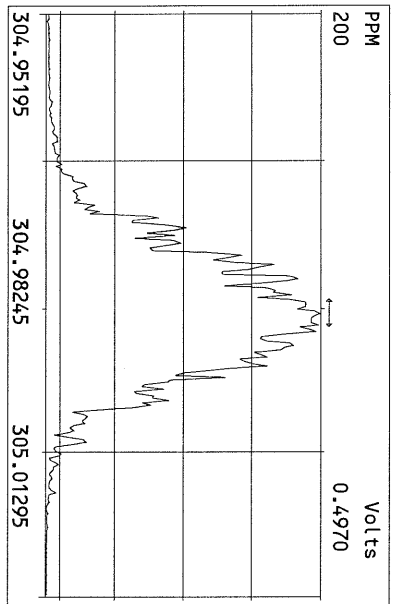
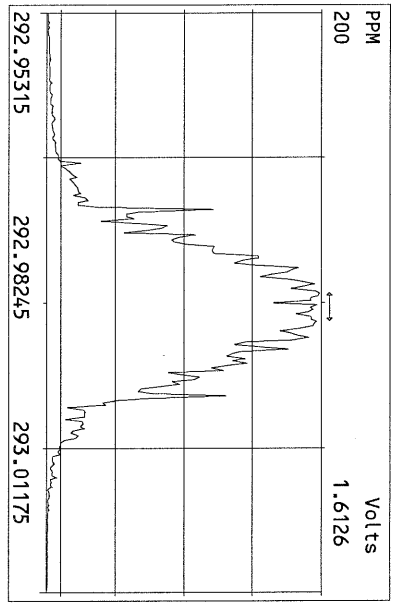


File:08APR11A Acq: 8-APR-2011 09:21:16 FAL-1 Frontier Analytical Laboratory
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FAL ID:ST040811A1



File:08APR11A Acq: 8-APR-2011 09:21:16 FAL-1 Frontier Analytical Laboratory
317.9389 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:TCDF
FAL ID:ST040811A1





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: 08APR11A Sam:6

Analysis Date: 8-APR-11 Time: 12:16:35

| | 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 | 0.77 | 0.65-0.89 | y | 8.54 | 8.40 - 12.0 |
| LABELED COMPOUNDS | | | | | | |
| 13C-2,3,7,8-TCDF | M/M+2 | 0.80 | 0.65-0.89 | y | 95.0 | 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: 4/8/11

FAL ID: ST040811A2 Filename: 08APR11A Sam:6 Acquired: 8-APR-11 12:16:35 ICal: tcdfal1-2-18-11
Client ID: 1613 CS3 100511J ConCal: ST040811A1 EndCal: ST040811A2
Results: 6678TCDF GC Column: DB225 Amount: 1.000

| Name | Resp | RA | RT | RRF | Conc | Qual | Fac | Noise | DL | #Hom | Rec |
|------------------|----------|--------|-------|------|------|------|------|-------|----|------|------|
| 2,3,7,8-TCDF | 1.56e+07 | 0.77 y | 19:18 | 1.16 | 8.54 | | 2.50 | - | - | 1 | |
| 13C-2,3,7,8-TCDF | 1.58e+08 | 0.80 y | 19:16 | 1.05 | 95.0 | | | | | | 95.0 |
| 13C-1,2,3,4-TCDF | 1.58e+08 | 0.79 y | 16:43 | - | 89.5 | | | | | | |

Analyst:  Date: 4/8/11

Frontier Analytical Laboratory - Acquisition Log

Run Name:08APR11A Instrument: FAL1 GC: DB225 Experiment:TCDF

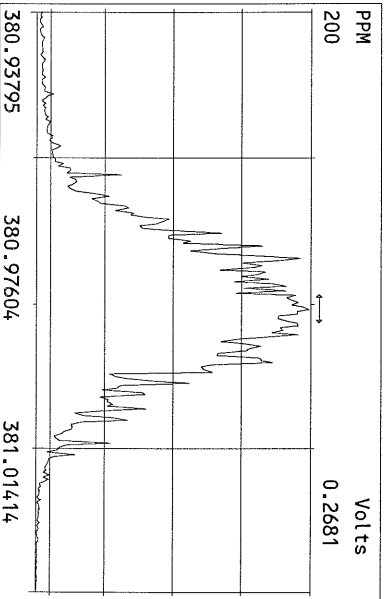
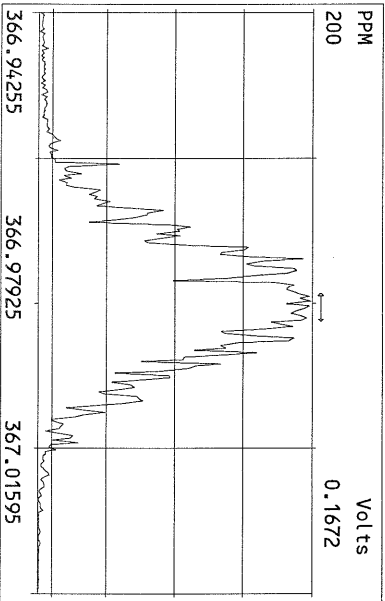
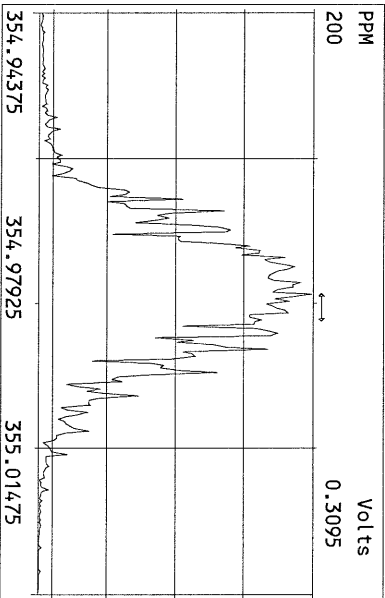
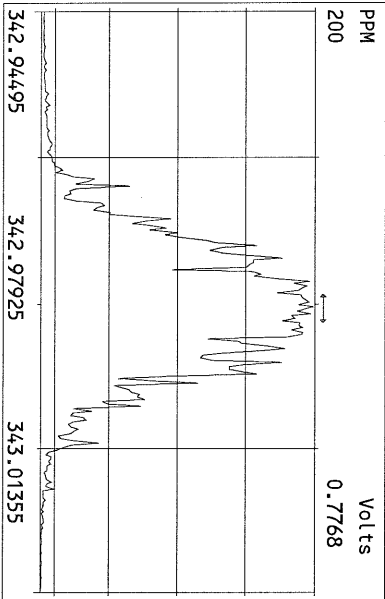
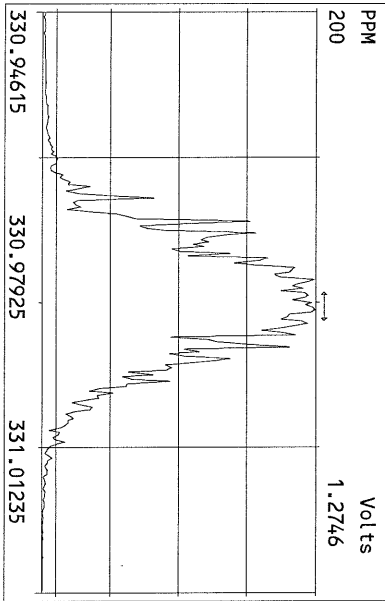
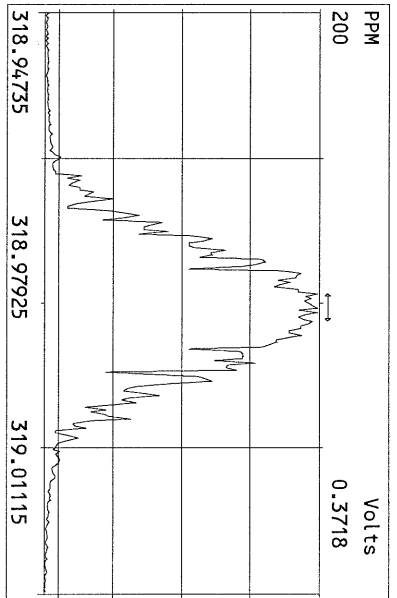
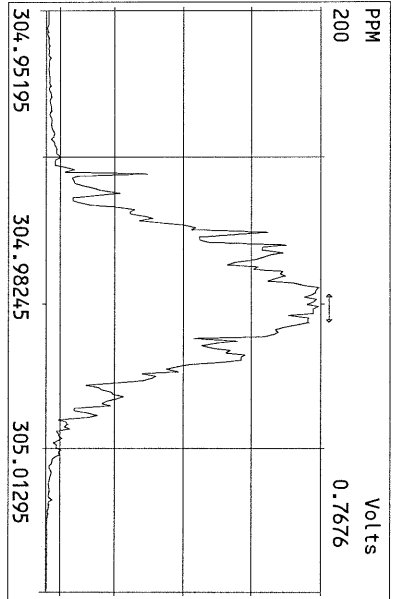
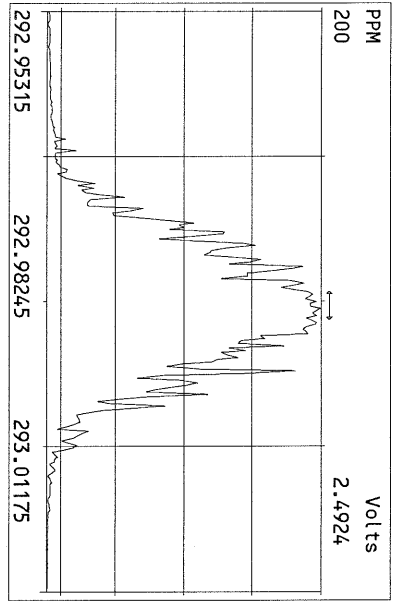
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|-------------|------------------|-----------------------|-------------------|------------|------------|---------|
| 08APR11A 1 | ST040811A1 | 1613 CS3 100511J | 8-APR-11 09:21:16 | ST040811A1 | ST040811A2 | TC |
| 08APR11A 2 | 6624-001-0001-SA | Lot #018690 | 8-APR-11 09:56:19 | ST040811A1 | ST040811A2 | TC |
| 08APR11A 3 | 6678-004-0001-SA | LL-SED2-0-56-031511 | 8-APR-11 10:31:24 | ST040811A1 | ST040811A2 | TC |
| 08APR11A 4 | 6678-007-0001-SA | LL-SED2-0-56-031511-D | 8-APR-11 11:06:27 | ST040811A1 | ST040811A2 | TC |
| 08APR11A 5 | 6678-001-0001-SA | LL-SED3-0-36-031511 | 8-APR-11 11:41:32 | ST040811A1 | ST040811A2 | TC |
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TC

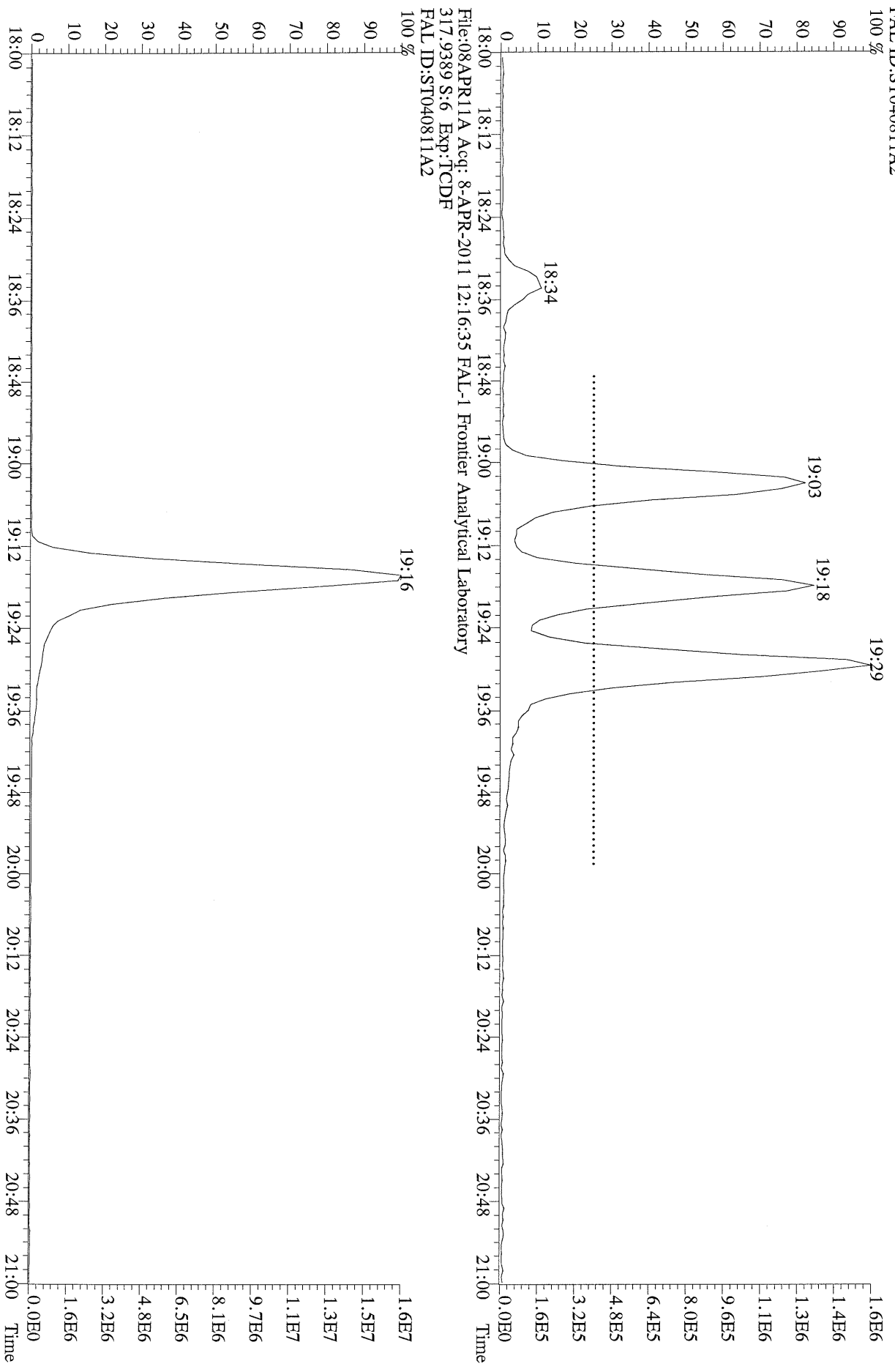
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Date: _____

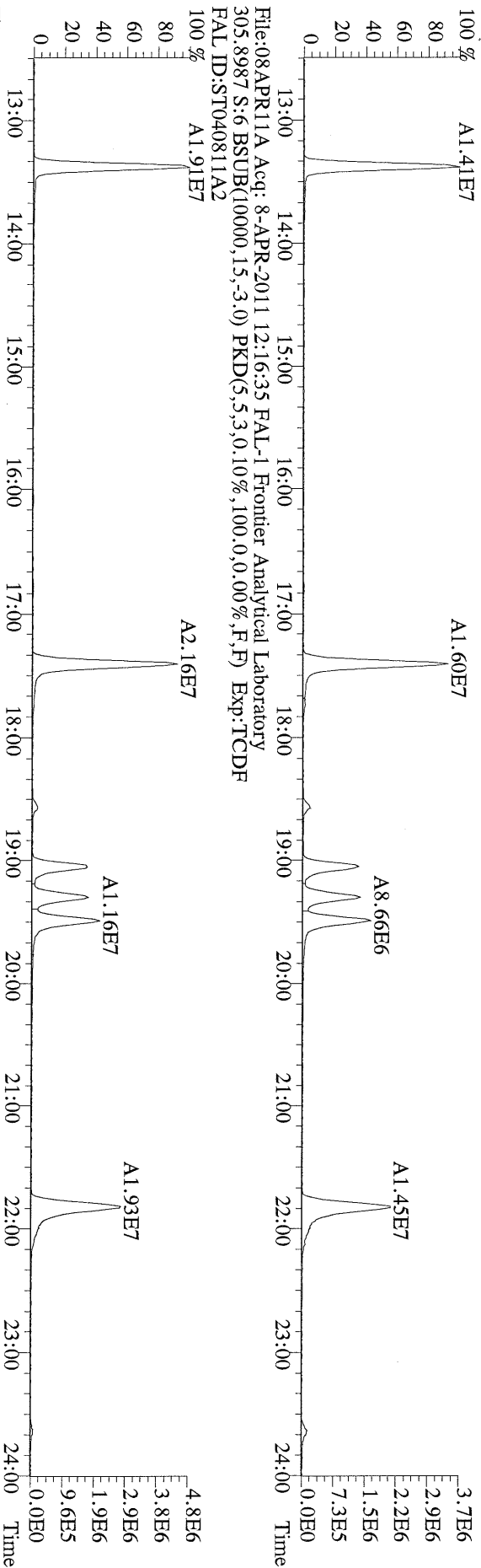


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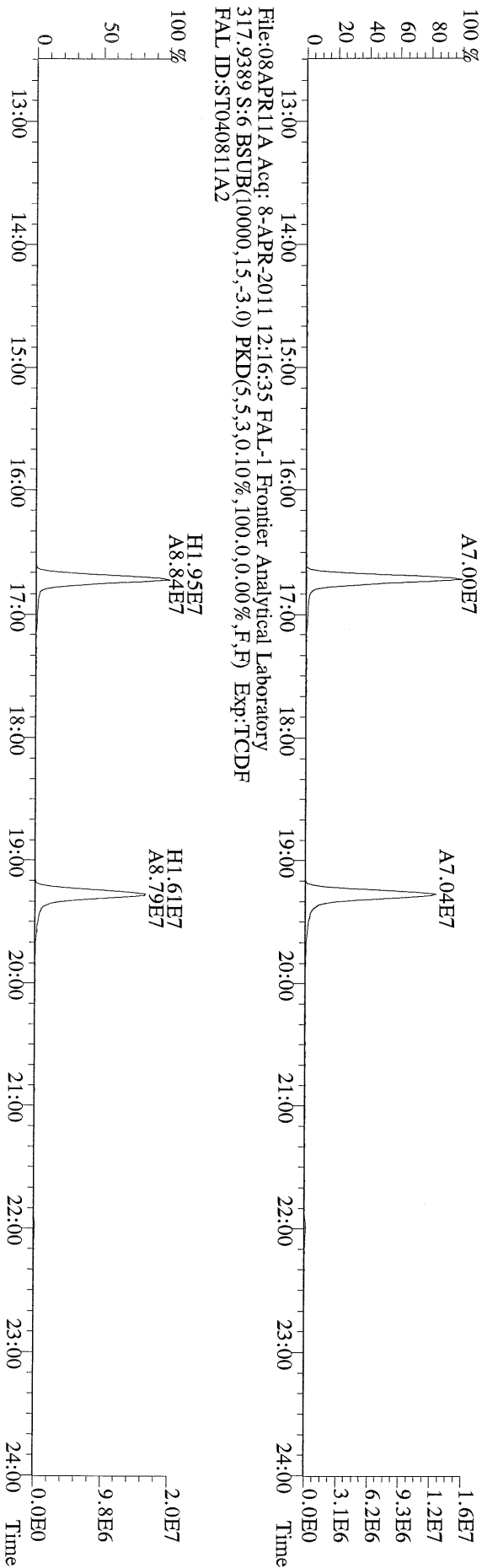


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317.9389 S:6 Exp:TCDF
FAL ID:ST040811A2

File:08APR11A Acq: 8-APR-2011 12:16:35 FAL-1 Frontier Analytical Laboratory
303.9016 S:6 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:TCDF
FAL ID:ST040811A2



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FAL ID:ST040811A2



File:08APR11A Acq: 8-APR-2011 12:16:35 FAL-1 Frontier Analytical Laboratory
317.9389 S:6 BSUB(10000,15,-3.0) PKD(5,5,3,0,10%,100,0,0,0.00%,F,F) Exp:TCDF
FAL ID:ST040811A2

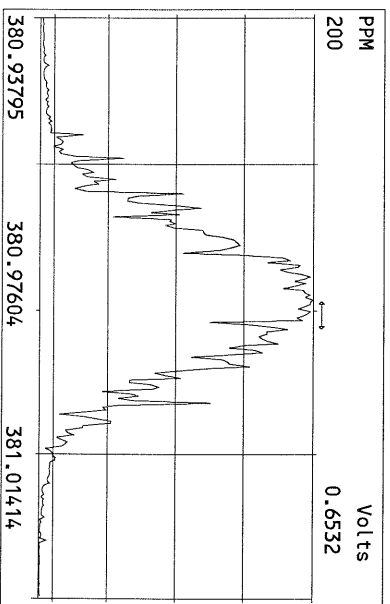
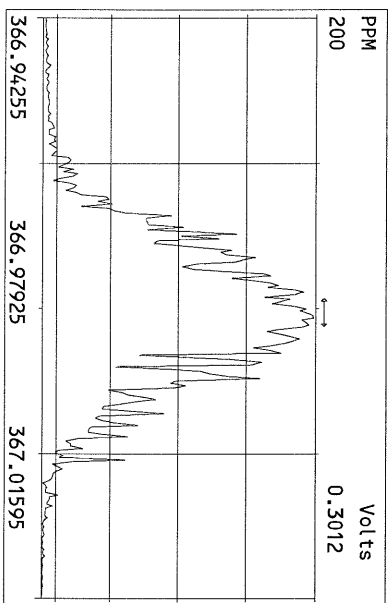
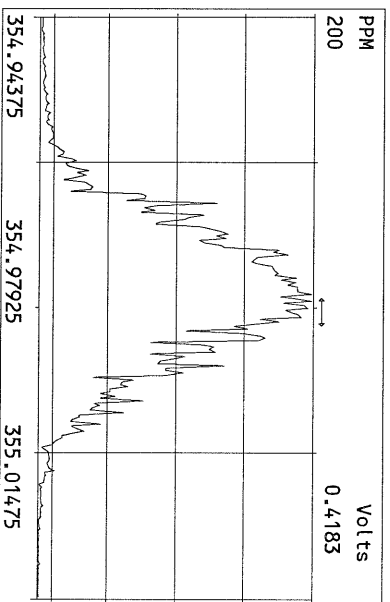
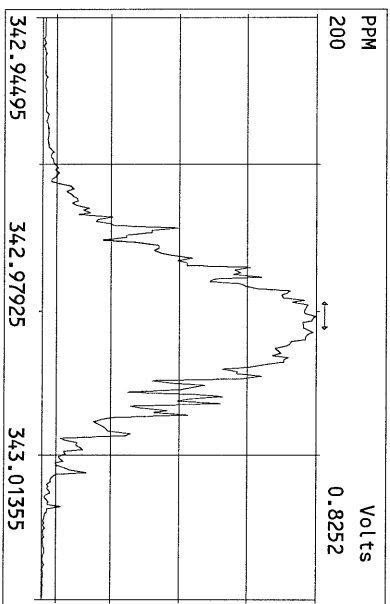
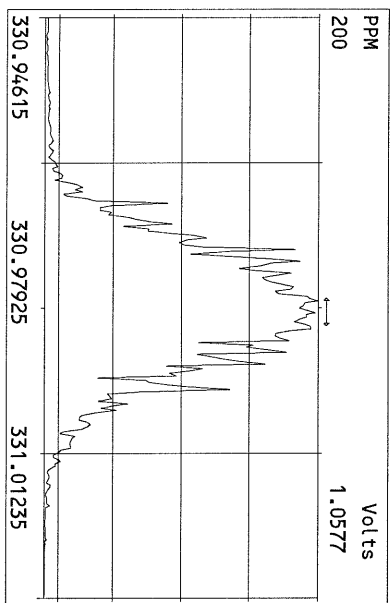
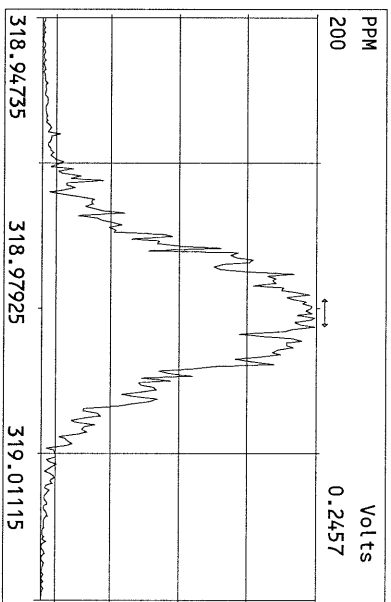
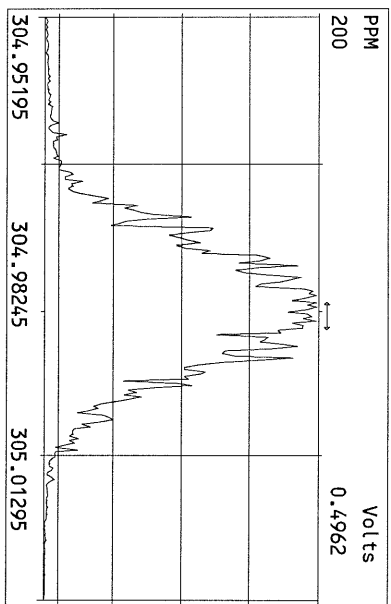
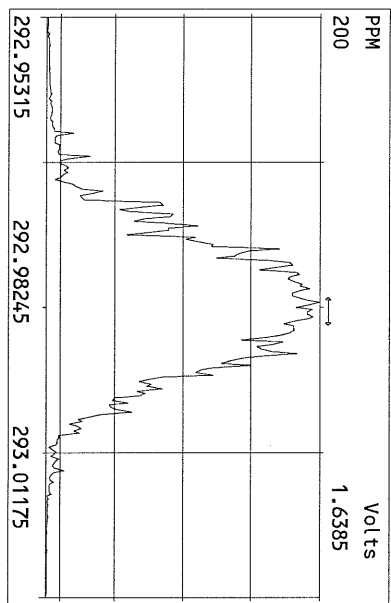


Table of Contents: ARI Job SN54

Client: Floyd Snider

Project: POS-LL Lora Lake - Subsurface Sediment

| | Page From: | Page To: |
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| General Chemistry Raw Data | | |
| Analyst Notes and Raw Data | <u>691</u> | <u>718</u> |
| Geotechnical Raw Data | | |
| Analyst Notes and Raw Data | <u>719</u> | <u>720</u> |

 PC
Signature

March-24-2011
Date



Analytical Resources, Incorporated
Analytical Chemists and Consultants

April 21, 2011

Erin Breckel
Floyd-Snyder Inc.
601 Union Street, Suite 600
Seattle, WA 98101-2341

RE: Client Project: Lora Lake Subsurface Sediment, POS-LL
ARI Job No: SN54

Dear Ms. Breckel:

Please find enclosed the original Chain-of-Custody (COC) record, sample receipt documentation, and the final data package for samples from the project referenced above.

Sample receipt and detail of these analyses are discussed in the Case Narrative.

An electronic copy of this package will remain on file with ARI. Should you have any questions or problems, please feel free to contact me at your convenience.

Sincerely,

ANALYTICAL RESOURCES, INC.

Cheronne Oreiro
Project Manager

-For-

Susan D. Dunning
Director, Client Services
sue@arilabs.com
206-695-6207

Enclosures

cc: eFile SN54

Chain of Custody Documentation

ARI Job ID: SN54

Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: **SNS4** Turn-around Requested: **Standard**

ARI Client Company: **Floyd Snider** Phone: **206-2512-2078**

Client Contact: **Erin Breckel**

Client Project Name: **Lower Laine - Subsurface Sediment**

Client Project #: **POS-LL** Samplers: **E-B/A/M**

Page: **1** of **1**

Date: **3/15/11** Ice Present? **Y**

No. of Coolers: **1** Cooler Temps: **1.5**



Analytical Resources, Incorporated
Analytical Chemists and Consultants
4611 South 134th Place, Suite 100
Tukwila, WA 98168
206-695-6200 206-695-6201 (fax)

| Sample ID | Date | Time | Matrix | No. Containers | Analysis Requested | | | | | Notes/Comments | |
|--|---------|-------|----------|----------------|----------------------------|-----------------------------|---------------------|---------------|------------|--|-------------------------|
| | | | | | CADH (B70-Site Code) | DG (B04) | Metals (AS, Pb, Cr) | Dioxin (1613) | TOC (PSEP) | | |
| LL-SED 3-0-36-031511 | 3/15/11 | 12:30 | Sediment | 4 | X | X | X | X | X | | |
| LL-SED 3-36-141-031511 | | 12:40 | | 4 | X | X | X | X | X | Hold - will call if we decide to analyze | |
| LL-SED 3-141-167-031511 | | 12:50 | | 4 | X | X | X | X | X | | |
| LL-SED 2-0-56-031511 | | 14:15 | | 4 | X | X | X | X | X | | |
| LL-SED 2-56-112-031511 | | 14:25 | | 8 | X | X | X | X | X | MS/MSD Volume | |
| LL-SED 2-112-168-031511 | | 14:55 | | 4 | X | X | X | X | X | | |
| LL-SED 2-0-56-031511-D | | 14:15 | | 4 | X | X | X | X | X | | |
| LL-SED 1-0-56-031511 | | 18:00 | ↓ | 4 | X | X | X | X | X | | |
| Comments/Special Instructions | | | | | Received by (Signature) | Relinquished by (Signature) | | | | | Received by (Signature) |
| Samples frozen, need to be thawed and homogenized. | | | | | Erin Breckel | | | | | | Jennifer Millsap |
| | | | | | Printed Name | Printed Name | | | | | Company |
| | | | | | Company: ARI | | | | | Company | |
| | | | | | Date & Time: 3/16/11 10:10 | | | | | Date & Time: 3/16/11 10:10 | |

Limits of Liability: ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, not withstanding any provision to the contrary in any contract, purchase order or co-signed agreement between ARI and the Client.

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.

SNS4-00000



Cooler Receipt Form

ARI Client: floyd Snider

Project Name: Lora Lake- Subsurface Sediment

COC No(s): _____ (NA)

Delivered by: Fed-Ex UPS Courier Hand Delivered Other: _____

Assigned ARI Job No: SN54

Tracking No: _____ (NA)

Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO

Were custody papers included with the cooler? YES NO

Were custody papers properly filled out (ink, signed, etc.) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry)..... 1.5

If cooler temperature is out of compliance fill out form 00070F Temp Gun ID#: 90041619

Cooler Accepted by: ML Date: 3/16/11 Time: 1000

Complete custody forms and attach all shipping documents

Log-In Phase:

Was a temperature blank included in the cooler? YES NO

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other: _____

Was sufficient ice used (if appropriate)? NA YES NO

Were all bottles sealed in individual plastic bags? YES NO

Did all bottles arrive in good condition (unbroken)? YES NO

Were all bottle labels complete and legible? YES NO

Did the number of containers listed on COC match with the number of containers received? YES NO

Did all bottle labels and tags agree with custody papers? YES NO

Were all bottles used correct for the requested analyses? YES NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO

Were all VOC vials free of air bubbles? NA YES NO

Was sufficient amount of sample sent in each bottle? YES NO

Date VOC Trip Blank was made at ARI..... NA _____

Was Sample Split by ARI : NA YES Date/Time: _____ Equipment: _____ Split by: _____

Samples Logged by: JM Date: 3/17/11 Time: 945

**** Notify Project Manager of discrepancies or concerns ****

| Sample ID on Bottle | Sample ID on COC | Sample ID on Bottle | Sample ID on COC |
|---------------------|------------------|---------------------|------------------|
| | | | |
| | | | |
| | | | |

Additional Notes, Discrepancies, & Resolutions:

Samples LL-SED3-36-141-031511 + LL-SED3-141-167-031511 were split & preserved for subside.

By: JM Date: 3/17/11

| | | | |
|-----------------------------------|------------------------------|--|--|
| <p>Small Air Bubbles ~2mm</p> | <p>Peabubbles 2-4 mm</p> | <p>LARGE Air Bubbles > 4 mm</p> | <p>Small → "sm"</p> <p>Peabubbles → "pb"</p> <p>Large → "lg"</p> <p>Headspace → "hs"</p> |
|-----------------------------------|------------------------------|--|--|

Subject: Lora Lake samples

From: Erin Breckel <Erin.Breckel@floydsnider.com>

Date: Wed, 16 Mar 2011 16:56:01 -0700

To: 'Sue Dunnihoo' <sue@arilabs.com>

CC: Megan McCullough <Megan.McCullough@floydsnider.com>

Sue -

For both samples LL-SED3-36-141-031511 and LL-SED3-141-167-031511 can you please add a sulfide analysis to these samples?

I'm out of the office for the next two days, but if you have any questions regarding these samples you can talk to Megan McCullough here.

Thanks for your help!

Erin

Erin Breckel

FLOYD | SNIDER

Strategy * Science * Engineering

Two Union Square

601 Union Street, Suite 600

Seattle, WA 98101

tel: 206.292.2078 fax: 206.682.7867

www.floydsnider.com

Subject: Lora Lake samples
From: Erin Breckel <Erin.Breckel@floydsnider.com>
Date: Wed, 16 Mar 2011 11:43:07 -0700
To: 'Sue Dunnihoo' <sue@arilabs.com>

Hi Sue –

We dropped off our Lora Lake sediment samples this morning and had two samples listed as hold on the COC until we determined whether to analyze them or not. We just met and decided to go ahead and analyze them both. They are samples LL-SED3-36-141-031511 and LL-SED3-141-167-031511. Could you please let the lab know?

Thanks,
Erin

Erin Breckel
FLOYD | SNIDER
Strategy * Science * Engineering
Two Union Square
601 Union Street, Suite 600
Seattle, WA 98101
tel: 206.292.2078 fax: 206.682.7867
www.floydsnider.com

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: SN54



Case Narrative

Client: Floyd Snider
Project: Lora Lake Subsurface Sediment, POS-LL
ARI Job No.: SN54

Sample receipt

Analytical Resources, Inc. (ARI) accepted sediment samples on March 16, 2011 under ARI job SN54. The cooler temperature measured by IR thermometer following ARI SOP was 1.5°C. For details regarding sample receipt, please refer to the enclosed Cooler Receipt Form.

Samples **LL-SED3-0-36-031511**, **LL-SED2-0-56-031511**, **LL-SED2-0-56-031511-D**, and **LL-SED1-0-56-031511** were centrifuged prior to analysis.

Dioxin/Furan analyses were subcontracted to Frontier Analytical Laboratory in El Dorado Hills, CA. The dioxin data on CD as generated by Frontier is forwarded with this package.

SIM PAHs by SW8270D

The samples were initially screened to determine if a response was present that would require modifications in the extraction process. Based on the screen, the initial extraction weights were reduced for samples **LL-SED3-0-36-031511**, **LL-SED2-0-56-031511**, and **LL-SED2-0-56-031511-D**. The samples and associated laboratory QC were extracted and analyzed within the method recommended holding times.

Initial and continuing calibrations were within method requirements.

The internal standard areas of Naphthalene-d8 fell outside the control limits low for samples **LL-SED2-112-168-031511**, **LL-SED2-0-56-031511-D**, and **LL-SED1-0-56-031511**. The internal standard area of Phenanthrene-d10 fell outside the control limits low for sample **LL-SED1-0-56-031511**. These internal standards are not associated with requested compounds. No corrective action was taken.

The internal standard areas of Perylene-d12 fell outside the control limits low for samples **LL-SED2-112-168-031511**, **LL-SED2-0-56-031511-D**, and **LL-SED1-0-56-031511**. The samples were re-analyzed at dilutions and the internal standard areas were comparable to the initial analysis. No corrective action was taken.

The surrogate percent recoveries were within control limits.

The method blank was clean at the reporting limit.



The LCS percent recovery of Benzo(a)pyrene fell outside the control limits low for LCS-032611. All other percent recoveries were within control limits. No corrective action was taken.

The matrix spike and matrix spike duplicate percent recoveries were within advisory control limits.

In response to comments from NELAP and DOD auditors, ARI will now report the 'total' benzofluoranthenes rather than the individual compounds. This total will include the response of the b, k and j isomers.

Pentachlorophenol by SW8041

The samples and associated laboratory QC were extracted and analyzed within the method recommended holding times.

Initial and continuing calibrations were within method requirements.

The surrogate percent recoveries were within control limits.

The method blank was clean at the reporting limit. The LCS percent recovery was within control limits.

The matrix spike and matrix spike duplicate percent recoveries were within advisory control limits.

Total Arsenic and Lead by SW846 6010B

The samples and associated laboratory QC were digested and analyzed within the method recommended holding time.

The method blank was clean at the reporting limits. The LCS percent recoveries were within control limits.

The matrix spike percent recoveries and duplicate RPDs were within control limits.

General Chemistry

The samples and associated laboratory QC were prepared and analyzed within the method recommended holding time.

The method blanks were clean at the reporting limits. The LCS percent recoveries were within control limits.



The SRM percent recovery was within limits.

The matrix spike percent recovery of sulfide fell outside the control limits low for sample **LL-SED3-141-167-031511**. All other quality control parameters were met. No corrective action was taken.

The replicate RPD/RSDs were within limits.

Geotechnical Parameters

A laboratory-specific case narrative follows.



Analytical Resources, Incorporated

Analytical Chemists and Consultants

Client: Floyd-Snider

ARI Project No.: SN54

Client Project: Lora Lake – Subsurface Sediment

Client Project No.: POS-LLA

Case Narrative

1. Four samples were submitted for preparation on March 16, 2011, and were in good condition. Each sample was comprised of two, partially full 8oz jars.
2. The samples were submitted for separation of solids by means of centrifuging according to modified Corp of Engineers draft interim guide lines.
3. The samples were centrifuged in a pre-cooled centrifuge (4°C) at 1,000xg for 30 minutes.
4. After centrifuging of the sample, the supernatant liquid was decanted. The solids for each sample were combined, homogenized and spooned into a labeled 8 oz jar.
5. There were no other anomalies in the sample or methods on this project.

Approved by: *Guerra Suato*
Title: *Geotechnical Laboratory Manager*

Date: *3/22/11*

Sample ID Cross Reference Report



ARI Job No: SN54
Client: Floyd Snider
Project Event: POS-LL
Project Name: Lora Lake - Subsurface Sediment

| Sample ID | ARI Lab ID | ARI LIMS ID | Matrix | Sample Date/Time | VTSR |
|---------------------------|------------|-------------|----------|------------------|----------------|
| 1. LL-SED3-0-36-031511 | SN54A | 11-5925 | Sediment | 03/15/11 12:30 | 03/16/11 10:10 |
| 2. LL-SED3-36-141-031511 | SN54B | 11-5926 | Sediment | 03/15/11 12:40 | 03/16/11 10:10 |
| 3. LL-SED3-141-167-031511 | SN54C | 11-5927 | Sediment | 03/15/11 12:50 | 03/16/11 10:10 |
| 4. LL-SED2-0-56-031511 | SN54D | 11-5928 | Sediment | 03/15/11 14:15 | 03/16/11 10:10 |
| 5. LL-SED2-56-112-031511 | SN54E | 11-5929 | Sediment | 03/15/11 14:25 | 03/16/11 10:10 |
| 6. LL-SED2-112-168-031511 | SN54F | 11-5930 | Sediment | 03/15/11 14:55 | 03/16/11 10:10 |
| 7. LL-SED2-0-56-031511-D | SN54G | 11-5931 | Sediment | 03/15/11 14:15 | 03/16/11 10:10 |
| 8. LL-SED1-0-56-031511 | SN54H | 11-5932 | Sediment | 03/15/11 18:00 | 03/16/11 10:10 |

Printed 03/17/11



Data Reporting Qualifiers

Effective 2/14/2011

Inorganic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but \geq the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is ≤ 5 times the Reporting Limit and the replicate control limit defaults to ± 1 RL instead of the normal 20% RPD

Organic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Flagged value is not within established control limits
- B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
- J Estimated concentration when the value is less than ARI's established reporting limits
- D The spiked compound was not detected due to sample extract dilution
- E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte.
- Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria ($< 20\%$ RSD, $< 20\%$ Drift or minimum RRF).



- S Indicates an analyte response that has saturated the detector. The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- M2 The sample contains PCB congeners that do not match any standard Aroclor pattern. The PCBs are identified and quantified as the Aroclor whose pattern most closely matches that of the sample. The reported value is an estimate.
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit.
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by $\geq 40\%$ RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions. **(Dioxin/Furan analysis only)**



Geotechnical Data

- A The total of all fines fractions. This flag is used to report total fines when only sieve analysis is requested and balances total grain size with sample weight.

- F Samples were frozen prior to particle size determination

- SM Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations

- SS Sample did not contain the proportion of "fines" required to perform the pipette portion of the grain size analysis

- W Weight of sample in some pipette aliquots was below the level required for accurate weighting

LCS SOLUTIONS

| LABL SOLN ID | TEST | CONC. UG/ML | SOLVENT | EXP. |
|--------------|--------|--------------|------------|------------------|
| 1 | 1820-2 | PCB 1660 | 20 | ACETONE 01/25/12 |
| 2# | NA | BCOC PEST | 10 | ACETONE NA |
| 3 | 1793-3 | PEST | 01/02/10 | ACETONE 12/15/11 |
| 4 | 1806-2 | LOW PEST | .1/.2/1 | ACETONE 12/15/11 |
| 5 | 1779-1 | EPH | 1500 | MECL2 11/11/11 |
| 6 | 1791-5 | PCP | 12.5/125 | ACETONE 12/10/11 |
| 7 | 1834-4 | ABN | 100 | MEOH 08/21/11 |
| 8 | 1785-3 | TBT | 2.5 | MECL2 11/27/11 |
| 9 | 1786-3 | PORE TBT | .125/.25 | MECL2 11/27/11 |
| 10 | 1790-1 | ABN ACID | 100/200 | MEOH 06/07/11 |
| 11 | 1777-2 | TPHD | 15000 | ACETONE 11/01/11 |
| 12 | 1790-2 | ABN BASE | 200 | MEOH 06/07/11 |
| 13 | 1716-2 | LOW PCB | 2 | ACETONE 03/30/11 |
| 14 | 1822-2 | LOW ABN ACID | 10/20 | MEOH 06/07/11 |
| 15 | 1814-2 | SIM PNA | 15/75 | MEOH 01/04/12 |
| 16 | 1834-5 | 1,4-DIOXANE | 100 | MEOH 08/25/11 |
| 17 | 1772-3 | 1248 PCB | 10 | ACETONE 05/01/11 |
| 18 | 1814-3 | LOW SIM PNA | 1.5 | ACETONE 01/04/12 |
| 19 | 1815-2 | AK103 | 7500 | ACETONE 06/02/11 |
| 20 | 1775-3 | PNA | 100 | ACETONE 08/14/11 |
| 21 | 1725-1 | SKY/BHT | 100 | MEOH 03/18/11 |
| 22 | 1781-1 | HERB | 05 to 4000 | MEOH 04/15/11 |
| 23 | 1822-3 | LW ABN BASE | 20 | MEOH 06/07/11 |
| 24 | 1822-4 | LOW ABN | 10 | ACETONE 10/01/11 |
| 25# | NA | DIPHENYL | 100 | MEOH NA |
| 26 | 1823-1 | OP-PEST | 25 | MEOH 07/01/11 |
| 27 | NA | STEROLS | 200 | MEOH NA |
| 28# | 1807-1 | ADD. PEST | 2 | ACETONE 08/31/11 |
| 29# | NA | DECANES | 100 | MEOH NA |

LCS SOLUTIONS

| | | | | | |
|----|--------|-----------------------------|--------|---------|----------|
| 30 | NA | EDB/DBCP | 0.2 | MEOH | NA |
| 31 | 1835-2 | TERPINEOL | 100 | MEOH | 09/02/11 |
| 32 | NA | GUAIACOL | 50-200 | ACETONE | NA |
| 33 | NA | RETENE | 100 | MEOH | NA |
| 34 | NA | CONGENERS | 2.5 | ACETONE | NA |
| 35 | NA | ALKYL PNA A | 10 | MEOH | NA |
| 36 | NA | ALKYL PNA B | 10 | MEOH | NA |
| 37 | 1773-1 | CAR/PERY | 100 | ACETONE | 10/14/11 |
| 50 | 1757-4 | FULL RESIN | 250 | ACETONE | 08/14/11 |
| 51 | 1772-1 | DDTS | 0.01 | ACETONE | 04/24/11 |
| 52 | NA | 1232 PCB | 20 | ACETONE | NA |
| 53 | 1780-1 | DALAPON | 50 | MEOH | 05/07/11 |
| 54 | 1753-1 | T-CHLORDANE | 10 | ACETONE | 07/21/11 |
| 55 | 1753-2 | TOXAPHENE | 50 | ACETONE | 07/21/11 |
| | | | | | |
| | | #=PROJECT SPECIFIC SOLUTION | | | |
| | | *=REVERIFIED SOLUTION | | | |
| | | | | | |
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SURRE SOLUTIONS

| LABEL | SOLN ID | TEST | CONC. UG/ML | SOLVENT | EXP. |
|----------------------|---------|------------|-------------|---------|----------|
| A | 1824-2 | ABN | 100/150 | MEOH | 07/22/11 |
| B | 1834-6 | SIM PNA | 15/75 | ACETONE | 10/05/11 |
| C | 1705-4 | SIM ABN | 25/37.5 | MEOH | 03/08/11 |
| D | 1795-4 | LOW PCB | 0.2 | ACETONE | 12/16/11 |
| E | 1771-3 | HERB | 62.5 | MEOH | 10/06/11 |
| F | 1791-3 | PCP | 12.5 | ACETONE | 12/09/11 |
| G | 1824-1 | d8-DIOXANE | 100 | MEOH | 08/14/11 |
| H | 1723-2 | OP-PEST | 25 | MEOH | 04/02/11 |
| I | 1835-1 | LOW S. PNA | 1.5 | ACETONE | 10/05/11 |
| J | 1787-2 | TBT-PORE | 0.125 | MECL2 | 11/27/11 |
| K | 1795-2 | MED PCB | 20 | ACETONE | 12/16/11 |
| L | 1785-4 | TBT | 2.5 | MECL2 | 11/27/11 |
| M | 1767-1 | EPH | 1500 | MECL2 | 06/02/11 |
| N | 1795-3 | PCB | 2 | ACETONE | 12/16/11 |
| O | 1821-3 | TPH | 450 | MECL2 | 09/07/11 |
| P | 1813-2 | HCID | 2250 | MECL2 | 08/05/11 |
| Q | NA | EDB | 1 | MEOH | NA |
| R | 1757-3 | RESIN ACID | 250 | ACETONE | 08/14/11 |
| S* | NA | PBDE | .25 | MEOH | NA |
| T | 1768-2 | ALKYL PNA | 10 | MEOH | 07/22/11 |
| U | NA | CONGENER | 2.5 | ACETONE | NA |
| V | 1791-4 | LOW PCP | 1.25 | ACETONE | 12/09/11 |
| *reverified solution | | | | | |
| | | | | | |
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**Spike Recovery Control Limits for Polycyclic Aromatic Hydrocarbons
Selected Ion Monitoring (SIM) EPA Method SW-846-8270D-Modified ^(1,7)**
Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

| Sample Matrix | Water | | Soil | |
|--|-------------------------|--------------------------|-------------------------|--------------------------|
| Sample Volume / Final Volume | 500 mL to 0.5 mL | | 7.5 g / 0.5 mL | |
| | Control Limits | ME Limits ⁽²⁾ | Control Limits | ME Limits ⁽²⁾ |
| LCS Spike Recovery ⁽⁶⁾ | | | | |
| Napthalene | 39 - 100 | 30 - 102 | 37 - 100 | 27 - 107 |
| 2-Methylnapthalene | 39 - 100 | 31 - 100 | 37 - 100 | 28 - 100 |
| 1-Methylnapthalene | 30 - 160 ⁽³⁾ | 30 - 160 ⁽³⁾ | 30 - 160 ⁽³⁾ | 30 - 160 ⁽³⁾ |
| Acenaphthylene | 37 - 100 | 27 - 111 | 35 - 100 | 26 - 102 |
| Acenaphthene | 42 - 100 | 33 - 107 | 39 - 100 | 31 - 100 |
| Dibenzofuran | 46 - 100 | 38 - 101 | 39 - 100 | 31 - 100 |
| Fluorene | 49 - 101 | 40 - 110 | 42 - 100 | 33 - 106 |
| Phenanthrene | 55 - 101 | 47 - 109 | 47 - 100 | 38 - 108 |
| Anthracene | 47 - 102 | 38 - 111 | 41 - 106 | 30 - 117 |
| Fluoranthene | 60 - 106 | 52 - 114 | 52 - 109 | 43 - 119 |
| Pyrene | 55 - 110 | 46 - 119 | 47 - 111 | 36 - 122 |
| Benz(a)anthracene | 56 - 104 | 48 - 112 | 47 - 114 | 36 - 125 |
| Chrysene | 58 - 104 | 50 - 112 | 51 - 106 | 42 - 115 |
| Benzofluoranthene(s) (Total) | 30 - 160 ⁽⁸⁾ | 30 - 160 ⁽⁸⁾ | 30 - 160 ⁽⁸⁾ | 30 - 160 ⁽⁸⁾ |
| Benzo(a)pyrene | 32 - 110 | 19 - 123 | 44 - 111 | 33 - 122 |
| Indeno(1,2,3-cd)pyrene | 50 - 114 | 39 - 125 | 41 - 114 | 29 - 126 |
| Dibenzo(a,h)anthracene | 42 - 121 | 29 - 134 | 42 - 116 | 30 - 128 |
| Benzo(g,h,i)perylene | 50 - 113 | 40 - 124 | 37 - 115 | 27 - 107 |
| | | | | |
| MB / LCS Surrogate Recovery | | | | |
| d10-2-Methylnaphthalene | 36 - 101 | (4) | 35 - 100 | (4) |
| d14-Dibenzo(a,h)anthracene | 42 - 121 | (4) | 37 - 120 | (4) |
| | | | | |
| Sample Surrogate Recovery | | | | |
| d10-2-Methylnaphthalene | 30 - 106 | (4) | 34 - 100 | (4) |
| d14-Dibenzo(a,h)anthracene | 10 - 130 | (4) | 10 - 117 | (4) |

(1) ARI's Control limits calculated using all available spike recovery data from 1/1/08 through 12/31/08.

(2) **ME = A marginal exceedance** defined in the NELAC Standard ⁽⁵⁾ as beyond the LCS-CL but still within the ME limits. ME limits are between 3 and 4 standard deviations around the mean. A maximum of one marginal exceedance is acceptable. Two or more marginal exceedances require corrective action.

(3) 30 - 160 are default, advisory control limits used when there is insufficient data to calculate historic control limits. **DO NOT** use these limits as the sole reason to reject the data from a batch of analyses.

(4) Marginal Exceedances not allowed for surrogate standards.

(5) **2003 NELAC Standard (EPA/600/R-04/003), July 2003**, Chapter 5, pages 251-252.

(6) Laboratory Control Sample (LCS) spike recovery control limits also used as advisory control limits for sample matrix spike (MS) analyzes. MS recovery values are advisory and not used to assess the acceptability of an analytical batch.

(7) Highlighted control limits (**bold font**) adjusted to demonstrate that ARI does not use control limits < 10 for the lower limit or < 100 for the upper limit.

(8) Default limits pending generation of historic limits for total benzofluoranthrenes (7/29/10)



Spike Recovery Control Limits for Chlorinated Phenols
EPA Method SW-846-8041^(1,2)
Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

| | ARI's Calculated Control Limits | |
|--|---------------------------------|-----------------|
| | Water | Soil / Sediment |
| Sample Matrix: | Water | Soil / Sediment |
| Sample Amount / Final Volume: | 500 / 50 mL | 10 g / 25 mL |
| LCS Spike Recovery⁽³⁾ | | |
| Pentachlorophenol | 27 - 115 | 10 - 162 |
| | | |
| Method Blank/LCS Surrogate Recovery | | |
| 2,4,6-Tribromophenol | 40 - 130 | 50 - 115 |
| | | |
| Sample Surrogate Recovery | | |
| 2,4,6-Tribromophenol | 11 - 156 | 10 - 146 |
| | | |

(1) ARI's Control limits calculated using all available spike recovery data from 1/1/08 through 12/1/08.

(2) Highlighted control limits (**bold font**) adjusted to demonstrate that ARI does not use control limits < 10.

(3) Laboratory Control Sample (LCS) spike recovery control limits also used as advisory control limits for sample matrix spike (MS) analyzes. MS recovery values are advisory and not used to assess the acceptability of an analytical batch.



Summary of Laboratory Control Limits Metals Analyses (All Methods & Sample Matrices)

Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

| Element | Matrix Spike Recovery | LCS Recovery | Replicate RPD |
|-----------|-----------------------|--------------|---------------|
| Aluminum | 75 - 125 | 80 - 120 | ≤ 20% |
| Antimony | 75 - 125 | 80 - 120 | ≤ 20% |
| Arsenic | 75 - 125 | 80 - 120 | ≤ 20% |
| Barium | 75 - 125 | 80 - 120 | ≤ 20% |
| Beryllium | 75 - 125 | 80 - 120 | ≤ 20% |
| Boron | 75 - 125 | 80 - 120 | ≤ 20% |
| Cadmium | 75 - 125 | 80 - 120 | ≤ 20% |
| Calcium | 75 - 125 | 80 - 120 | ≤ 20% |
| Chromium | 75 - 125 | 80 - 120 | ≤ 20% |
| Cobalt | 75 - 125 | 80 - 120 | ≤ 20% |
| Copper | 75 - 125 | 80 - 120 | ≤ 20% |
| Iron | 75 - 125 | 80 - 120 | ≤ 20% |
| Lead | 75 - 125 | 80 - 120 | ≤ 20% |
| Magnesium | 75 - 125 | 80 - 120 | ≤ 20% |
| Manganese | 75 - 125 | 80 - 120 | ≤ 20% |
| Mercury | 75 - 125 | 80 - 120 | ≤ 20% |
| Nickel | 75 - 125 | 80 - 120 | ≤ 20% |
| Potassium | 75 - 125 | 80 - 120 | ≤ 20% |
| Selenium | 75 - 125 | 80 - 120 | ≤ 20% |
| Silica | 75 - 125 | 80 - 120 | ≤ 20% |
| Silver | 75 - 125 | 80 - 120 | ≤ 20% |
| Sodium | 75 - 125 | 80 - 120 | ≤ 20% |
| Strontium | 75 - 125 | 80 - 120 | ≤ 20% |
| Thallium | 75 - 125 | 80 - 120 | ≤ 20% |
| Vanadium | 75 - 125 | 80 - 120 | ≤ 20% |
| Zinc | 75 - 125 | 80 - 120 | ≤ 20% |



| Spike Recovery Control Limits for Conventional Wet Chemistry | | |
|---|-----------------------------|-----------------|
| Effective 5/1/09 | | |
| Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. http://www.arilabs.com/portal/downloads/ARI-CLs.zip | | |
| | ARI's Control Limits | |
| Sample Matrix: | Water | Soil / Sediment |
| Matrix Spike Recoveries | % Recovery | % Recovery |
| Ammonia | 75 - 125 | 75 - 125 |
| Bromide | 75 - 125 | 75 - 125 |
| Chloride | 75 - 125 | 75 - 125 |
| Cyanide | 75 - 125 | 75 - 125 |
| Ferrous Iron | 75 - 125 | 75 - 125 |
| Fluoride | 75 - 125 | 75 - 125 |
| Formaldehyde | 75 - 125 | 75 - 125 |
| Hexane Extractable Material | -- - -- | 78 - 114 |
| Hexavalent Chromium | 75 - 125 | 75 - 125 |
| Nitrate/Nitrite | 75 - 125 | 75 - 125 |
| Oil and Grease | 75 - 125 | 75 - 125 |
| Phenol | 75 - 125 | 75 - 125 |
| Phosphorous | 75 - 125 | 75 - 125 |
| Sulfate | 75 - 125 | 75 - 125 |
| Sulfide | 75 - 125 | 75 - 125 |
| Total Kjeldahl Nitrogen | 75 - 125 | 75 - 125 |
| Total Organic Carbon | 75 - 125 | 75 - 125 |
| Duplicate RPDs | | |
| Acidity | ±20% | ±20% |
| Alkalinity | ±20% | ±20% |
| BOD | ±20% | ±20% |
| Cation Exchange | ±20% | ±20% |
| COD | ±20% | ±20% |
| Conductivity | ±20% | ±20% |
| Salinity | ±20% | ±20% |
| Solids | ±20% | ±20% |
| Turbidity | ±20% | ±20% |

**SIM PAH Analysis
Report and Summary QC Forms**

ARI Job ID: SN54

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
Page 1 of 1

Sample ID: LL-SED3-0-36-031511
SAMPLE

Lab Sample ID: SN54A
LIMS ID: 11-5925
Matrix: Sediment
Data Release Authorized: *MM*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/01/11 18:13
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 5.92 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 1.00
Percent Moisture: 75.4%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|-----|---------|
| 56-55-3 | Benzo (a) anthracene | 8.4 | 62 |
| 218-01-9 | Chrysene | 8.4 | 140 |
| 50-32-8 | Benzo (a) pyrene | 8.4 | 88 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 8.4 | 68 |
| 53-70-3 | Dibenz (a,h) anthracene | 8.4 | < 8.4 U |
| TOTBFA | Total Benzofluoranthenes | 8.4 | 180 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 64.0%
d14-Dibenzo (a,h) anthracen 57.7%

Sample ID: LL-SED3-36-141-031511
SAMPLE

Lab Sample ID: SN54B
LIMS ID: 11-5926
Matrix: Sediment
Data Release Authorized: *MM*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/01/11 18:41
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.07 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 88.3%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | < 15 U |
| 218-01-9 | Chrysene | 15 | 18 |
| 50-32-8 | Benzo(a)pyrene | 15 | < 15 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | < 15 U |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | < 15 U |
| TOTBFA | Total Benzofluoranthenes | 15 | < 15 U |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 74.0%
d14-Dibenzo(a,h)anthracen 55.0%

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED3-141-167-031511
SAMPLE

Lab Sample ID: SN54C
LIMS ID: 11-5927
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/01/11 19:09
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.06 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 87.9%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | < 15 U |
| 218-01-9 | Chrysene | 15 | < 15 U |
| 50-32-8 | Benzo(a)pyrene | 15 | < 15 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | < 15 U |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | < 15 U |
| TOTBFA | Total Benzofluoranthenes | 15 | < 15 U |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 65.0%
d14-Dibenzo(a,h)anthracen 51.0%

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED2-0-56-031511
SAMPLE

Lab Sample ID: SN54D
LIMS ID: 11-5928
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/01/11 19:37
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 3.29 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 1.00
Percent Moisture: 85.9%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo (a) anthracene | 15 | 130 |
| 218-01-9 | Chrysene | 15 | 300 |
| 50-32-8 | Benzo (a) pyrene | 15 | 170 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 15 | 110 |
| 53-70-3 | Dibenz (a,h) anthracene | 15 | 30 |
| TOTBFA | Total Benzofluoranthenes | 15 | 470 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 70.3%
d14-Dibenzo(a,h)anthracen 53.7%

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED2-56-112-031511
SAMPLE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 11:59
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.15 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | < 15 U |
| 218-01-9 | Chrysene | 15 | < 15 U |
| 50-32-8 | Benzo(a)pyrene | 15 | < 15 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | < 15 U |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | < 15 U |
| TOTBFA | Total Benzofluoranthenes | 15 | < 15 U |

Reported in µg/kg (ppb)

SIM Semivolatle Surrogate Recovery

d10-2-Methylnaphthalene 72.0%
d14-Dibenzo(a,h)anthracen 55.0%

ORGANICS ANALYSIS DATA SHEET
 PNAs by SIM SW8270D-SIM GC/MS
 Page 1 of 1

Sample ID: LL-SED2-112-168-031511
 SAMPLE

Lab Sample ID: SN54F
 LIMS ID: 11-5930
 Matrix: Sediment
 Data Release Authorized: *MMW*
 Reported: 04/06/11

QC Report No: SN54-Floyd Snider
 Project: Lora Lake - Subsurface Sediment
 Event: POS-LL
 Date Sampled: 03/15/11
 Date Received: 03/16/11

Date Extracted: 03/26/11
 Date Analyzed: 04/04/11 15:08
 Instrument/Analyst: NT12/JZ
 GPC Cleanup: No
 Silica Gel Cleanup: Yes
 Alumina Cleanup: No

Sample Amount: 10.09 g-dry-wt
 Final Extract Volume: 0.5 mL
 Dilution Factor: 3.00
 Percent Moisture: 87.7%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | < 15 U |
| 218-01-9 | Chrysene | 15 | < 15 U |
| 50-32-8 | Benzo(a)pyrene | 15 | < 15 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | < 15 U |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | < 15 U |
| TOTBFA | Total Benzofluoranthenes | 15 | < 15 U |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

| | |
|---------------------------|-------|
| d10-2-Methylnaphthalene | 84.0% |
| d14-Dibenzo(a,h)anthracen | 64.0% |

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED2-112-168-031511
DILUTION

Lab Sample ID: SN54F
LIMS ID: 11-5930
Matrix: Sediment
Data Release Authorized: *mm*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 19:24
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.09 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 30.0
Percent Moisture: 87.7%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|-----|---------|
| 56-55-3 | Benzo(a)anthracene | 150 | < 150 U |
| 218-01-9 | Chrysene | 150 | < 150 U |
| 50-32-8 | Benzo(a)pyrene | 150 | < 150 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 150 | < 150 U |
| 53-70-3 | Dibenz(a,h)anthracene | 150 | < 150 U |
| TOTBFA | Total Benzofluoranthenes | 150 | < 150 U |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

| | |
|---------------------------|-------|
| d10-2-Methylnaphthalene | 90.0% |
| d14-Dibenzo(a,h)anthracen | 60.0% |

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED2-0-56-031511-D
SAMPLE

Lab Sample ID: SN54G
LIMS ID: 11-5931
Matrix: Sediment
Data Release Authorized: *MM*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 15:36
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 2.15 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 1.00
Percent Moisture: 88.7%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo (a) anthracene | 23 | 270 |
| 218-01-9 | Chrysene | 23 | 620 |
| 50-32-8 | Benzo (a) pyrene | 23 | 400 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 23 | 280 |
| 53-70-3 | Dibenz (a,h) anthracene | 23 | 74 |
| TOTBFA | Total Benzofluoranthenes | 23 | 1,100 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 75.0%
d14-Dibenzo (a,h) anthracen 59.0%

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
Page 1 of 1

Sample ID: LL-SED2-0-56-031511-D
DILUTION

Lab Sample ID: SN54G
LIMS ID: 11-5931
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 17:22
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 2.15 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 88.7%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo (a) anthracene | 70 | 240 |
| 218-01-9 | Chrysene | 70 | 570 |
| 50-32-8 | Benzo (a) pyrene | 70 | 330 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 70 | 270 |
| 53-70-3 | Dibenz (a,h) anthracene | 70 | 96 |
| TOTBFA | Total Benzofluoranthenes | 70 | 1,000 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

| | |
|-----------------------------|-------|
| d10-2-Methylnaphthalene | 69.0% |
| d14-Dibenzo (a,h) anthracen | 59.0% |

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED1-0-56-031511
SAMPLE

Lab Sample ID: SN54H
LIMS ID: 11-5932
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 16:03
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.56 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 1.00
Percent Moisture: .45.6%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|-----|--------|
| 56-55-3 | Benzo (a) anthracene | 4.7 | 37 |
| 218-01-9 | Chrysene | 4.7 | 81 |
| 50-32-8 | Benzo (a) pyrene | 4.7 | 55 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 4.7 | 38 |
| 53-70-3 | Dibenz (a,h) anthracene | 4.7 | 5.8 |
| TOTBFA | Total Benzofluoranthenes | 4.7 | 140 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

| | |
|-----------------------------|-------|
| d10-2-Methylnaphthalene | 72.7% |
| d14-Dibenzo (a,h) anthracen | 62.0% |

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
Page 1 of 1

Sample ID: LL-SED1-0-56-031511
DILUTION

Lab Sample ID: SN54H
LIMS ID: 11-5932
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 17:50
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.56 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 45.6%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo (a) anthracene | 14 | 31 |
| 218-01-9 | Chrysene | 14 | 70 |
| 50-32-8 | Benzo (a) pyrene | 14 | 46 |
| 193-39-5 | Indeno (1,2,3-cd) pyrene | 14 | 34 |
| 53-70-3 | Dibenz (a,h) anthracene | 14 | < 14 U |
| TOTBEA | Total Benzofluoranthenes | 14 | 130 |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 62.0%
d14-Dibenzo (a,h) anthracen 58.0%

SIM SW8270 SURROGATE RECOVERY SUMMARY

Matrix: Sediment

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL

| <u>Client ID</u> | <u>MNP</u> | <u>DBA</u> | <u>TOT OUT</u> |
|---------------------------|------------|------------|----------------|
| LL-SED3-0-36-031511 | 64.0% | 57.7% | 0 |
| LL-SED3-36-141-031511 | 74.0% | 55.0% | 0 |
| LL-SED3-141-167-031511 | 65.0% | 51.0% | 0 |
| LL-SED2-0-56-031511 | 70.3% | 53.7% | 0 |
| MB-032611 | 63.0% | 74.3% | 0 |
| LCS-032611 | 58.3% | 75.3% | 0 |
| LCSD-032611 | 59.0% | 70.3% | 0 |
| LL-SED2-56-112-031511 | 72.0% | 55.0% | 0 |
| LL-SED2-56-112-031511 MS | 67.0% | 54.0% | 0 |
| LL-SED2-56-112-031511 MSD | 83.0% | 57.0% | 0 |
| LL-SED2-112-168-031511 | 84.0% | 64.0% | 0 |
| LL-SED2-112-168-031511 DL | 90.0% | 60.0% | 0 |
| LL-SED2-0-56-031511-D | 75.0% | 59.0% | 0 |
| LL-SED2-0-56-031511-D DL | 69.0% | 59.0% | 0 |
| LL-SED1-0-56-031511 | 72.7% | 62.0% | 0 |
| LL-SED1-0-56-031511 DL | 62.0% | 58.0% | 0 |

LCS/MB LIMITS QC LIMITS

(MNP) = d10-2-Methylnaphthalene (35-100) (34-100)
(DBA) = d14-Dibenzo(a,h)anthracene (37-120) (10-117)

Prep Method: SW3550C
Log Number Range: 11-5925 to 11-5932

ORGANICS ANALYSIS DATA SHEET
PNAs by SW8270D-SIM GC/MS
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Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted MS/MSD: 03/26/11
Date Analyzed MS: 04/04/11 12:37
MSD: 04/04/11 13:05
Instrument/Analyst MS: NT12/JZ
MSD: NT12/JZ

Sample Amount MS: 10.1 g-dry-wt
MSD: 10.1 g-dry-wt
Final Extract Volume MS: 0.50 mL
MSD: 0.50 mL
Dilution Factor MS: 3.00
MSD: 3.00

| Analyte | Sample | MS | Spike | | MSD | Spike | | RPD |
|--------------------------|----------|------|----------|----------|------|-----------|----------|-------|
| | | | Added-MS | Recovery | | Added-MSD | Recovery | |
| Benzo(a)anthracene | < 14.8 U | 95.0 | 149 | 63.8% | 118 | 148 | 79.7% | 21.6% |
| Chrysene | < 14.8 U | 97.5 | 149 | 65.4% | 114 | 148 | 77.0% | 15.6% |
| Benzo(a)pyrene | < 14.8 U | 89.9 | 149 | 60.3% | 116 | 148 | 78.4% | 25.4% |
| Indeno(1,2,3-cd)pyrene | < 14.8 U | 81.6 | 149 | 54.8% | 86.5 | 148 | 58.4% | 5.8% |
| Dibenz(a,h)anthracene | < 14.8 U | 88.4 | 149 | 59.3% | 93.6 | 148 | 63.2% | 5.7% |
| Total Benzofluoranthenes | < 14.8 U | 196 | 298 | 65.8% | 225 | 297 | 75.8% | 13.8% |

Reported in µg/kg (ppb)

RPD calculated using sample concentrations per SW846.

Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 12:37
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.08 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | --- |
| 218-01-9 | Chrysene | 15 | --- |
| 50-32-8 | Benzo(a)pyrene | 15 | --- |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | --- |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | --- |
| TOTBFA | Total Benzofluoranthenes | 15 | --- |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 67.0%
d14-Dibenzo(a,h)anthracen 54.0%

ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE DUPLICATE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/26/11
Date Analyzed: 04/04/11 13:05
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.11 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 3.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|----|--------|
| 56-55-3 | Benzo(a)anthracene | 15 | --- |
| 218-01-9 | Chrysene | 15 | --- |
| 50-32-8 | Benzo(a)pyrene | 15 | --- |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 15 | --- |
| 53-70-3 | Dibenz(a,h)anthracene | 15 | --- |
| TOTBFA | Total Benzofluoranthenes | 15 | --- |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

| | |
|---------------------------|-------|
| d10-2-Methylnaphthalene | 83.0% |
| d14-Dibenzo(a,h)anthracen | 57.0% |

ORGANICS ANALYSIS DATA SHEET
PNAs by SW8270D-SIM GC/MS
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Sample ID: LCS-032611
LAB CONTROL SAMPLE

Lab Sample ID: LCS-032611
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: NA
Date Received: NA

Date Extracted: 03/26/11

Sample Amount LCS: 10.0 g-dry-wt
LCSD: 10.0 g-dry-wt

Date Analyzed LCS: 04/01/11 16:49
LCSD: 04/01/11 17:17
Instrument/Analyst LCS: NT12/JZ
LCSD: NT12/JZ

Final Extract Volume LCS: 0.50 mL
LCSD: 0.50 mL
Dilution Factor LCS: 1.00
LCSD: 1.00

| Analyte | LCS | Spike | | LCS | LCSD | Spike | | RPD |
|--------------------------|------|-----------|----------|------|------|------------|----------|-----|
| | | Added-LCS | Recovery | | | Added-LCSD | Recovery | |
| Benzo(a)anthracene | 108 | 150 | 72.0% | 104 | 150 | 69.3% | 3.8% | |
| Chrysene | 122 | 150 | 81.3% | 123 | 150 | 82.0% | 0.8% | |
| Benzo(a)pyrene | 70.6 | 150 | 47.1% | 62.2 | 150 | 41.5% | 12.7% | |
| Indeno(1,2,3-cd)pyrene | 108 | 150 | 72.0% | 108 | 150 | 72.0% | 0.0% | |
| Dibenz(a,h)anthracene | 116 | 150 | 77.3% | 112 | 150 | 74.7% | 3.5% | |
| Total Benzofluoranthenes | 242 | 300 | 80.7% | 234 | 300 | 78.0% | 3.4% | |

Reported in µg/kg (ppb)

RPD calculated using sample concentrations per SW846.

SIM Semivolatile Surrogate Recovery

| | LCS | LCSD |
|---------------------------|-------|-------|
| d10-2-Methylnaphthalene | 58.3% | 59.0% |
| d14-Dibenzo(a,h)anthracen | 75.3% | 70.3% |

4B
SEMIVOLATILE METHOD BLANK SUMMARY

BLANK NO.

| |
|----------|
| SN54MBS2 |
|----------|

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE - SUBSURFA

Lab File ID: 04011104

Date Extracted: 03/26/11

Instrument ID: NT12

Date Analyzed: 04/01/11

Matrix: SOLID

Time Analyzed: 1621

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

| | CLIENT SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED |
|----|----------------------|------------------|----------------|------------------|
| | ===== | ===== | ===== | ===== |
| 01 | SN54LCSS2 | SN54LCSS2 | 04011105 | 04/01/11 |
| 02 | SN54LCSDS2 | SN54LCSDS2 | 04011106 | 04/01/11 |
| 03 | LL-SED3-0-36-031 | SN54A | 04011108 | 04/01/11 |
| 04 | LL-SED3-36-141-0 | SN54B | 04011109 | 04/01/11 |
| 05 | LL-SED3-141-167- | SN54C | 04011110 | 04/01/11 |
| 06 | LL-SED2-0-56-031 | SN54D | 04011111 | 04/01/11 |
| 07 | LL-SED2-56-112-0 | SN54E | 04041103 | 04/04/11 |
| 08 | LL-SED2-56-112- | SN54EMS | 04041104 | 04/04/11 |
| 09 | LL-SED2-56-112- | SN54EMSD | 04041105 | 04/04/11 |
| 10 | LL-SED2-112-168- | SN54F | 04041108 | 04/04/11 |
| 11 | LL-SED2-0-56-031 | SN54G | 04041109 | 04/04/11 |
| 12 | LL-SED1-0-56-031 | SN54H | 04041110 | 04/04/11 |
| 13 | LL-SED2-0-56-031 | SN54G | 04041112 | 04/04/11 |
| 14 | LL-SED1-0-56-031 | SN54H | 04041113 | 04/04/11 |
| 15 | LL-SED2-112-168- | SN54F | 04041114 | 04/04/11 |
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ORGANICS ANALYSIS DATA SHEET
PNAs by SIM SW8270D-SIM GC/MS
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Sample ID: MB-032611
METHOD BLANK

Lab Sample ID: MB-032611
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/06/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
Event: POS-LL
Date Sampled: NA
Date Received: NA

Date Extracted: 03/26/11
Date Analyzed: 04/01/11 16:21
Instrument/Analyst: NT12/JZ
GPC Cleanup: No
Silica Gel Cleanup: Yes
Alumina Cleanup: No

Sample Amount: 10.00 g-dry-wt
Final Extract Volume: 0.5 mL
Dilution Factor: 1.00
Percent Moisture: NA

| CAS Number | Analyte | RL | Result |
|------------|--------------------------|-----|---------|
| 56-55-3 | Benzo(a)anthracene | 5.0 | < 5.0 U |
| 218-01-9 | Chrysene | 5.0 | < 5.0 U |
| 50-32-8 | Benzo(a)pyrene | 5.0 | < 5.0 U |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | 5.0 | < 5.0 U |
| 53-70-3 | Dibenz(a,h)anthracene | 5.0 | < 5.0 U |
| TOTBFA | Total Benzofluoranthenes | 5.0 | < 5.0 U |

Reported in µg/kg (ppb)

SIM Semivolatile Surrogate Recovery

d10-2-Methylnaphthalene 63.0%
d14-Dibenzo(a,h)anthracen 74.3%

5B
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

Instrument ID: NT12

Project: LORA LAKE

DFTPP Injection Date: 03/31/11

DFTPP Injection Time: 2014

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51 | 10.0 - 80.0% of mass 198 | 19.7 |
| 68 | Less than 2.0% of mass 69 | 0.0 (0.0)1 |
| 69 | Mass 69 relative abundance | 78.7 |
| 70 | Less than 2.0% of mass 69 | 0.4 (0.5)1 |
| 127 | 10.0 - 80.0% of mass 198 | 54.0 |
| 197 | Less than 2.0% of mass 198 | 0.0 |
| 198 | Base Peak, 100% relative abundance | 100.0 |
| 199 | 5.0 to 9.0% of mass 198 | 7.8 |
| 275 | 10.0 - 60.0% of mass 198 | 26.9 |
| 365 | Greater than 1.0% of mass 198 | 4.71 |
| 441 | 0.0 - 24.0% of mass 442 | 9.4 (16.8)2 |
| 442 | 50.0 - 200.0% of mass 198 | 55.8 |
| 443 | 15.0 - 24.0% of mass 442 | 11.1 (19.9)2 |

1-Value is % mass 69

2-Value is % mass 442

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

| | CLIENT SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------------|------------------|----------------|------------------|------------------|
| 01 | IC0250331 | IC0250331 | 03311102 | 03/31/11 | 2027 |
| 02 | IC010331 | IC010331 | 03311103 | 03/31/11 | 2055 |
| 03 | IC050331 | IC050331 | 03311104 | 03/31/11 | 2123 |
| 04 | IC10331 | IC10331 | 03311105 | 03/31/11 | 2151 |
| 05 | IC50331 | IC50331 | 03311106 | 03/31/11 | 2219 |
| 06 | IC10331 | IC100331 | 03311107 | 03/31/11 | 2247 |
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5B
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

Instrument ID: NT12

Project: LORA LAKE

DFTPP Injection Date: 04/01/11

DFTPP Injection Time: 1457

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51 | 10.0 - 80.0% of mass 198 | 21.0 |
| 68 | Less than 2.0% of mass 69 | 0.0 (0.0)1 |
| 69 | Mass 69 relative abundance | 78.0 |
| 70 | Less than 2.0% of mass 69 | 0.2 (0.3)1 |
| 127 | 10.0 - 80.0% of mass 198 | 54.2 |
| 197 | Less than 2.0% of mass 198 | 0.0 |
| 198 | Base Peak, 100% relative abundance | 100.0 |
| 199 | 5.0 to 9.0% of mass 198 | 8.1 |
| 275 | 10.0 - 60.0% of mass 198 | 29.2 |
| 365 | Greater than 1.0% of mass 198 | 5.20 |
| 441 | 0.0 - 24.0% of mass 442 | 10.3 (16.2)2 |
| 442 | 50.0 - 200.0% of mass 198 | 63.5 |
| 443 | 15.0 - 24.0% of mass 442 | 12.8 (20.2)2 |

1-Value is % mass 69

2-Value is % mass 442

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

| | CLIENT SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------------|------------------|----------------|------------------|------------------|
| 01 | CC0401 | CC0401 | 04011102 | 04/01/11 | 1516 |
| 02 | SN54MBS2 | SN54MBS2 | 04011104 | 04/01/11 | 1621 |
| 03 | SN54LCSS2 | SN54LCSS2 | 04011105 | 04/01/11 | 1649 |
| 04 | SN54LCSDS2 | SN54LCSDS2 | 04011106 | 04/01/11 | 1717 |
| 05 | LL-SED3-0-36-031 | SN54A | 04011108 | 04/01/11 | 1813 |
| 06 | LL-SED3-36-141-0 | SN54B | 04011109 | 04/01/11 | 1841 |
| 07 | LL-SED3-141-167- | SN54C | 04011110 | 04/01/11 | 1909 |
| 08 | LL-SED2-0-56-031 | SN54D | 04011111 | 04/01/11 | 1937 |
| 09 | | | | | |
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5B
SEMIVOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

Instrument ID: NT12

Project: LORA LAKE

DFTPP Injection Date: 04/04/11

DFTPP Injection Time: 0949

| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 51 | 10.0 - 80.0% of mass 198 | 23.4 |
| 68 | Less than 2.0% of mass 69 | 0.0 (0.0)1 |
| 69 | Mass 69 relative abundance | 89.0 |
| 70 | Less than 2.0% of mass 69 | 0.6 (0.7)1 |
| 127 | 10.0 - 80.0% of mass 198 | 55.8 |
| 197 | Less than 2.0% of mass 198 | 0.0 |
| 198 | Base Peak, 100% relative abundance | 100.0 |
| 199 | 5.0 to 9.0% of mass 198 | 7.7 |
| 275 | 10.0 - 60.0% of mass 198 | 28.4 |
| 365 | Greater than 1.0% of mass 198 | 5.30 |
| 441 | 0.0 - 24.0% of mass 442 | 9.5 (15.9)2 |
| 442 | 50.0 - 200.0% of mass 198 | 59.7 |
| 443 | 15.0 - 24.0% of mass 442 | 11.8 (19.8)2 |

1-Value is % mass 69

2-Value is % mass 442

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

| | CLIENT SAMPLE NO. | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|----|----------------------|------------------|----------------|------------------|------------------|
| 01 | CC0404 | CC0404 | 04041102 | 04/04/11 | 1118 |
| 02 | LL-SED2-56-112-0 | SN54E | 04041103 | 04/04/11 | 1159 |
| 03 | LL-SED2-56-112- | SN54EMS | 04041104 | 04/04/11 | 1237 |
| 04 | LL-SED2-56-112- | SN54EMSD | 04041105 | 04/04/11 | 1305 |
| 05 | LL-SED2-112-168- | SN54F | 04041108 | 04/04/11 | 1508 |
| 06 | LL-SED2-0-56-031 | SN54G | 04041109 | 04/04/11 | 1536 |
| 07 | LL-SED1-0-56-031 | SN54H | 04041110 | 04/04/11 | 1603 |
| 08 | LL-SED2-0-56-031 | SN54G | 04041112 | 04/04/11 | 1722 |
| 09 | LL-SED1-0-56-031 | SN54H | 04041113 | 04/04/11 | 1750 |
| 10 | LL-SED2-112-168- | SN54F | 04041114 | 04/04/11 | 1924 |
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SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Instrument ID: NT12

Cont. Calib. Date: 04/01/11

Init. Calib. Date: 03/31/11

Cont. Calib. Time: 1516

| COMPOUND | CalAmt or ARF | CC Amt or RF | MIN RRF | CURVE TYPE | %D or Drift |
|----------------------------|------------------|-----------------|------------|---------------|----------------|
| ===== | ===== | ===== | ===== | ===== | ===== |
| Naphthalene | 0.966 | 0.992 | 0.700 | AVRG | 2.7 |
| 2-Methylnaphthalene | 0.537 | 0.538 | 0.400 | AVRG | 0.2 |
| Acenaphthylene | 1.750 | 1.659 | 0.900 | AVRG | -5.2 |
| Acenaphthene | 1.090 | 1.040 | 0.900 | AVRG | -4.6 |
| Dibenzofuran | 1.482 | 1.436 | 0.800 | AVRG | -3.1 |
| Fluorene | 1.220 | 1.207 | 0.900 | AVRG | -1.1 |
| Phenanthrene | 1.056 | 1.064 | 0.700 | AVRG | 0.8 |
| Anthracene | 1.060 | 1.010 | 0.700 | AVRG | -4.7 |
| Fluoranthene | 1.148 | 1.130 | 0.600 | AVRG | -1.6 |
| Pyrene | 1.077 | 1.059 | 0.600 | AVRG | -1.7 |
| Benzo(a)anthracene | 0.962 | 0.936 | 0.800 | AVRG | -2.7 |
| Chrysene | 0.940 | 0.907 | 0.700 | AVRG | -3.5 |
| Benzo(b)fluoranthene | 1.128 | 1.129 | 0.700 | AVRG | 0.1 |
| Benzo(k)fluoranthene | 1.183 | 1.061 | 0.700 | AVRG | -10.3 |
| Benzo(j)fluoranthene | 1.130 | 1.063 | 0.010 | AVRG | -5.9 |
| Benzo(a)pyrene | 1.035 | 1.015 | 0.700 | AVRG | -1.9 |
| Indeno(1,2,3-cd)pyrene | 1.259 | 1.196 | 0.500 | AVRG | -5.0 |
| Dibenzo(a,h)anthracene | 1.029 | 0.996 | 0.400 | AVRG | -3.2 |
| Benzo(g,h,i)perylene | 1.116 | 1.041 | 0.500 | AVRG | -6.7 |
| 1-methylnaphthalene | 0.571 | 0.564 | 0.010 | AVRG | -1.2 |
| Perylene | 0.876 | 0.837 | 0.010 | AVRG | -4.4 |
| ===== | ===== | ===== | ===== | ===== | ===== |
| 2-Methylnaphthalene-d10 | 0.611 | 0.599 | 0.010 | AVRG | -2.0 |
| Dibenzo(a,h)anthracene-d14 | 0.938 | 0.879 | 0.010 | AVRG | -6.3 |

<- Exceeds QC limit of 20% D

* RF less than minimum RF

SEMIVOLATILE 8270-D CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Instrument ID: NT12

Cont. Calib. Date: 04/04/11

Init. Calib. Date: 03/31/11

Cont. Calib. Time: 1118

| COMPOUND | CalAmt or ARF | CC Amt or RF | MIN RRF | CURVE TYPE | %D or Drift |
|----------------------------|------------------|-----------------|------------|---------------|----------------|
| Naphthalene | 0.966 | 0.941 | 0.700 | AVRG | -2.6 |
| 2-Methylnaphthalene | 0.537 | 0.566 | 0.400 | AVRG | 5.4 |
| Acenaphthylene | 1.750 | 1.648 | 0.900 | AVRG | -5.8 |
| Acenaphthene | 1.090 | 0.998 | 0.900 | AVRG | -8.4 |
| Dibenzofuran | 1.482 | 1.356 | 0.800 | AVRG | -8.5 |
| Fluorene | 1.220 | 1.159 | 0.900 | AVRG | -5.0 |
| Phenanthrene | 1.056 | 0.994 | 0.700 | AVRG | -5.9 |
| Anthracene | 1.060 | 1.037 | 0.700 | AVRG | -2.2 |
| Fluoranthene | 1.148 | 1.074 | 0.600 | AVRG | -6.4 |
| Pyrene | 1.077 | 1.156 | 0.600 | AVRG | 7.3 |
| Benzo(a)anthracene | 0.962 | 0.962 | 0.800 | AVRG | 0.0 |
| Chrysene | 0.940 | 0.893 | 0.700 | AVRG | -5.0 |
| Benzo(b)fluoranthene | 1.128 | 1.172 | 0.700 | AVRG | 3.9 |
| Benzo(k)fluoranthene | 1.183 | 1.096 | 0.700 | AVRG | -7.4 |
| Benzo(j)fluoranthene | 1.130 | 1.090 | 0.010 | AVRG | -3.5 |
| Benzo(a)pyrene | 1.035 | 1.004 | 0.700 | AVRG | -3.0 |
| Indeno(1,2,3-cd)pyrene | 1.259 | 1.198 | 0.500 | AVRG | -4.8 |
| Dibenzo(a,h)anthracene | 1.029 | 1.008 | 0.400 | AVRG | -2.0 |
| Benzo(g,h,i)perylene | 1.116 | 1.053 | 0.500 | AVRG | -5.6 |
| 1-methylnaphthalene | 0.571 | 0.581 | 0.010 | AVRG | 1.8 |
| Perylene | 0.876 | 0.820 | 0.010 | AVRG | -6.4 |
| 2-Methylnaphthalene-d10 | 0.611 | 0.657 | 0.010 | AVRG | 7.5 |
| Dibenzo(a,h)anthracene-d14 | 0.938 | 0.903 | 0.010 | AVRG | -3.7 |

<- Exceeds QC limit of 20% D

* RF less than minimum RF

8B
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Ical Midpoint ID: 03311102

Ical Date: 03/31/11

Instrument ID: NT12

Cont. Cal Date: 04/01/11

| | IS1 (NPT) AREA # | RT # | IS2 (ANT) AREA # | RT # | IS3 (PHN) AREA # | RT # |
|-----------------|---------------------|-------|---------------------|-------|---------------------|-------|
| ===== | ===== | ===== | ===== | ===== | ===== | ===== |
| ICAL MIDPT | 494112 | 4.92 | 280105 | 7.16 | 461353 | 9.10 |
| UPPER LIMIT | 988224 | | 560210 | | 922706 | |
| LOWER LIMIT | 247056 | | 140052 | | 230676 | |
| ===== | ===== | ===== | ===== | ===== | ===== | ===== |
| CCAL | 499536 | 4.85 | 294599 | 7.09 | 482436 | 9.03 |
| UPPER LIMIT | | 5.35 | | 7.59 | | 9.53 |
| LOWER LIMIT | | 4.35 | | 6.59 | | 8.53 |
| 01 SN54MBS2 | 362184 | 4.84 | 207575 | 7.08 | 332402 | 9.02 |
| 02 SN54LCSS2 | 366499 | 4.84 | 217644 | 7.08 | 347080 | 9.02 |
| 03 SN54LCSDS2 | 364394 | 4.84 | 218342 | 7.08 | 356245 | 9.02 |
| 04 LL-SED3-0-36 | 342514 | 4.84 | 202181 | 7.09 | 322755 | 9.03 |
| 05 LL-SED3-36-1 | 301159 | 4.84 | 187766 | 7.09 | 301621 | 9.02 |
| 06 LL-SED3-141- | 262380 | 4.85 | 167023 | 7.09 | 271187 | 9.03 |
| 07 LL-SED2-0-56 | 249435 | 4.84 | 163363 | 7.09 | 258445 | 9.03 |
| 08 | | | | | | |
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IS1 = Naphthalene-d8
IS2 = Acenaphthene-d10
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

* Values outside of QC limits.

8B
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Ical Midpoint ID: 03311102

Ical Date: 03/31/11

Instrument ID: NT12

Cont. Cal Date: 04/01/11

| | IS4 (CRY) | | IS5 (PRY) | | | |
|-----------------|-----------|-------|-----------|-------|--------|-------|
| | AREA # | RT # | AREA # | RT # | AREA # | RT # |
| ===== | ===== | ===== | ===== | ===== | ===== | ===== |
| ICAL MIDPT | 503160 | 13.80 | 442215 | 17.47 | | |
| UPPER LIMIT | 1006320 | | 884430 | | | |
| LOWER LIMIT | 251580 | | 221108 | | | |
| ===== | ===== | ===== | ===== | ===== | ===== | ===== |
| CCAL | 526425 | 13.68 | 439084 | 17.32 | | |
| UPPER LIMIT | | 14.18 | | 17.82 | | |
| LOWER LIMIT | | 13.18 | | 16.82 | | |
| 01 SN54MBS2 | 371865 | 13.67 | 319297 | 17.32 | | |
| 02 SN54LCSS2 | 381079 | 13.67 | 335486 | 17.32 | | |
| 03 SN54LCSDS2 | 383014 | 13.68 | 345990 | 17.32 | | |
| 04 LL-SED3-0-36 | 373966 | 13.73 | 298676 | 17.39 | | |
| 05 LL-SED3-36-1 | 322033 | 13.69 | 280959 | 17.35 | | |
| 06 LL-SED3-141- | 293472 | 13.69 | 247444 | 17.36 | | |
| 07 LL-SED2-0-56 | 339995 | 13.73 | 242975 | 17.38 | | |
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IS4 = Chrysene-d12
IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

* Values outside of QC limits.

8B
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Ical Midpoint ID: 03311102

Ical Date: 03/31/11

Instrument ID: NT12

Cont. Cal Date: 04/04/11

| | IS1 (NPT) AREA # | RT # | IS2 (ANT) AREA # | RT # | IS3 (PHN) AREA # | RT # |
|-----------------|---------------------|------|---------------------|------|---------------------|------|
| ICAL MIDPT | 494112 | 4.92 | 280105 | 7.16 | 461353 | 9.10 |
| UPPER LIMIT | 988224 | | 560210 | | 922706 | |
| LOWER LIMIT | 247056 | | 140052 | | 230676 | |
| CCAL | 401183 | 4.80 | 270252 | 7.04 | 455484 | 8.98 |
| UPPER LIMIT | | 5.30 | | 7.54 | | 9.48 |
| LOWER LIMIT | | 4.30 | | 6.54 | | 8.48 |
| 01 LL-SED2-56-1 | 327676 | 4.80 | 205193 | 7.04 | 322531 | 8.98 |
| 02 LL-SED2-56-1 | 303192 | 4.80 | 192063 | 7.04 | 293988 | 8.98 |
| 03 LL-SED2-56-1 | 262660 | 4.80 | 181702 | 7.04 | 299869 | 8.98 |
| 04 LL-SED2-112- | 223462* | 4.80 | 148879 | 7.05 | 248837 | 8.98 |
| 05 LL-SED2-0-56 | 210690* | 4.80 | 140224 | 7.04 | 232179 | 8.99 |
| 06 LL-SED1-0-56 | 208943* | 4.80 | 140371 | 7.05 | 227824* | 8.98 |
| 07 LL-SED2-0-56 | 261511 | 4.81 | 166500 | 7.05 | 238192 | 8.99 |
| 08 LL-SED1-0-56 | 269407 | 4.81 | 163401 | 7.05 | 239810 | 8.98 |
| 09 LL-SED2-112- | 305119 | 4.81 | 194866 | 7.05 | 291590 | 8.99 |
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| 24 | | | | | | |
| 25 | | | | | | |

IS1 = Naphthalene-d8
IS2 = Acenaphthene-d10
IS3 = Phenanthrene-d10

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal
RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

* Values outside of QC limits.

8B
SEMIVOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No: SN54

Project: LORA LAKE

Ical Midpoint ID: 03311102

Ical Date: 03/31/11

Instrument ID: NT12

Cont. Cal Date: 04/04/11

| | IS4 (CRY) AREA # | RT # | IS5 (PRY) AREA # | RT # | AREA # | RT # |
|-----------------|---------------------|-------|---------------------|-------|--------|------|
| ICAL MIDPT | 503160 | 13.80 | 442215 | 17.47 | | |
| UPPER LIMIT | 1006320 | | 884430 | | | |
| LOWER LIMIT | 251580 | | 221108 | | | |
| CCAL | 430701 | 13.60 | 334445 | 17.23 | | |
| UPPER LIMIT | | 14.10 | | 17.73 | | |
| LOWER LIMIT | | 13.10 | | 16.73 | | |
| 01 LL-SED2-56-1 | 332757 | 13.61 | 295718 | 17.26 | | |
| 02 LL-SED2-56-1 | 320215 | 13.61 | 271011 | 17.26 | | |
| 03 LL-SED2-56-1 | 326706 | 13.61 | 267412 | 17.26 | | |
| 04 LL-SED2-112- | 266485 | 13.62 | 175502* | 17.27 | | |
| 05 LL-SED2-0-56 | 292600 | 13.66 | 173734* | 17.31 | | |
| 06 LL-SED1-0-56 | 279926 | 13.65 | 163012* | 17.30 | | |
| 07 LL-SED2-0-56 | 263188 | 13.64 | 165436* | 17.28 | | |
| 08 LL-SED1-0-56 | 272758 | 13.63 | 171762* | 17.26 | | |
| 09 LL-SED2-112- | 253711 | 13.62 | 174659* | 17.27 | | |
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| 25 | | | | | | |

IS4 = Chrysene-d12
IS5 = Perylene-d12

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Cont. Cal
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Cont. Cal

* Values outside of QC limits.

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

ARI Job ID: SN54

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

Sample ID: LL-SED3-0-36-031511
SAMPLE

Lab Sample ID: SN54A
LIMS ID: 11-5925
Matrix: Sediment
Data Release Authorized: *W*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 03:51
Instrument/Analyst: ECD1/YZ

Sample Amount: 2.46 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 75.4%

| CAS Number | Analyte | RL | Result |
|--|----------------------|-------|--------|
| 87-86-5 | Pentachlorophenol | 25 | < 25 U |
| Reported in µg/kg (ppb) | | | |
| <u>Chlorophenol Surrogate Recovery</u> | | | |
| | 2,4,6-Tribromophenol | 91.2% | |

Sample ID: LL-SED3-36-141-031511
SAMPLE

Lab Sample ID: SN54B
LIMS ID: 11-5926
Matrix: Sediment
Data Release Authorized: *WV*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 04:27
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.17 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 88.3%

| CAS Number | Analyte | RL | Result |
|---------------------------------|----------------------|-------|--------|
| 87-86-5 | Pentachlorophenol | 53 | < 53 U |
| Reported in µg/kg (ppb) | | | |
| Chlorophenol Surrogate Recovery | | | |
| | 2,4,6-Tribromophenol | 71.2% | |

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

Sample ID: LL-SED3-141-167-031511
SAMPLE

Lab Sample ID: SN54C
LIMS ID: 11-5927
Matrix: Sediment
Data Release Authorized: *mw*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 05:03
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.22 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 87.9%

| CAS Number | Analyte | RL | Result |
|--|----------------------|-------|--------|
| 87-86-5 | Pentachlorophenol | 51 | < 51 U |
| Reported in µg/kg (ppb) | | | |
| <u>Chlorophenol Surrogate Recovery</u> | | | |
| | 2,4,6-Tribromophenol | 70.4% | |

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

Sample ID: LL-SED2-0-56-031511
SAMPLE

Lab Sample ID: SN54D
LIMS ID: 11-5928
Matrix: Sediment
Data Release Authorized: *MW*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 05:39
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.42 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 85.9%

| CAS Number | Analyte | RL | Result |
|------------|-------------------|----|--------|
| 87-86-5 | Pentachlorophenol | 44 | < 44 U |

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 88.8% |
|----------------------|-------|

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

Sample ID: LL-SED2-56-112-031511
SAMPLE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *MM*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 06:15
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.38 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|---------------------------------|----------------------|-------|--------|
| 87-86-5 | Pentachlorophenol | 45 | < 45 U |
| Reported in µg/kg (ppb) | | | |
| Chlorophenol Surrogate Recovery | | | |
| | 2,4,6-Tribromophenol | 87.2% | |

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

Sample ID: LL-SED2-112-168-031511
SAMPLE

Lab Sample ID: SN54F
LIMS ID: 11-5930
Matrix: Sediment
Data Release Authorized: *mw*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 10:20
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.23 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 87.7%

| CAS Number | Analyte | RL | Result |
|------------|-------------------|----|--------|
| 87-86-5 | Pentachlorophenol | 51 | < 51 U |

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 80.4% |
|----------------------|-------|

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

Sample ID: LL-SED2-0-56-031511-D
SAMPLE

Lab Sample ID: SN54G
LIMS ID: 11-5931
Matrix: Sediment
Data Release Authorized: *WW*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 10:56
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.14 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 88.7%

| CAS Number | Analyte | RL | Result |
|------------|-------------------|----|--------|
| 87-86-5 | Pentachlorophenol | 55 | < 55 U |

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 90.8% |
|----------------------|-------|

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

Sample ID: LL-SED1-0-56-031511
SAMPLE

Lab Sample ID: SN54H
LIMS ID: 11-5932
Matrix: Sediment
Data Release Authorized: *WV*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 11:32
Instrument/Analyst: ECD1/YZ

Sample Amount: 5.45 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 45.6%

| CAS Number | Analyte | RL | Result |
|--|----------------------|-------|---------|
| 87-86-5 | Pentachlorophenol | 110 | < 110 Y |
| Reported in µg/kg (ppb) | | | |
| Chlorophenol Surrogate Recovery | | | |
| | 2,4,6-Tribromophenol | 82.4% | |

SW8041 CHLOROPHENOLICS SURROGATE RECOVERY SUMMARY

Matrix: Sediment

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL

| Client ID | TBP | TOT OUT |
|---------------------------|-------|---------|
| LL-SED3-0-36-031511 | 91.2% | 0 |
| LL-SED3-36-141-031511 | 71.2% | 0 |
| LL-SED3-141-167-031511 | 70.4% | 0 |
| LL-SED2-0-56-031511 | 88.8% | 0 |
| MB-032811 | 67.6% | 0 |
| LCS-032811 | 69.4% | 0 |
| LL-SED2-56-112-031511 | 87.2% | 0 |
| LL-SED2-56-112-031511 MS | 78.4% | 0 |
| LL-SED2-56-112-031511 MSD | 75.2% | 0 |
| LL-SED2-112-168-031511 | 80.4% | 0 |
| LL-SED2-0-56-031511-D | 90.8% | 0 |
| LL-SED1-0-56-031511 | 82.4% | 0 |

LCS/MB LIMITS QC LIMITS

(TBP) = 2,4,6-Tribromophenol

(50-115)

(10-146)

Prep Method: SW3550B
Log Number Range: 11-5925 to 11-5932

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

Sample ID: LL-SED2-56-112-031511
MS/MSD

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *mmw*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted MS/MSD: 03/28/11

Sample Amount MS: 1.39 g-dry-wt
MSD: 1.38 g-dry-wt

Date Analyzed MS: 04/07/11 06:51
MSD: 04/07/11 07:27

Final Extract Volume MS: 25 mL
MSD: 25 mL

Instrument/Analyst MS: ECD1/YZ
MSD: ECD1/YZ

Dilution Factor MS: 1.00
MSD: 1.00

Percent Moisture: 86.2%

| Analyte | Sample | MS | Spike Added-MS | MS Recovery | MSD | Spike Added-MSD | MSD Recovery | RPD |
|-------------------|--------|-----|----------------|-------------|-----|-----------------|--------------|------|
| Pentachlorophenol | < 45.3 | 308 | 450 | 68.4% | 316 | 453 | 69.8% | 2.6% |

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

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

Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *mw*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 06:51
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.39 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|------------|-------------------|----|--------|
| 87-86-5 | Pentachlorophenol | 45 | --- |

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 78.4% |
|----------------------|-------|

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

Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE DUP

Lab Sample ID: SN54E
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *WVW*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/07/11 07:27
Instrument/Analyst: ECD1/YZ

Sample Amount: 1.38 g-dry-wt
Final Extract Volume: 25 mL
Dilution Factor: 1.00
Percent Moisture: 86.2%

| CAS Number | Analyte | RL | Result |
|------------|-------------------|----|--------|
| 87-86-5 | Pentachlorophenol | 45 | --- |

Reported in µg/kg (ppb)

Chlorophenol Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 75.2% |
|----------------------|-------|

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

Sample ID: LCS-032811
LAB CONTROL

Lab Sample ID: LCS-032811
LIMS ID: 11-5929
Matrix: Sediment
Data Release Authorized: *mm*
Reported: 04/07/11

QC Report No: SN54-Floyd Snider
Project: Lora Lake - Subsurface Sediment
POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Date Extracted: 03/28/11
Date Analyzed: 04/06/11 18:13
Instrument/Analyst: ECD1/YZ

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

| Analyte | Lab Control | Spike Added | Recovery |
|-------------------|-------------|-------------|----------|
| Pentachlorophenol | 50.2 | 62.5 | 80.3% |

Chlorophenols Surrogate Recovery

| | |
|----------------------|-------|
| 2,4,6-Tribromophenol | 69.4% |
|----------------------|-------|

Results reported in µg/kg

4
CHLOROPHENOL METHOD BLANK SUMMARY

SAMPLE NO.

| |
|----------|
| SN54MBS1 |
|----------|

| | |
|-------------------------------------|--------------------------------------|
| Lab Name: ANALYTICAL RESOURCES, INC | Client: FLOYD SNIDER |
| ARI Job No.: SN54 | Project: LORA LAKE |
| Lab Sample ID: SN54MBS1 | Lab File ID: 0401A005 |
| Matrix (soil/water) SOLID | Extraction: (SepF/Cont/Sonc) SW3550C |
| Sulfur Cleanup (Y/N) Y | Date Extracted: 03/28/11 |
| Date Analyzed (1): 04/06/11 | Date Analyzed (2): 04/06/11 |
| Time Analyzed (1): 1737 | Time Analyzed (2): 1737 |
| Instrument ID (1): ECD1 | Instrument ID (2): ECD1 |
| GC Column (1): ZB5 ID: 0.53 (mm) | 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 | SN54LCSS1 | SN54LCSS1 | 04/06/11 | 04/06/11 |
| 02 | LL-SED3-0-36 | SN54A | 04/07/11 | 04/07/11 |
| 03 | LL-SED3-36-1 | SN54B | 04/07/11 | 04/07/11 |
| 04 | LL-SED3-141- | SN54C | 04/07/11 | 04/07/11 |
| 05 | LL-SED2-0-56 | SN54D | 04/07/11 | 04/07/11 |
| 06 | LL-SED2-56-1 | SN54E | 04/07/11 | 04/07/11 |
| 07 | LL-SED2-56-1 | SN54EMS | 04/07/11 | 04/07/11 |
| 08 | LL-SED2-56-1 | SN54EMSD | 04/07/11 | 04/07/11 |
| 09 | LL-SED2-112- | SN54F | 04/07/11 | 04/07/11 |
| 10 | LL-SED2-0-56 | SN54G | 04/07/11 | 04/07/11 |
| 11 | LL-SED1-0-56 | SN54H | 04/07/11 | 04/07/11 |

ORGANICS ANALYSIS DATA SHEET

PCP by GC/ECD Method SW8041

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Sample ID: MB-032811

METHOD BLANK

Lab Sample ID: MB-032811

LIMS ID: 11-5929

Matrix: Sediment

Data Release Authorized: *mw*

Reported: 04/07/11

QC Report No: SN54-Floyd Snider

Project: Lora Lake - Subsurface Sediment

POS-LL

Date Sampled: NA

Date Received: NA

Date Extracted: 03/28/11

Date Analyzed: 04/06/11 17:37

Instrument/Analyst: ECD1/YZ

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 | 67.6% |
|----------------------|-------|

6D
 CHLOROPHENOL INITIAL CALIBRATION
 RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SN54

Project: LORA LAKE

GC Column: ZB5 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/01/11

| COMPOUND | RT OF STANDARDS | | | | | | MEAN RT | RT WINDOW | |
|----------------------|-----------------|-------|-------|-------|-------|-------|------------|-----------|-------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | FROM | TO |
| Pentachlorophenol | 21.82 | 21.81 | 21.80 | 21.79 | 21.79 | 21.78 | 21.80 | 21.72 | 21.86 |
| 2,4,6-Trichloropheno | 13.41 | 13.41 | 13.41 | 13.41 | 13.41 | 13.41 | 13.41 | 13.34 | 13.48 |
| 2,3,6-Trichloropheno | 14.38 | 14.38 | 14.38 | 14.38 | 14.38 | 14.38 | 14.38 | 14.31 | 14.45 |
| 2,4,5-Trichloropheno | | 16.12 | 16.10 | 16.08 | 16.08 | 16.07 | 16.09 | 16.01 | 16.15 |
| 2,3,4-Trichloropheno | 17.29 | 17.39 | 17.37 | 17.35 | 17.33 | 17.32 | 17.35 | 17.28 | 17.42 |
| 2,3,5,6-Tetrachlorop | 17.73 | 17.73 | 17.73 | 17.72 | 17.72 | 17.72 | 17.72 | 17.65 | 17.79 |
| 2,3,4,5-Tetrachlorop | 20.59 | 20.57 | 20.56 | 20.54 | 20.53 | 20.53 | 20.56 | 20.48 | 20.61 |
| 2,4-Dichlorophenol | 12.43 | 12.44 | 12.43 | 12.43 | 12.43 | 12.43 | 12.43 | 12.36 | 12.50 |
| 2,4,6-Tribromophenol | 19.69 | 19.68 | 19.66 | 19.65 | 19.64 | 19.64 | 19.66 | 19.58 | 19.72 |

<-

6D
 CHLOROPHENOL INITIAL CALIBRATION
 RETENTION TIME WINDOWS

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SN54

Project: LORA LAKE

GC Column: ZB35 ID: 0.53 (mm)

Instrument ID: ECD1

Calibration Date: 04/01/11

| COMPOUND | RT OF STANDARDS | | | | | | MEAN RT | RT WINDOW | |
|----------------------|-----------------|-------|-------|-------|-------|-------|------------|-----------|-------|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | FROM | TO |
| Pentachlorophenol | 22.65 | 22.64 | 22.64 | 22.63 | 22.63 | 22.62 | 22.64 | 22.56 | 22.70 |
| 2,4,6-Trichloropheno | 13.78 | 13.78 | 13.78 | 13.78 | 13.78 | 13.78 | 13.78 | 13.71 | 13.85 |
| 2,3,6-Trichloropheno | 15.19 | 15.19 | 15.19 | 15.18 | 15.18 | 15.18 | 15.18 | 15.11 | 15.25 |
| 2,4,5-Trichloropheno | 17.19 | 17.17 | 17.16 | 17.15 | 17.15 | 17.14 | 17.16 | 17.08 | 17.22 |
| 2,3,4-Trichloropheno | 18.76 | 18.75 | 18.74 | 18.73 | 18.71 | 18.71 | 18.74 | 18.66 | 18.80 |
| 2,3,5,6-Tetrachlorop | 18.42 | 18.42 | 18.41 | 18.41 | 18.41 | 18.41 | 18.41 | 18.34 | 18.48 |
| 2,3,4,5-Tetrachlorop | 21.85 | 21.84 | 21.83 | 21.82 | 21.81 | 21.81 | 21.83 | 21.75 | 21.89 |
| 2,4-Dichlorophenol | 13.43 | 13.43 | 13.43 | 13.43 | 13.43 | 13.43 | 13.43 | 13.36 | 13.50 |
| 2,4,6-Tribromophenol | 20.89 | 20.89 | 20.88 | 20.87 | 20.86 | 20.86 | 20.88 | 20.80 | 20.94 |

6E
 CHLOROPHENOL INITIAL CALIBRATION
 CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm) Instrument ID: ECD1
 Calibration Date: 04/01/11

| COMPOUND | CALIBRATION FACTORS | | | | | | R ² / %RSD | CT |
|-----------------------|---------------------|-------|-------|-------|-------|-------|--------------------------|----|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | |
| Pentachlorophenol | 29405 | 27955 | 27704 | 22662 | 24213 | 22067 | 12.0 | A |
| 2,4,6-Trichlorophenol | 23973 | 19056 | 17104 | 13948 | 14398 | 12938 | 0.9928 | L |
| 2,3,6-Trichlorophenol | 17697 | 15903 | 15568 | 12834 | 13478 | 12207 | 14.4 | A |
| 2,4,5-Trichlorophenol | | 8437 | 8744 | 7336 | 7536 | 6502 | 11.6 | A |
| 2,3,4-Trichlorophenol | 15137 | 12572 | 13596 | 10067 | 9677 | 7988 | 0.9986 | Q |
| 2,3,5,6-Tetrachloroph | 28937 | 25200 | 24487 | 20185 | 21240 | 19116 | 15.9 | A |
| 2,3,4,5-Tetrachloroph | 19018 | 18483 | 17317 | 14350 | 15450 | 13546 | 13.7 | A |
| 2,4-Dichlorophenol | 1167 | 995 | 930 | 711 | 692 | 595 | 0.9988 | Q |
| 2,4,6-Tribromophenol | 24452 | 23977 | 22590 | 18840 | 20483 | 18745 | 11.7 | A |
| AVE RSD | | | | | | | 20.6 | |

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/ecdl.i/PCP20110401.b/ical-1.b/0401A009.d
 LVL 2: /chem2/ecdl.i/PCP20110401.b/ical-1.b/0401A010.d
 LVL 3: /chem2/ecdl.i/PCP20110401.b/ical-1.b/0401A011.d
 LVL 4: /chem2/ecdl.i/PCP20110401.b/ical-1.b/0401A008.d
 LVL 5: /chem2/ecdl.i/PCP20110401.b/ical-1.b/0401A012.d
 LVL 6: /chem2/ecdl.i/PCP20110401.b/ical-1.b/0401A013.d

6E
 CHLOROPHENOL INITIAL CALIBRATION
 CALIBRATION FACTORS

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB35 ID: 0.53 (mm) Instrument ID: ECD1
 Calibration Date: 04/01/11

| COMPOUND | CALIBRATION FACTORS | | | | | | R ² / %RSD | CT |
|-----------------------|---------------------|-------|-------|-------|-------|-------|--------------------------|----|
| | LVL 1 | LVL 2 | LVL 3 | LVL 4 | LVL 5 | LVL 6 | | |
| Pentachlorophenol | 27844 | 27891 | 28971 | 25017 | 28058 | 26455 | 5.1 | A |
| 2,4,6-Trichlorophenol | 18067 | 16685 | 17022 | 14481 | 16094 | 14858 | 8.4 | A |
| 2,3,6-Trichlorophenol | 16623 | 15623 | 15797 | 13291 | 14677 | 13675 | 8.7 | A |
| 2,4,5-Trichlorophenol | 9223 | 8624 | 8724 | 7196 | 7623 | 6835 | 11.9 | A |
| 2,3,4-Trichlorophenol | 10128 | 10897 | 11196 | 8917 | 9732 | 8688 | 10.2 | A |
| 2,3,5,6-Tetrachloroph | 25256 | 23960 | 24545 | 20670 | 22769 | 21310 | 7.9 | A |
| 2,3,4,5-Tetrachloroph | 16901 | 17108 | 17518 | 14662 | 16006 | 14923 | 7.4 | A |
| 2,4-Dichlorophenol | 1052 | 934 | 903 | 724 | 742 | 641 | 18.6 | A |
| 2,4,6-Tribromophenol | 21409 | 20597 | 22269 | 19519 | 22526 | 21660 | 5.2 | A |
| AVE RSD | | | | | | | 9.3 | |

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/ecdl.i/PCP20110401.b/ical-2.b/0401A009.d
 LVL 2: /chem2/ecdl.i/PCP20110401.b/ical-2.b/0401A010.d
 LVL 3: /chem2/ecdl.i/PCP20110401.b/ical-2.b/0401A011.d
 LVL 4: /chem2/ecdl.i/PCP20110401.b/ical-2.b/0401A008.d
 LVL 5: /chem2/ecdl.i/PCP20110401.b/ical-2.b/0401A012.d
 LVL 6: /chem2/ecdl.i/PCP20110401.b/ical-2.b/0401A013.d

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No. (PCP): Date Analyzed :04/06/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :1701

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|-----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 21.79 | 21.72 | 21.86 | 22.9 | 25.0 | -8.4 |
| 2,4,6-Trichlorophenol | 13.41 | 13.34 | 13.48 | 27.7 | 25.0 | 10.8 |
| 2,3,6-Trichlorophenol | 14.38 | 14.31 | 14.45 | 22.6 | 25.0 | -9.6 |
| 2,4,5-Trichlorophenol | 16.08 | 16.01 | 16.15 | 23.8 | 25.0 | -4.8 |
| 2,3,4-Trichlorophenol | 17.33 | 17.28 | 17.42 | 24.5 | 25.0 | -2.0 |
| 2,3,5,6-Tetrachlorophenol | 17.72 | 17.65 | 17.79 | 22.5 | 25.0 | -10.0 |
| 2,3,4,5-Tetrachlorophenol | 20.54 | 20.48 | 20.61 | 22.5 | 25.0 | -10.0 |
| 2,4-Dichlorophenol | 12.43 | 12.36 | 12.50 | 256 | 250 | 2.4 |
| 2,4,6-Tribromophenol (surr) | 19.65 | 19.58 | 19.72 | 23.0 | 25.0 | -8.0 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 7.3

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB35 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP): Date Analyzed :04/06/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :1701

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 22.63 | 22.56 | 22.70 | 22.3 | 25.0 | -10.8 |
| 2,4,6-Trichlorophenol | 13.77 | 13.71 | 13.85 | 22.0 | 25.0 | -12.0 |
| 2,3,6-Trichlorophenol | 15.18 | 15.11 | 15.25 | 22.0 | 25.0 | -12.0 |
| 2,4,5-Trichlorophenol | 17.15 | 17.08 | 17.22 | 22.3 | 25.0 | -10.8 |
| 2,3,4-Trichlorophenol | 18.72 | 18.66 | 18.80 | 22.9 | 25.0 | -8.4 |
| 2,3,5,6-Tetrachlorophenol | 18.41 | 18.34 | 18.48 | 22.1 | 25.0 | -11.6 |
| 2,3,4,5-Tetrachlorophenol | 21.81 | 21.75 | 21.89 | 22.8 | 25.0 | -8.8 |
| 2,4-Dichlorophenol | 13.43 | 13.36 | 13.50 | 222 | 250 | -11.2 |
| 2,4,6-Tribromophenol (surr | 20.86 | 20.80 | 20.94 | 21.8 | 25.0 | -12.8 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 10.9

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP): Date Analyzed :04/07/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :0015

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 21.81 | 21.72 | 21.86 | 21.9 | 25.0 | -12.4 |
| 2,4,6-Trichlorophenol | 13.42 | 13.34 | 13.48 | 27.3 | 25.0 | 9.2 |
| 2,3,6-Trichlorophenol | 14.39 | 14.31 | 14.45 | 22.6 | 25.0 | -9.6 |
| 2,4,5-Trichlorophenol | 16.10 | 16.01 | 16.15 | 23.8 | 25.0 | -4.8 |
| 2,3,4-Trichlorophenol | 17.36 | 17.28 | 17.42 | 21.6 | 25.0 | -13.6 |
| 2,3,5,6-Tetrachlorophenol | 17.73 | 17.65 | 17.79 | 22.5 | 25.0 | -10.0 |
| 2,3,4,5-Tetrachlorophenol | 20.56 | 20.48 | 20.61 | 23.3 | 25.0 | -6.8 |
| 2,4-Dichlorophenol | 12.44 | 12.36 | 12.50 | 234 | 250 | -6.4 |
| 2,4,6-Tribromophenol (surr | 19.66 | 19.58 | 19.72 | 23.2 | 25.0 | -7.2 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 8.9

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER

ARI Job No.: SN54 Project: LORA LAKE

GC Column: ZB35 ID: 0.53 (mm)

Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP): Date Analyzed :04/07/11

Lab Sample ID (PCP): PCP CCAL Time Analyzed :0015

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|-----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 22.64 | 22.56 | 22.70 | 23.6 | 25.0 | -5.6 |
| 2,4,6-Trichlorophenol | 13.79 | 13.71 | 13.85 | 22.7 | 25.0 | -9.2 |
| 2,3,6-Trichlorophenol | 15.19 | 15.11 | 15.25 | 22.5 | 25.0 | -10.0 |
| 2,4,5-Trichlorophenol | 17.16 | 17.08 | 17.22 | 23.0 | 25.0 | -8.0 |
| 2,3,4-Trichlorophenol | 18.74 | 18.66 | 18.80 | 23.6 | 25.0 | -5.6 |
| 2,3,5,6-Tetrachlorophenol | 18.42 | 18.34 | 18.48 | 22.7 | 25.0 | -9.2 |
| 2,3,4,5-Tetrachlorophenol | 21.83 | 21.75 | 21.89 | 23.2 | 25.0 | -7.2 |
| 2,4-Dichlorophenol | 13.44 | 13.36 | 13.50 | 22.7 | 25.0 | -9.2 |
| 2,4,6-Tribromophenol (surr) | 20.88 | 20.80 | 20.94 | 23.0 | 25.0 | -8.0 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 8.0

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP): Date Analyzed :04/07/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :0315

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|-----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 21.81 | 21.72 | 21.86 | 22.0 | 25.0 | -12.0 |
| 2,4,6-Trichlorophenol | 13.42 | 13.34 | 13.48 | 27.4 | 25.0 | 9.6 |
| 2,3,6-Trichlorophenol | 14.39 | 14.31 | 14.45 | 22.8 | 25.0 | -8.8 |
| 2,4,5-Trichlorophenol | 16.10 | 16.01 | 16.15 | 24.7 | 25.0 | -1.2 |
| 2,3,4-Trichlorophenol | 17.36 | 17.28 | 17.42 | 21.8 | 25.0 | -12.8 |
| 2,3,5,6-Tetrachlorophenol | 17.74 | 17.65 | 17.79 | 23.2 | 25.0 | -7.2 |
| 2,3,4,5-Tetrachlorophenol | 20.56 | 20.48 | 20.61 | 23.1 | 25.0 | -7.6 |
| 2,4-Dichlorophenol | 12.45 | 12.36 | 12.50 | 256 | 250 | 2.4 |
| 2,4,6-Tribromophenol (surr) | 19.67 | 19.58 | 19.72 | 23.4 | 25.0 | -6.4 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 7.6

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB35 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No. (PCP): Date Analyzed :04/07/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :0315

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|-------|
| ===== | ===== | FROM | TO | ===== | ===== | ===== |
| Pentachlorophenol | 22.65 | 22.56 | 22.70 | 23.9 | 25.0 | -4.4 |
| 2,4,6-Trichlorophenol | 13.79 | 13.71 | 13.85 | 23.0 | 25.0 | -8.0 |
| 2,3,6-Trichlorophenol | 15.20 | 15.11 | 15.25 | 22.6 | 25.0 | -9.6 |
| 2,4,5-Trichlorophenol | 17.17 | 17.08 | 17.22 | 23.2 | 25.0 | -7.2 |
| 2,3,4-Trichlorophenol | 18.74 | 18.66 | 18.80 | 23.2 | 25.0 | -7.2 |
| 2,3,5,6-Tetrachlorophenol | 18.42 | 18.34 | 18.48 | 22.3 | 25.0 | -10.8 |
| 2,3,4,5-Tetrachlorophenol | 21.83 | 21.75 | 21.89 | 23.4 | 25.0 | -6.4 |
| 2,4-Dichlorophenol | 13.44 | 13.36 | 13.50 | 229 | 250 | -8.4 |
| 2,4,6-Tribromophenol (surr | 20.88 | 20.80 | 20.94 | 23.3 | 25.0 | -6.8 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 7.6

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP): Date Analyzed :04/07/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed :0944

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|-------|
| | | FROM | TO | | | |
| Pentachlorophenol | 21.81 | 21.72 | 21.86 | 21.4 | 25.0 | -14.4 |
| 2,4,6-Trichlorophenol | 13.42 | 13.34 | 13.48 | 28.1 | 25.0 | 12.4 |
| 2,3,6-Trichlorophenol | 14.39 | 14.31 | 14.45 | 23.3 | 25.0 | -6.8 |
| 2,4,5-Trichlorophenol | 16.09 | 16.01 | 16.15 | 24.5 | 25.0 | -2.0 |
| 2,3,4-Trichlorophenol | 17.35 | 17.28 | 17.42 | 21.1 | 25.0 | -15.6 |
| 2,3,5,6-Tetrachlorophenol | 17.73 | 17.65 | 17.79 | 22.4 | 25.0 | -10.4 |
| 2,3,4,5-Tetrachlorophenol | 20.55 | 20.48 | 20.61 | 22.2 | 25.0 | -11.2 |
| 2,4-Dichlorophenol | 12.45 | 12.36 | 12.50 | 266 | 250 | 6.4 |
| 2,4,6-Tribromophenol (surr | 19.66 | 19.58 | 19.72 | 23.0 | 25.0 | -8.0 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 9.7

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SN54

Project: LORA LAKE

GC Column: ZB35 ID: 0.53 (mm)

Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No.(PCP):

Date Analyzed :04/07/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :0944

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|------|
| | | FROM | TO | | | |
| Pentachlorophenol | 22.64 | 22.56 | 22.70 | 23.2 | 25.0 | -7.2 |
| 2,4,6-Trichlorophenol | 13.79 | 13.71 | 13.85 | 24.1 | 25.0 | -3.6 |
| 2,3,6-Trichlorophenol | 15.20 | 15.11 | 15.25 | 23.6 | 25.0 | -5.6 |
| 2,4,5-Trichlorophenol | 17.16 | 17.08 | 17.22 | 23.4 | 25.0 | -6.4 |
| 2,3,4-Trichlorophenol | 18.73 | 18.66 | 18.80 | 23.9 | 25.0 | -4.4 |
| 2,3,5,6-Tetrachlorophenol | 18.42 | 18.34 | 18.48 | 23.6 | 25.0 | -5.6 |
| 2,3,4,5-Tetrachlorophenol | 21.83 | 21.75 | 21.89 | 22.8 | 25.0 | -8.8 |
| 2,4-Dichlorophenol | 13.44 | 13.36 | 13.50 | 236 | 250 | -5.6 |
| 2,4,6-Tribromophenol (surr | 20.88 | 20.80 | 20.94 | 23.3 | 25.0 | -6.8 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 6.0

7E
CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm)
 Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No. (PCP): Date Analyzed : 04/07/11
 Lab Sample ID (PCP): PCP CCAL Time Analyzed : 1244

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|------|
| | | FROM | TO | | | |
| Pentachlorophenol | 21.80 | 21.72 | 21.86 | 23.0 | 25.0 | -8.0 |
| 2,4,6-Trichlorophenol | 13.42 | 13.34 | 13.48 | 28.9 | 25.0 | 15.6 |
| 2,3,6-Trichlorophenol | 14.38 | 14.31 | 14.45 | 24.1 | 25.0 | -3.6 |
| 2,4,5-Trichlorophenol | 16.09 | 16.01 | 16.15 | 24.6 | 25.0 | -1.6 |
| 2,3,4-Trichlorophenol | 17.34 | 17.28 | 17.42 | 22.7 | 25.0 | -9.2 |
| 2,3,5,6-Tetrachlorophenol | 17.72 | 17.65 | 17.79 | 23.2 | 25.0 | -7.2 |
| 2,3,4,5-Tetrachlorophenol | 20.54 | 20.48 | 20.61 | 23.2 | 25.0 | -7.2 |
| 2,4-Dichlorophenol | 12.44 | 12.36 | 12.50 | 274 | 250 | 9.6 |
| 2,4,6-Tribromophenol (surr | 19.65 | 19.58 | 19.72 | 23.3 | 25.0 | -6.8 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 7.6

7E
 CHLOROPHENOL CALIBRATION VERIFICATION SUMMARY

Lab Name: ANALYTICAL RESOURCES, INC

Client: FLOYD SNIDER

ARI Job No.: SN54

Project: LORA LAKE

GC Column: ZB35 ID: 0.53 (mm)

Init. Calib. Date(s): 04/01/11 04/01/11

Client Sample No. (PCP):

Date Analyzed :04/07/11

Lab Sample ID (PCP): PCP CCAL

Time Analyzed :1244

| PCP MIX COMPOUND | RT | RT WINDOW | | CALC AMOUNT | NOM AMOUNT | %D |
|----------------------------|-------|-----------|-------|----------------|---------------|------|
| | | FROM | TO | | | |
| Pentachlorophenol | 22.63 | 22.56 | 22.70 | 24.5 | 25.0 | -2.0 |
| 2,4,6-Trichlorophenol | 13.78 | 13.71 | 13.85 | 25.5 | 25.0 | 2.0 |
| 2,3,6-Trichlorophenol | 15.19 | 15.11 | 15.25 | 24.7 | 25.0 | -1.2 |
| 2,4,5-Trichlorophenol | 17.15 | 17.08 | 17.22 | 24.4 | 25.0 | -2.4 |
| 2,3,4-Trichlorophenol | 18.72 | 18.66 | 18.80 | 24.6 | 25.0 | -1.6 |
| 2,3,5,6-Tetrachlorophenol | 18.42 | 18.34 | 18.48 | 24.5 | 25.0 | -2.0 |
| 2,3,4,5-Tetrachlorophenol | 21.82 | 21.75 | 21.89 | 23.9 | 25.0 | -4.4 |
| 2,4-Dichlorophenol | 13.44 | 13.36 | 13.50 | 250 | 250 | 0.0 |
| 2,4,6-Tribromophenol (surr | 20.87 | 20.80 | 20.94 | 24.1 | 25.0 | -3.6 |
| | | | | | | |
| | | | | | | |

AVERAGE %D = 2.1

8
CHLOROPHENOL ANALYTICAL SEQUENCE

Lab Name: ANALYTICAL RESOURCES, INC Client: FLOYD SNIDER
 ARI Job No.: SN54 Project: LORA LAKE
 GC Column: ZB5 ID: 0.53 (mm) Instrument ID: ECD1
 Init. Calib. Date(s): 04/01/11 04/01/11

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

| MEAN SURROGATE RT FROM INITIAL CALIBRATION | | | | | |
|--|--------------|----------|----------|-------|-------|
| S1 : 19.65 | | | | | |
| CLIENT | LAB | DATE | TIME | S1 | |
| SAMPLE NO. | SAMPLE ID | ANALYZED | ANALYZED | RT | # |
| ===== | ===== | ===== | ===== | ===== | ===== |
| 01 | PCP D | 04/01/11 | 1629 | 19.65 | |
| 02 | PCP A | 04/01/11 | 1706 | 19.69 | |
| 03 | PCP B | 04/01/11 | 1742 | 19.68 | |
| 04 | PCP C | 04/01/11 | 1818 | 19.66 | |
| 05 | PCP E | 04/01/11 | 1854 | 19.64 | |
| 06 | PCP F | 04/01/11 | 1931 | 19.64 | |
| 07 | PCP CCAL | 04/06/11 | 1701 | 19.65 | |
| 08 | SN54MBS1 | 04/06/11 | 1737 | 19.65 | |
| 09 | SN54LCSS1 | 04/06/11 | 1813 | 19.65 | |
| 10 | PCP CCAL | 04/07/11 | 0015 | 19.66 | |
| 11 | PCP CCAL | 04/07/11 | 0315 | 19.67 | |
| 12 | LL-SED3-0-36 | 04/07/11 | 0351 | 19.65 | |
| 13 | LL-SED3-36-1 | 04/07/11 | 0427 | 19.65 | |
| 14 | LL-SED3-141- | 04/07/11 | 0503 | 19.65 | |
| 15 | LL-SED2-0-56 | 04/07/11 | 0539 | 19.65 | |
| 16 | LL-SED2-56-1 | 04/07/11 | 0615 | 19.65 | |
| 17 | LL-SED2-56-1 | 04/07/11 | 0651 | 19.65 | |
| 18 | LL-SED2-56-1 | 04/07/11 | 0727 | 19.65 | |
| 19 | PCP CCAL | 04/07/11 | 0944 | 19.66 | |
| 20 | LL-SED2-112- | 04/07/11 | 1020 | 19.65 | |
| 21 | LL-SED2-0-56 | 04/07/11 | 1056 | 19.64 | |
| 22 | LL-SED1-0-56 | 04/07/11 | 1132 | 19.64 | |
| 23 | PCP CCAL | 04/07/11 | 1244 | 19.65 | |

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.: SN54 Project: LORA LAKE
 GC Column: ZB35 ID: 0.53 (mm) Instrument ID: ECD1
 Init. Calib. Date(s): 04/01/11 04/01/11

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

| MEAN SURROGATE RT FROM INITIAL CALIBRATION S1 : 20.87 | | | | | |
|--|------------------|------------------|------------------|----------|-------|
| CLIENT SAMPLE NO. | LAB SAMPLE ID | DATE ANALYZED | TIME ANALYZED | S1 RT | # |
| ===== | ===== | ===== | ===== | ===== | ===== |
| 01 | PCP D | 04/01/11 | 1629 | 20.87 | |
| 02 | PCP A | 04/01/11 | 1706 | 20.89 | |
| 03 | PCP B | 04/01/11 | 1742 | 20.89 | |
| 04 | PCP C | 04/01/11 | 1818 | 20.88 | |
| 05 | PCP E | 04/01/11 | 1854 | 20.86 | |
| 06 | PCP F | 04/01/11 | 1931 | 20.86 | |
| 07 | PCP CCAL | 04/06/11 | 1701 | 20.86 | |
| 08 | SN54MBS1 | 04/06/11 | 1737 | 20.87 | |
| 09 | SN54LCSS1 | 04/06/11 | 1813 | 20.87 | |
| 10 | PCP CCAL | 04/07/11 | 0015 | 20.88 | |
| 11 | PCP CCAL | 04/07/11 | 0315 | 20.88 | |
| 12 | LL-SED3-0-36 | 04/07/11 | 0351 | 20.87 | |
| 13 | LL-SED3-36-1 | 04/07/11 | 0427 | 20.88 | |
| 14 | LL-SED3-141- | 04/07/11 | 0503 | 20.87 | |
| 15 | LL-SED2-0-56 | 04/07/11 | 0539 | 20.87 | |
| 16 | LL-SED2-56-1 | 04/07/11 | 0615 | 20.87 | |
| 17 | LL-SED2-56-1 | 04/07/11 | 0651 | 20.87 | |
| 18 | LL-SED2-56-1 | 04/07/11 | 0727 | 20.87 | |
| 19 | PCP CCAL | 04/07/11 | 0944 | 20.88 | |
| 20 | LL-SED2-112- | 04/07/11 | 1020 | 20.87 | |
| 21 | LL-SED2-0-56 | 04/07/11 | 1056 | 20.87 | |
| 22 | LL-SED1-0-56 | 04/07/11 | 1132 | 20.87 | |
| 23 | PCP CCAL | 04/07/11 | 1244 | 20.87 | |

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: SN54

Cover Page

INORGANIC ANALYSIS DATA PACKAGE



CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

| CLIENT ID | ARI ID | ARI LIMS ID | REPREP |
|---------------------|------------|-------------|--------|
| LL-SED3-0-36-03151 | SN54A | 11-5925 | |
| LL-SED3-36-141-031 | SN54B | 11-5926 | , |
| LL-SED3-141-167-03 | SN54C | 11-5927 | |
| LL-SED2-0-56-03151 | SN54D | 11-5928 | |
| LL-SED2-56-112-031 | SN54E | 11-5929 | |
| LL-SED2-56-112-031D | SN54EDUP | 11-5929 | |
| LL-SED2-56-112-031S | SN54ESPK | 11-5929 | |
| LL-SED2-112-168-03 | SN54F | 11-5930 | |
| PBS | SN54MB1 | 11-5930 | |
| LCSS | SN54MB1SPK | 11-5930 | |
| LL-SED2-0-56-03151 | SN54G | 11-5931 | |
| LL-SED1-0-56-03151 | SN54H | 11-5932 | |

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: Jay Kuhn

Name: Jay Kuhn

Date: 3/23/11

Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED3-0-36-031511

SAMPLE

Lab Sample ID: SN54A


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5925

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 11.3%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 40 | 70 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 450 | |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED3-36-141-031511
SAMPLE

Lab Sample ID: SN54B


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5926

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 12.0%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 40 | 40 | U |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 20 | U |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED3-141-167-031511

SAMPLE

Lab Sample ID: SN54C

QC Report No: SN54-Floyd Snider

LIMS ID: 11-5927

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 11.8%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 40 | 50 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 20 | U |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED2-0-56-031511
SAMPLE

Lab Sample ID: SN54D


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5928

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 8.7%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 60 | 70 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 350 | |

U-Analyte undetected at given RL
RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS


Page 1 of 1

Sample ID: LL-SED2-56-112-031511
SAMPLE

Lab Sample ID: SN54E

LIMS ID: 11-5929

Matrix: Sediment

Data Release Authorized 

Reported: 03/23/11

QC Report No: SN54-Floyd Snider

Project: Lora Lake - Subsurface Sediment

POS-LL

Date Sampled: 03/15/11

Date Received: 03/16/11

Percent Total Solids: 13.0%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 40 | 80 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 10 | 10 | U |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

**Sample ID: LL-SED2-56-112-031511
MATRIX SPIKE**

Lab Sample ID: SN54E


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5929

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

MATRIX SPIKE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Sample | Spike | Spike Added | % Recovery | Q |
|----------------|------------------------|---------------|--------------|--------------------|-------------------|----------|
| Arsenic | 6010B | 80 | 1,620 | 1,440 | 107% | |
| Lead | 6010B | 10 U | 1,460 | 1,440 | 101% | |

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-SED2-56-112-031511
 DUPLICATE

Lab Sample ID: SN54E
 LIMS ID: 11-5929
 Matrix: Sediment
 Data Release Authorized 
 Reported: 03/23/11

QC Report No: SN54-Floyd Snider
 Project: Lora Lake - Subsurface Sediment
 POS-LL
 Date Sampled: 03/15/11
 Date Received: 03/16/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Sample | Duplicate | RPD | Control Limit | Q |
|---------|-----------------|--------|-----------|-------|---------------|---|
| Arsenic | 6010B | 80 | 120 | 40.0% | +/- 40 | L |
| Lead | 6010B | 10 U | 10 U | 0.0% | +/- 10 | 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: LL-SED2-112-168-031511

SAMPLE

Lab Sample ID: SN54F


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5930

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 12.2%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 40 | 80 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 20 | U |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED2-0-56-031511-D
SAMPLE

Lab Sample ID: SN54G

LIMS ID: 11-5931

Matrix: Sediment

Data Release Authorized: *MJ*

Reported: 03/23/11

QC Report No: SN54-Floyd Snider

Project: Lora Lake - Subsurface Sediment

POS-LL

Date Sampled: 03/15/11

Date Received: 03/16/11

Percent Total Solids: 10.3%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 50 | 80 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 20 | 380 | |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LL-SED1-0-56-031511

SAMPLE

Lab Sample ID: SN54H


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5932

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: 03/15/11

Reported: 03/23/11

Date Received: 03/16/11

Percent Total Solids: 65.1%

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 7 | 9 | |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 3 | 29 | |

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: SN54LCS


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5930

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: NA

Reported: 03/23/11

Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

| Analyte | Analysis Method | Spike Found | Spike Added | % Recovery | Q |
|----------------|------------------------|--------------------|--------------------|-------------------|----------|
| Arsenic | 6010B | 204 | 200 | 102% | |
| Lead | 6010B | 200 | 200 | 100% | |

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: SN54MB


QC Report No: SN54-Floyd Snider

LIMS ID: 11-5930

Project: Lora Lake - Subsurface Sediment

Matrix: Sediment

POS-LL

Data Release Authorized: 

Date Sampled: NA

Reported: 03/23/11

Date Received: NA

Percent Total Solids: NA

| Prep Meth | Prep Date | Analysis Method | Analysis Date | CAS Number | Analyte | RL | mg/kg-dry | Q |
|-----------|-----------|-----------------|---------------|------------|---------|----|-----------|---|
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7440-38-2 | Arsenic | 5 | 5 | U |
| 3050B | 03/21/11 | 6010B | 03/22/11 | 7439-92-1 | Lead | 2 | 2 | U |

U-Analyte undetected at given RL

RL-Reporting Limit

Calibration Verification

CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

UNITS:ug/L

SDG: SN54

| ANALYTE | EL | M | RUN | ICVTV | ICV | %R | CCVTV | CCV1 | %R | CCV2 | %R | CCV3 | %R | CCV4 | %R | CCV5 | %R |
|---------|----|-----|----------|--------|---------|-------|--------|---------|-------|---------|-------|---------|------|---------|------|---------|-------|
| Arsenic | AS | ICP | IP032271 | 2000.0 | 2041.47 | 102.1 | 2000.0 | 2030.01 | 101.5 | 2025.39 | 101.3 | 1990.25 | 99.5 | 1990.20 | 99.5 | 2020.93 | 101.0 |
| Lead | PB | ICP | IP032271 | 2000.0 | 1963.04 | 98.2 | 2000.0 | 1945.12 | 97.3 | 1948.85 | 97.4 | 1921.07 | 96.1 | 1964.92 | 98.2 | 1952.91 | 97.6 |

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

Calibration Verification



CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

UNITS: ug/L

| ANALYTE | EL | M | RUN | CCVTV | CCV6 | %R | CCV7 | %R | CCV8 | %R | CCV9 | %R | CCV10 | %R | CCV11 | %R |
|---------|----|-----|----------|--------|---------|-------|---------|-------|---------|------|---------|------|-------|----|-------|----|
| Arsenic | AS | ICP | IP032271 | 2000.0 | 2003.51 | 100.2 | 2010.32 | 100.5 | 1996.69 | 99.8 | 1979.69 | 99.0 | | | | |
| Lead | PB | ICP | IP032271 | 2000.0 | 1946.37 | 97.3 | 1936.73 | 96.8 | 1925.16 | 96.3 | 1949.07 | 97.5 | | | | |

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

CRDL Standard

CLIENT: Floyd Snider
 PROJECT: Lora Lake - Subsurfa
 SDG: SN54



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 | IP032271 | 50.0 | | 51.05 | 102.1 | 50.19 | 100.4 | | | | | | | | |
| Lead | PB | ICP | IP032271 | 20.0 | | 20.46 | 102.3 | 20.07 | 100.4 | | | | | | | | |

SN54 : 00100

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

Calibration Blanks

CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54



UNITS: ug/L

| ANALYTE | EL METH | RUN | CRDL | IDL | ICB | CCB1 | CCB2 | CCB3 | CCB4 | CCB5 |
|---------|---------|----------|------|------|------|------|------|------|------|------|
| Arsenic | AS ICP | IP032271 | 10.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 |
| Lead | PB ICP | IP032271 | 3.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 |

Calibration Blanks

CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54



UNITS: ug/L

| ANALYTE | EL | METH | RUN | CRDL | IDL | CCB6 | CCB7 | CCB8 | CCB9 | CCB10 | CCB11 | C |
|---------|----|------|----------|------|------|------|------|------|------|-------|-------|---|
| Arsenic | AS | ICP | IP032271 | 10.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 | 50.0 | U |
| Lead | PB | ICP | IP032271 | 3.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 | 20.0 | U |

SN54 : 00102

ICP Interference Check Sample



CLIENT: Floyd Snider

ICS SOURCE: I.V.

PROJECT: Lora Lake - Subsurfa

RUNID: IP032271

SDG: SN54

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 | 202617.0 | 203157.9 | 101.6 | 202283.2 | 202085.4 | 101.0 | | | |
| Antimony | 1000 | 1000 | 3.1 | 1040.0 | 104.0 | 4.2 | 1027.4 | 102.7 | | | |
| Arsenic | 1000 | 1000 | 13.0 | 1018.8 | 101.9 | 15.7 | 1004.9 | 100.5 | | | |
| Barium | 1000 | 1000 | 0.3 | 983.7 | 98.4 | 1.0 | 964.7 | 96.5 | | | |
| Beryllium | 1000 | 1000 | 0.1 | 1004.8 | 100.5 | 0.0 | 980.8 | 98.1 | | | |
| Boron | | | -6.2 | -5.9 | | -6.6 | | -6.3 | | | |
| Cadmium | 1000 | 1000 | 1.1 | 995.3 | 99.5 | 1.3 | 997.6 | 99.8 | | | |
| Calcium | 100000 | 100000 | 102961.0 | 102500.8 | 102.5 | 101481.1 | 100984.3 | 101.0 | | | |
| Chromium | 1000 | 1000 | -0.7 | 1012.9 | 101.3 | 0.3 | 997.7 | 99.8 | | | |
| Cobalt | 1000 | 1000 | 1.3 | 925.2 | 92.5 | 1.2 | 923.9 | 92.4 | | | |
| Copper | 1000 | 1000 | -0.4 | 1006.9 | 100.7 | 0.0 | 1044.2 | 104.4 | | | |
| Iron | 200000 | 200000 | 200623.7 | 200411.7 | 100.2 | 197183.3 | 196283.1 | 98.1 | | | |
| Lead | 1000 | 1000 | -3.0 | 952.7 | 95.3 | -1.8 | 942.5 | 94.3 | | | |
| Magnesium | 100000 | 100000 | 105993.8 | 102095.3 | 102.1 | 104500.5 | 100418.6 | 100.4 | | | |
| Manganese | 1000 | 1000 | 0.7 | 939.9 | 94.0 | 0.7 | 920.5 | 92.1 | | | |
| Molybdenum | | | 4.3 | 4.9 | | 4.3 | | 4.5 | | | |
| Nickel | 1000 | 1000 | 3.6 | 974.6 | 97.5 | 3.1 | 953.0 | 95.3 | | | |
| Potassium | | | 3.6 | 549.3 | | 24.3 | | 533.4 | | | |
| Selenium | 1000 | 1000 | 29.6 | 1026.2 | 102.6 | 24.0 | 1013.5 | 101.4 | | | |
| Silicon | | | -6.3 | -6.3 | | -8.7 | | -8.4 | | | |
| Silver | 1000 | 1000 | -0.2 | 1009.7 | 101.0 | -0.2 | 1031.8 | 103.2 | | | |
| Sodium | | | 9.9 | 13.8 | | 6.2 | | 10.3 | | | |
| Strontium | | | 3.7 | 3.7 | | 3.7 | | 3.7 | | | |
| Thallium | 1000 | 1000 | -0.9 | 932.2 | 93.2 | 0.6 | 923.3 | 92.3 | | | |
| Tin | | | 0.1 | -0.2 | | -0.9 | | -0.6 | | | |
| Titanium | | | -1.1 | 0.4 | | -0.3 | | 0.0 | | | |
| Vanadium | 1000 | 1000 | 0.4 | 964.1 | 96.4 | 0.5 | 989.4 | 98.9 | | | |
| Zinc | 1000 | 1000 | 3.4 | 983.4 | 98.3 | 1.8 | 960.7 | 96.1 | | | |

FORM IV

SN54 : 00103

IDLs and ICP Linear Ranges



CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

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 | 4/1/2010 | 30000.0 | 2/3/2011 |
| Lead | PB | ICP | OPTIMA ICP 2 | 220.35 | | 3 | 20.0 | 4/1/2010 | 300000.0 | 2/3/2011 |

ICP Interelement Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

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 | 0.0000000 |
| Cobalt | 228.62 | 0.0000000 | 0.0000000 | 0.0227510 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | -0.0451581 |
| Copper | 324.75 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | -0.1969920 | -0.0283867 | 0.0000000 | 0.0000000 |
| 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 | -1.9460200 | 1.1789000 | 0.0588763 |
| Magnesium | 279.08 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | -0.8349460 | 0.0000000 | 0.4579600 |
| Manganese | 257.61 | 0.0056707 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | -0.0082982 |
| Molybdenum | 202.03 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0185703 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 |
| 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 | 2.4468400 | 0.3572340 | 0.0000000 | -0.1350510 |
| Titanium | 334.90 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 |
| Vanadium | 292.40 | 0.0000000 | 0.0000000 | 0.0000000 | 0.0000000 | 0.1735400 | 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 Interement Correction Factors



CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

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 | 232.40 | 0.000000 | -0.161968 | -0.951633 | 0.000000 | 0.000000 | 0.000000 | 0.620997 | 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 |

FORM XI

Preparation Log



CLIENT: Floyd Snider

ANALYSIS METHOD: ICP

PROJECT: Lora Lake - Subsurfa

ARI PREP CODE: SWC

SDG: SN54

PREPDATE: 3/21/2011

| CLIENT ID | ARI ID | MASS (g) | INITIAL VOLUME (mL) | FINAL VOLUME (mL) |
|---------------------|------------|----------|---------------------|-------------------|
| LL-SED3-0-36-03151 | SN54A | 1.037 | 0.0 | 50.0 |
| LL-SED3-36-141-031 | SN54B | 1.021 | 0.0 | 50.0 |
| LL-SED3-141-167-03 | SN54C | 1.082 | 0.0 | 50.0 |
| LL-SED2-0-56-03151 | SN54D | 1.022 | 0.0 | 50.0 |
| LL-SED2-56-112-031 | SN54E | 1.068 | 0.0 | 50.0 |
| LL-SED2-56-112-031D | SN54EDUP | 1.066 | 0.0 | 50.0 |
| LL-SED2-56-112-031S | SN54ESPK | 1.071 | 0.0 | 50.0 |
| LL-SED2-112-168-03 | SN54F | 1.064 | 0.0 | 50.0 |
| LL-SED2-0-56-03151 | SN54G | 1.040 | 0.0 | 50.0 |
| LL-SED1-0-56-03151 | SN54H | 1.032 | 0.0 | 50.0 |
| PBS | SN54MB1 | 1.000 | 0.0 | 50.0 |
| LCSS | SN54MB1SPK | 1.000 | 0.0 | 50.0 |



Analysis Run Log

CLIENT: Floyd Snider

PROJECT: Lora Lake - Subsurfa

SDG: SN54

INSTRUMENT ID: OPTIMA ICP 2

RUNID: IP032271 METHOD: ICP

START DATE: 3/22/2011

END DATE: 3/22/2011

| 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 | S0 | 1.00 | 11084 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| S2 | S2 | 1.00 | 11125 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| S3 | S3 | 1.00 | 11143 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| S4 | S4 | 1.00 | 11170 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| S5 | S5 | 1.00 | 11191 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ICV | ICV | 1.00 | 11233 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X | |
| ICB | ICB | 1.00 | 11264 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ZZZZZ | ZZZZZ | 1.00 | 11304 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| S0 | S0 | 1.00 | 11345 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCV | CCV1 | 1.00 | 11385 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCB | CCB1 | 1.00 | 11420 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CRI | CRII | 1.00 | 11461 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ICSA | ICSAI | 1.00 | 11500 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ICSAB | ICSABI | 1.00 | 11542 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCV | CCV2 | 1.00 | 11580 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCB | CCB2 | 1.00 | 12011 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ZZZZZ | SN04MB1 | 2.00 | 12052 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN74MB1 | 2.00 | 12092 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN74C-L | 10.00 | 12133 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN74C | 2.00 | 12172 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN04B | 2.00 | 12212 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN04ADUP | 2.00 | 12244 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN04A | 2.00 | 12282 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN04ASP | 2.00 | 12320 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN04APOST | 2.00 | 12350 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| ZZZZZ | SN74MB1SPK | 2.00 | 12385 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| CCV | CCV3 | 1.00 | 12424 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCB | CCB3 | 1.00 | 12455 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CRI | CRIF | 1.00 | 12500 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ICSA | ICSAF | 1.00 | 12540 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ICSAB | ICSABF | 1.00 | 12581 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCV | CCV4 | 1.00 | 13021 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| CCB | CCB4 | 1.00 | 13052 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ZZZZZ | SN04C | 2.00 | 13093 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |
| ZZZZZ | SN04D | 2.00 | 13125 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | X |

**General Chemistry Analysis
Report and Summary QC Forms**

ARI Job ID: SN54

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: *[Signature]*
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11


Client ID: LL-SED3-0-36-031511
ARI ID: 11-5925 SN54A

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|-------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 12.80 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb, 1981 | Percent | 0.146 | 8.32 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Client ID: LL-SED3-36-141-031511
ARI ID: 11-5926 SN54B

| Analyte | Date | Method | Units | RL | Sample |
|------------------------|----------------------|-------------|---------|-------|----------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 12.10 |
| Preserved Total Solids | 03/17/11 031711#1 | EPA 160.3 | Percent | 0.01 | 12.30 |
| Sulfide | 03/21/11 032111#1 | EPA 376.2 | mg/kg | 7.85 | < 7.85 U |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb, 1981 | Percent | 0.198 | 25.7 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized
Reported: 03/31/11

A handwritten signature in black ink, appearing to be 'F. Snider', written over the 'Data Release Authorized' text.

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Client ID: LL-SED3-141-167-031511


ARI ID: 11-5927 SN54C

| Analyte | Date | Method | Units | RL | Sample |
|------------------------|----------------------|------------|---------|-------|----------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 12.70 |
| Preserved Total Solids | 03/17/11 031711#1 | EPA 160.3 | Percent | 0.01 | 12.30 |
| Sulfide | 03/21/11 032111#1 | EPA 376.2 | mg/kg | 7.91 | < 7.91 U |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb,1981 | Percent | 0.200 | 21.1 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11


Client ID: LL-SED2-0-56-031511
ARI ID: 11-5928 SN54D

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|-------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 7.00 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb, 1981 | Percent | 0.198 | 17.2 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Client ID: LL-SED2-56-112-031511


ARI ID: 11-5929 SN54E

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 13.80 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb,1981 | Percent | 0.216 | 26.1 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11


Client ID: LL-SED2-112-168-031511
ARI ID: 11-5930 SN54F

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|-------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 11.90 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb, 1981 | Percent | 0.202 | 22.9 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

Client ID: LL-SED2-0-56-031511-D
ARI ID: 11-5931 SN54G

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 5.70 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb,1981 | Percent | 0.200 | 12.1 |

RL Analytical reporting limit
U Undetected at reported detection limit

SAMPLE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized:
Reported: 03/31/11

A handwritten signature in black ink, appearing to be 'F. Snider', written over the 'Data Release Authorized' text.

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11


Client ID: LL-SED1-0-56-031511
ARI ID: 11-5932 SN54H

| Analyte | Date | Method | Units | RL | Sample |
|----------------------|----------------------|-------------|---------|-------|--------|
| Total Solids | 03/22/11 032211#1 | EPA 160.3 | Percent | 0.01 | 41.30 |
| Total Organic Carbon | 03/29/11 032911#1 | Plumb, 1981 | Percent | 0.020 | 4.22 |

RL Analytical reporting limit
U Undetected at reported detection limit

MS/MSD RESULTS-CONVENTIONALS
SN54-Floyd Snider




Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

| Analyte | Date | Units | Sample | Spike | Spike Added | Recovery |
|--|----------|---------|--------|-------|-------------|----------|
| ARI ID: SN54C Client ID: LL-SED3-141-167-031511 | | | | | | |
| Sulfide | 03/21/11 | mg/kg | < 7.91 | 729 | 1,160 | 62.8% |
| ARI ID: SN54E Client ID: LL-SED2-56-112-031511 | | | | | | |
| Total Organic Carbon | 03/29/11 | Percent | 26.1 | 57.6 | 31.4 | 100.4% |

REPLICATE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: 03/15/11
Date Received: 03/16/11

| Analyte | Date | Units | Sample | Replicate(s) | RPD/RSD |
|--|----------|---------|--------|----------------|---------|
| ARI ID: SN54B Client ID: LL-SED3-36-141-031511 | | | | | |
| Preserved Total Solids | 03/17/11 | Percent | 12.30 | 12.40 | 0.8% |
| ARI ID: SN54C Client ID: LL-SED3-141-167-031511 | | | | | |
| Sulfide | 03/21/11 | mg/kg | < 7.91 | < 7.96 | NA |
| ARI ID: SN54E Client ID: LL-SED2-56-112-031511 | | | | | |
| Total Solids | 03/22/11 | Percent | 13.80 | 13.80 13.90 | 0.4% |
| Total Organic Carbon | 03/29/11 | Percent | 26.1 | 22.2 18.9 | 16.1% |

LAB CONTROL RESULTS-CONVENTIONALS
SN54-Floyd Snider




Matrix: Sediment
Data Release Authorized: *[Signature]*
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: NA
Date Received: NA

| Analyte/Method | QC ID | Date | Units | LCS | Spike Added | Recovery |
|-------------------------------------|-------|----------|---------|-------|-------------|----------|
| Sulfide EPA 376.2 | PREP | 03/21/11 | mg/kg | 7.30 | 7.28 | 100.3% |
| Total Organic Carbon Plumb, 1981 | ICVL | 03/29/11 | Percent | 0.094 | 0.100 | 94.0% |

METHOD BLANK RESULTS-CONVENTIONALS
SN54-Floyd Snider




Matrix: Sediment
Data Release Authorized 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: NA
Date Received: NA

| Analyte | Date | Units | Blank |
|------------------------|----------|---------|-----------|
| Total Solids | 03/22/11 | Percent | < 0.01 U |
| Preserved Total Solids | 03/17/11 | Percent | < 0.01 U |
| Sulfide | 03/21/11 | mg/kg | < 1.00 U |
| Total Organic Carbon | 03/29/11 | Percent | < 0.020 U |

STANDARD REFERENCE RESULTS-CONVENTIONALS
SN54-Floyd Snider



Matrix: Sediment
Data Release Authorized: 
Reported: 03/31/11

Project: Lora Lake - Subsurface Sedim
Event: POS-LL
Date Sampled: NA
Date Received: NA

| Analyte/SRM ID | Date | Units | SRM | True Value | Recovery |
|------------------------------------|----------|---------|------|------------|----------|
| Total Organic Carbon NIST 1941B | 03/29/11 | Percent | 2.44 | 2.99 | 81.6% |

Total Solids

ARI Job ID: SN54

Extractions Total Solids-exttts
Data By: Yen Luu
Created: 3/17/11

Worklist: 6028
Analyst: RVR
Comments:

Oven ID: _____

Balance ID: _____

Samples In: Date: _____ Time: _____ Temp: _____ Analyst: _____

Samples Out: Date: _____ Time: _____ Temp: _____ Analyst: _____

| ARI ID CLIENT ID | Tare Wt (g) | Wet Wt (g) | Dry Wt (g) | % Solids | pH |
|---|----------------|---------------|---------------|----------|----|
| 1. SN54A 11-5925 LL-SED3-0-36-031511 | 1.18 | 3.58 | 1.77 | 24.6 | NR |
| 2. SN54B 11-5926 LL-SED3-36-141-031511 | 1.18 | 11.12 | 2.34 | 11.7 | NR |
| 3. SN54C 11-5927 LL-SED3-141-167-031511 | 1.17 | 11.44 | 2.41 | 12.1 | NR |
| 4. SN54D 11-5928 LL-SED2-0-56-031511 | 1.16 | 4.36 | 1.61 | 14.1 | NR |
| 5. SN54E 11-5929 LL-SED2-56-112-031511 | 1.17 | 12.14 | 2.68 | 13.8 | NR |
| 6. SN54F 11-5930 LL-SED2-112-168-031511 | 1.16 | 11.79 | 2.47 | 12.3 | NR |
| 7. SN54G 11-5931 LL-SED2-0-56-031511-D | 1.19 | 3.58 | 1.46 | 11.3 | NR |
| 8. SN54H 11-5932 LL-SED1-0-56-031511 | 1.16 | 3.64 | 2.51 | 54.4 | NR |

SN54 : 00126

Extractions Total Solids-exttts
Data By: Yen Luu
Created: 3/17/11

Worklist: 6028
Analyst: YL
Comments:

Oven ID: 015

Balance ID: 21754520

Samples In: Date: 3/23/11 Time: 19:25 Temp: 102 Analyst: WC

Samples Out: Date: 3/24/11 Time: 09:10 Temp: 105° Analyst: RF

| ARI ID CLIENT ID | Tare Wt (g) | Wet Wt (g) | Dry Wt (g) | % Solids | pH |
|---|----------------|---------------|---------------|----------|----|
| 1. SN54A 11-5925 LL-SED3-0-36-031511 | <u>1.18</u> | <u>3.58g</u> | <u>1.77</u> | | NR |
| 2. SN54B 11-5926 LL-SED3-36-141-031511 | <u>1.18g</u> | <u>11.12g</u> | <u>2.34</u> | | NR |
| 3. SN54C 11-5927 LL-SED3-141-167-031511 | <u>1.17g</u> | <u>11.44g</u> | <u>2.41</u> | | NR |
| 4. SN54D 11-5928 LL-SED2-0-56-031511 | <u>1.16g</u> | <u>4.36g</u> | <u>1.61</u> | | NR |
| 5. SN54E 11-5929 LL-SED2-56-112-031511 | <u>1.17g</u> | <u>12.14g</u> | <u>2.68</u> | | NR |
| 6. SN54F 11-5930 LL-SED2-112-168-031511 | <u>1.16g</u> | <u>11.79g</u> | <u>2.47</u> | | NR |
| 7. SN54G 11-5931 LL-SED2-0-56-031511-D | <u>1.19g</u> | <u>3.58g</u> | <u>1.46</u> | | NR |
| 8. SN54H 11-5932 LL-SED1-0-56-031511 | <u>1.16g</u> | <u>3.64g</u> | <u>2.51</u> | | NR |

Solids Data Entry Report
Date: 03/22/11

Checked by: KM Date: 3/22/11
Data Analyst: DM

Solids Determination performed on 03/21/11 by KM

| JOB | SAMPLE | CLIENTID | TAREWEIGHT | SAMPDISH | DRYWEIGHT | SOLIDS |
|------|--------|---------------------|------------|----------|-----------|--------|
| SN54 | A | LL-SED3-0-36-031511 | 1.005 | 10.326 | 2.055 | 11.26 |
| SN54 | B | LL-SED3-36-141-0315 | 0.973 | 10.574 | 2.128 | 12.03 |
| SN54 | C | LL-SED3-141-167-031 | 0.965 | 10.092 | 2.043 | 11.81 |
| SN54 | D | LL-SED2-0-56-031511 | 0.991 | 10.236 | 1.796 | 8.71 |
| SN54 | E | LL-SED2-56-112-0315 | 0.993 | 10.314 | 2.200 | 12.95 |
| SN54 | F | LL-SED2-112-168-031 | 0.977 | 10.205 | 2.104 | 12.21 |
| SN54 | G | LL-SED2-0-56-031511 | 0.947 | 10.622 | 1.946 | 10.33 |
| SN54 | H | LL-SED1-0-56-031511 | 0.952 | 10.096 | 6.902 | 65.07 |

SN54 : 00128



Total Solids Bench Sheet

Laboratory Section Metals

Oven Identification: 07

Balance ID: 068755

Samples in Oven: Date: 3/21/11 Time: 1040 Temp: 102°C Analyst: KM

Removed from Oven: Date: 3-22-11 Time: 0910 Temp: 103°C Analyst: DM

Source of Total Solids Data If From A Different Lab: _____

| ARI Sample ID | Tare Weight (g) | Tare + Sample Wet (g) | Tare + Sample Dry (g) | Date & Time Last Weight | Final Weighting >12 hrs ¹ |
|--|-----------------|-----------------------|-----------------------|-------------------------|--------------------------------------|
| SN80 E | 0.960 | 10.101 | 9.052 | - | ✓ |
| SN54 A | 1.005 | 10.326 | 2.055 | - | ✓ |
| " B | 0.973 | 10.574 | 2.128 | - | ✓ |
| " C | 0.965 | 10.092 | 2.043 | - | ✓ |
| " D | 0.991 | 10.236 | 1.796 | - | ✓ |
| " E | 0.993 | 10.314 | 2.200 | - | ✓ |
| " F | 0.977 | 10.205 | 2.104 | - | ✓ |
| " G | 0.947 | 10.622 | 1.946 | - | ✓ |
| " H | 0.952 | 10.096 | 6.902 | - | ✓ |
| <div style="position: absolute; top: 50%; left: 50%; transform: translate(-50%, -50%); opacity: 0.5;"> <p>KM</p> <p>3/21/11</p> </div> | | | | | |

1) Place a check mark in this column if samples have dried > 12 but < 24 hours. When samples have been at 104°C < 12 hours, constant weight must be verified as described in SOP 10023S. Use a 2nd bench sheet for additional weightings.

**SIM PAH Raw Data
Extraction Bench Sheets and Notes**

ARI Job ID: SN54



Preparation Test SIM PNA # 1

ARI Job No(s) SN89, SN54

In-House (5ppb)

Batch set up by: SH

| Bottle # | Extraction Requirements | Verify Client ID | Volume Extracted (dry wt) | Sonic Horn ID + Check | KD 80-85°C | | TurboVap ①②③ | (REQ) | TurboVap ①②③ | Final Effective Volume | Volume to Lab | Comments |
|----------|-------------------------|---------------------|---------------------------|-----------------------|-----------------|----------------------------------|-----------------|-------|-----------------|------------------------|---------------|-----------|
| | | | | | Hex X X 2 100°C | (Opt) Silica Gel Clean (1:1) Y N | | | | | | |
| | <u>SN89</u> MBS | Date <u>3/26/11</u> | 10.00g | 3 | | | | | | 0.5mL | 0.5mL | |
| | SBS | | ↓ | 4 | | | | | | ↓ | ↓ | |
| | SBS Dup. | | ↓ | 7 | | | | | | ↓ | ↓ | |
| | QLS | | ↓ | 8 | | | | | | ↓ | ↓ | |
| 1 | A | checked | 7.15 | 9 | | | | | | | | see notes |
| 1 | B | | 14.25 | 10 | | | | | | | | ↓ |
| 2 | <u>SN54</u> A | | 24.07 | 11 | | | | | | | | ↓ |
| 4 | B | | 86.10 | 12 | | | | | | | | |
| 4 | C | | 83.15 | 3 | | | | | | | | |
| 2 | D | | 23.35 | 4 | | | | | | | | see notes |
| 3 | E | | 73.52 | 7 | | | | | | | | |
| 5 | Ems | | 73.08 | 8 | | | | | | | | |
| 5 | Emsd | | 73.29 | 9 | | | | | | | | |
| 3 | F | | 82.00 | 10 | | | | | | | | |
| 3 | G | | 19.06 | 11 | | | | | | | | see notes |
| 2 | H | | 19.41 | 12 | ↓ | ↓ | ↓ | ↓ | ↓ | ↓ | ↓ | |

Analyst/Date AR 03/26/11 → AL 3-28-11 RP/YL 3/29/11 ES 3-30-11

| Standard | Standard ID | Volume | Expiration Date | Analyst | Witness |
|-----------|----------------|--------|-----------------|---------|---------|
| Surrogate | B ₂ | 100µL | 10/25/11 | AR | TH |
| Spike | 15 | 100µL | 11/4/12 | AR | TH |
| QLS Spike | 4 | 50µL | 11/4/12 | AR | TH |

Extraction Time: 15:20 Balance ID: 2175452a

SPECIAL INSTRUCTIONS: 1. Weigh into 100mL beakers. 2. Extract 2X with 1:1 DCM/Acetone. Plus 1 X DCM only. 3. Collect into 200mL E-Flask with 5-10g sodium sulfate in the bottom + small funnel with neutral glasswool. **NO SODIUM SULFATE IN FUNNEL.** 4. KD (small drying column) to 8mL at 80-85°. 5. Exchange (2 X with 10mL) to Hexane at 100°. 6. TurboVap. 7. Silica Clean-up ~~Opt~~-Any color=REQUIRED (All ~~opt~~). 8. TurboVap (if Silica Clean). 9. Vial in DCM.

A. Need Total Solids Y (N) B. Archive (Freeze Y) N

SN89 only



ARI Job No.: SN54

Client ID: Floyd Snider

Parameter: SIM PNA

Client Project: Lora lake-Subsurface Sediment

| Note problems, concerns, corrective actions | Analyst/Date |
|--|------------------------|
| Screens: Soil/Sediment/Solid/Other: | |
| <input type="checkbox"/> No Anomalies (standard soil/sediment) | |
| <input checked="" type="checkbox"/> Wet sediment/sludge= <u>Wet (A-H)</u> | |
| <input type="checkbox"/> Standing Water Decanted= | |
| <input type="checkbox"/> Standing Water Homogenized (Shared samples)= | |
| <input type="checkbox"/> Clay (Difficult to homogenize/Mixed with Kitchen Aid)= | |
| <input type="checkbox"/> Rocks/Organics= | |
| <input type="checkbox"/> Oily, obvious fuel/sulfur odors= | |
| <input checked="" type="checkbox"/> Other (Details)= <u>Samples A, B and C are provided in amber jars.</u> | <u>WC 3/23/11</u> ↓ |
| Aqueous: | |
| <input type="checkbox"/> No Anomalies | |
| <input type="checkbox"/> Turbid/Color= | |
| <input type="checkbox"/> Particulates= | |
| <input type="checkbox"/> Emulsions= | |
| <input type="checkbox"/> Other (Details)= | |
| <input checked="" type="checkbox"/> Other Notes/Comments= <u>Samples A, D, G and H were received as mostly water (> 80%). As per project manager, volume for these samples was taken from the metals lab. Two jars each were centrifuged and combined by the Geotech lab.</u> | <u>JH 3/26/11</u> |
| <u>GCMS analyst, reduced extraction weights for samples A, D and G, based on sample pre-screens.</u> | <u>JH 3/26/11</u> |
| <u>At vialing filtered samples SN54 A-H with GDX stacked with D45, split samples SN54 A-H at ~ 1.25ml for SPE clean up, samples were dark and viscous, after clean up combined samples into turbo tubes for final blow down.</u> | <u>3-30-11 TS</u> |

**SIM PAH Raw Data
Initial Calibration**

ARI Job ID: SN54



GC/MS SVOA Analyst Notes / Corrective Action Log

ARI Project ID: CUSM Client ID: _____

ARI SOP: 801S(SIM-PNA) 802S(Butyl Tins) 804S(SVOA-8270D) 805S(op-Pest)

Parameter(s): Sim-PNA

Instrument: NT-2 NT-4 NT-6 NT-8 NT11, (NT12)

Curve Date: 3/31/11 Analysis Start Date: 3/31/11

| | | | |
|--------------------------------|----------------------|-----------------------------------|----------------------|
| DFTPP Tune Meets Criteria? | <u>YES</u> / NO | Internal Standard Meets Criteria? | <u>YES</u> / NO |
| DDT Breakdown <20%? | <u>YES</u> / NO / NA | Method Blank In Control? | YES / NO <u>(NA)</u> |
| Peak Tailing Factor ≤2? | <u>YES</u> / NO / NA | LCS / LCSD Recovery In Control? | <u>YES</u> / NO |
| ICal acceptable? | <u>YES</u> / NO | CCal acceptable? | <u>YES</u> / NO |
| Q flag applied? | YES / NO | Q flag applied? | YES / NO |
| Surrogate Recovery in Control? | <u>YES</u> / NO | Special Analysis Criteria Met? | YES / NO / <u>NA</u> |
| Manual Integrations for ICal? | <u>YES</u> / NO | Manual Integrations for Samples? | Yes / NO <u>(NA)</u> |

Detail problems, corrective actions and/or other pertinent information below (use reverse side when necessary):

*no perylene added in 1CV.
(1CV for perylene will be performed if needed).*

Additional Details on Reverse: Yes / No

Analyst: [Signature] Date: 04/06/11

Reviewer: [Signature] Date: 4/5/11

Analytical Resources, Inc.

INITIAL CALIBRATION DATA

Start Cal Date : 31-MAR-2011 20:27
 End Cal Date : 31-MAR-2011 22:47
 Quant Method : ISTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP RTE
 Method file : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Cal Date : 01-Apr-2011 17:23 jianqing
 Curve Type : Average

Calibration File Names:

- Level 1: /chem1/nt12.i/20110331.b/03311103.d
- Level 2: /chem1/nt12.i/20110331.b/03311104.d
- Level 3: /chem1/nt12.i/20110331.b/03311105.d
- Level 4: /chem1/nt12.i/20110331.b/03311102.d
- Level 5: /chem1/nt12.i/20110331.b/03311106.d
- Level 6: /chem1/nt12.i/20110331.b/03311107.d

B 04/01/11

| Compound | 0.10000 Level 1 | 0.50000 Level 2 | 1.000 Level 3 | 2.500 Level 4 | 5.000 Level 5 | 10.000 Level 6 | RRF | % RSD |
|---------------------------|--------------------|--------------------|------------------|------------------|------------------|-------------------|---------|--------|
| 28 Naphthalene | 1.02984 | 1.04229 | 0.96546 | 0.98620 | 0.93861 | 0.83191 | 0.96572 | 7.891 |
| 32 2-Methylnaphthalene | 0.63113 | 0.57566 | 0.54305 | 0.52380 | 0.50615 | 0.44017 | 0.53666 | 12.045 |
| 105 1-methylnaphthalene | 0.73051 | 0.60922 | 0.56684 | 0.54282 | 0.52262 | 0.45464 | 0.57111 | 16.350 |
| 40 Acenaphthylene | 1.96509 | 1.84469 | 1.78684 | 1.69133 | 1.69760 | 1.51530 | 1.75014 | 8.772 |
| 44 Acenaphthene | 1.32697 | 1.09842 | 1.09926 | 1.02851 | 1.04015 | 0.94739 | 1.09012 | 11.813 |
| 46 Dibenzofuran | 1.75677 | 1.55840 | 1.48369 | 1.41226 | 1.41246 | 1.26907 | 1.48211 | 11.143 |
| 49 Fluorene | 1.42196 | 1.24282 | 1.22580 | 1.16986 | 1.18176 | 1.07992 | 1.22035 | 9.339 |
| 60 Phenanthrene | 1.25697 | 1.09982 | 1.05963 | 1.00114 | 0.98489 | 0.93079 | 1.05554 | 10.890 |
| 61 Anthracene | 1.23908 | 1.09624 | 1.06084 | 1.06111 | 1.01593 | 0.88579 | 1.05983 | 10.814 |
| 64 Fluoranthene | 1.34085 | 1.21122 | 1.14001 | 1.10737 | 1.09067 | 0.99664 | 1.14779 | 10.242 |
| 65 Pyrene | 1.26093 | 1.10447 | 1.07602 | 1.03032 | 1.04090 | 0.94833 | 1.07683 | 9.707 |
| 68 Benzo(a)anthracene | 1.11299 | 0.94652 | 0.94287 | 0.93557 | 0.95613 | 0.87492 | 0.96150 | 8.284 |
| 71 Chrysene | 1.08380 | 0.95515 | 0.93453 | 0.90160 | 0.91903 | 0.84421 | 0.93972 | 8.517 |
| 74 Benzo(b)fluoranthene | 1.22881 | 1.13978 | 1.14663 | 1.09249 | 1.10623 | 1.05109 | 1.12751 | 5.365 |
| 75 Benzo(k)fluoranthene | 1.41742 | 1.20751 | 1.14974 | 1.13283 | 1.11431 | 1.07459 | 1.18273 | 10.400 |
| 188 Benzo(j)fluoranthene | 1.37699 | 1.13978 | 1.08633 | 1.08271 | 1.08914 | 1.00588 | 1.13014 | 11.354 |
| 76 Benzo(a)pyrene | 1.25288 | 1.00999 | 0.99588 | 0.99323 | 1.01691 | 0.94227 | 1.03519 | 10.608 |
| 78 Indeno(1,2,3-cd)pyrene | 1.37479 | 1.31242 | 1.22066 | 1.20629 | 1.25293 | 1.18555 | 1.25877 | 5.722 |
| 79 Dibenzo(a,h)anthracene | 1.25074 | 0.93800 | 0.99289 | 1.00459 | 1.01450 | 0.97604 | 1.02946 | 10.848 |
| 80 Benzo(g,h,i)perylene | 1.36701 | 1.12986 | 1.04412 | 1.05448 | 1.08217 | 1.02109 | 1.11645 | 11.492 |
| 99 Perylene | 1.04566 | 0.88472 | 0.85494 | 0.83735 | 0.83808 | 0.79215 | 0.87548 | 10.122 |

Analytical Resources, Inc.

INITIAL CALIBRATION DATA

Start Cal Date : 31-MAR-2011 20:27
 End Cal Date : 31-MAR-2011 22:47
 Quant Method : ISTD
 Origin : Disabled
 Target Version : 3.50
 Integrator : HP RTE
 Method file : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Cal Date : 01-Apr-2011 17:23 jianqing
 Curve Type : Average

| Compound | 0.10000 Level 1 | 0.50000 Level 2 | 1.000 Level 3 | 2.500 Level 4 | 5.000 Level 5 | 10.000 Level 6 | RRF | % RSD |
|-----------------------------------|--------------------|--------------------|------------------|------------------|------------------|-------------------|---------|--------|
| 58 Pentachlorophenol | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ |
| 187 Total Benzofluoranthenes | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ |
| 152 Benzo(e)pyrene | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ |
| \$ 190 2-Methylnaphthalene-d10 | 0.72188 | 0.66040 | 0.61951 | 0.58941 | 0.57524 | 0.50250 | 0.61149 | 12.308 |
| \$ 191 Dibenzo(a,h)anthracene-d14 | 1.10300 | 0.94416 | 0.88878 | 0.90105 | 0.92161 | 0.87241 | 0.93850 | 8.994 |
| \$ 55 2,4,6-Tribromophenol | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ |

Analytical Resources, Inc.
RETENTION TIME SUMMARY REPORT

Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
Batch File: /chem1/nt12.i/20110331.b
Inst ID: nt12.i

ID: RT01 RT02 RT03 RT04 RT05 RT06
FILENAME: 03311102 03311103 03311104 03311105 03311106 03311107
INJ. DATE: 31-MAR-2011 31-MAR-2011 31-MAR-2011 31-MAR-2011 31-MAR-2011 31-MAR-2011
INJ. TIME: 20:27 20:55 21:23 21:51 22:19 22:47

| Compound | RT01 | RT02 | RT03 | RT04 | RT05 | RT06 | EXPEC RT | RT WINDOW | AVG RT | STD DEV |
|-------------------------------|--------|--------|--------|--------|--------|--------|----------|---------------|--------|---------|
| * 27 Naphthalene-d8 | 4.920 | 4.917 | 4.917 | 4.920 | 4.920 | 4.917 | 4.917 | 4.817-5.017 | 4.919 | 0.002 |
| 28 Naphthalene | 4.945 | 4.945 | 4.949 | 4.945 | 4.948 | 4.945 | 4.945 | 4.845-5.045 | 4.946 | 0.002 |
| \$ 190 2-Methylnaphthalene-d1 | 5.655 | 5.655 | 5.655 | 5.655 | 5.655 | 5.655 | 5.655 | 5.555-5.755 | 5.655 | 0.000 |
| 32 2-Methylnaphthalene | 5.700 | 5.700 | 5.700 | 5.696 | 5.699 | 5.700 | 5.700 | 5.600-5.800 | 5.699 | 0.001 |
| 105 1-methylnaphthalene | 5.892 | 5.889 | 5.892 | 5.892 | 5.892 | 5.892 | 5.892 | 5.792-5.992 | 5.892 | 0.001 |
| 40 Acenaphthylene | 7.025 | 7.025 | 7.025 | 7.025 | 7.022 | 7.025 | 7.025 | 6.925-7.125 | 7.024 | 0.001 |
| * 42 Acenaphthene-d10 | 7.164 | 7.161 | 7.161 | 7.164 | 7.164 | 7.164 | 7.164 | 7.064-7.264 | 7.163 | 0.002 |
| 44 Acenaphthene | 7.208 | 7.211 | 7.208 | 7.208 | 7.208 | 7.208 | 7.208 | 7.108-7.308 | 7.208 | 0.001 |
| 46 Dibenzofuran | 7.353 | 7.350 | 7.353 | 7.353 | 7.353 | 7.353 | 7.353 | 7.253-7.453 | 7.353 | 0.001 |
| 49 Fluorene | 7.808 | 7.804 | 7.808 | 7.807 | 7.807 | 7.808 | 7.808 | 7.708-7.908 | 7.807 | 0.001 |
| * 59 Phenanthrene-d10 | 9.098 | 9.098 | 9.098 | 9.098 | 9.101 | 9.098 | 9.098 | 8.998-9.198 | 9.099 | 0.001 |
| 60 Phenanthrene | 9.133 | 9.130 | 9.130 | 9.133 | 9.133 | 9.133 | 9.133 | 9.033-9.233 | 9.132 | 0.002 |
| 61 Anthracene | 9.164 | 9.164 | 9.164 | 9.164 | 9.164 | 9.164 | 9.164 | 9.064-9.264 | 9.164 | 0.000 |
| 64 Fluoranthene | 10.827 | 10.821 | 10.827 | 10.824 | 10.824 | 10.827 | 10.827 | 10.727-10.927 | 10.825 | 0.003 |
| 65 Pyrene | 11.301 | 11.298 | 11.294 | 11.298 | 11.297 | 11.301 | 11.301 | 11.201-11.401 | 11.298 | 0.002 |
| 68 Benzo (a) anthracene | 13.690 | 13.686 | 13.683 | 13.689 | 13.693 | 13.693 | 13.693 | 13.593-13.793 | 13.689 | 0.004 |
| * 69 Chrysene-d12 | 13.803 | 13.806 | 13.803 | 13.803 | 13.806 | 13.806 | 13.806 | 13.706-13.906 | 13.805 | 0.002 |

Handwritten signature and date: 04/01/11

Reviewer 1
Reviewer 2

Date: _____
Date: _____

Analytical Resources, Inc.
RETENTION TIME SUMMARY REPORT

Method File: /chem1/nt12.1/20110331.b/SIMPNA0331.m
Batch File: /chem1/nt12.1/20110331.b
Inst ID: nt12.1

| Compound | RT01 | RT02 | RT03 | RT04 | RT05 | RT06 | EXPEC RT | RT WINDOW | AVG RT | STD DEV |
|---------------------------------|--------|--------|--------|--------|--------|--------|----------|---------------|--------|---------|
| 71 Chrysene | 13.873 | 13.873 | 13.866 | 13.872 | 13.876 | 13.873 | 13.873 | 13.773-13.973 | 13.872 | 0.003 |
| 74 Benzo (b) Fluoranthene | 16.306 | 16.302 | 16.299 | 16.302 | 16.305 | 16.312 | 16.312 | 16.212-16.412 | 16.304 | 0.004 |
| 75 Benzo (k) Fluoranthene | 16.362 | 16.359 | 16.359 | 16.359 | 16.359 | 16.369 | 16.369 | 16.269-16.469 | 16.361 | 0.004 |
| 188 Benzo (j) Fluoranthene | 16.435 | 16.425 | 16.429 | 16.432 | 16.435 | 16.444 | 16.444 | 16.344-16.544 | 16.433 | 0.007 |
| 76 Benzo (a) pyrene | 17.262 | 17.255 | 17.255 | 17.252 | 17.262 | 17.268 | 17.268 | 17.168-17.368 | 17.259 | 0.006 |
| * 77 Perylene-d12 | 17.470 | 17.464 | 17.464 | 17.467 | 17.470 | 17.473 | 17.473 | 17.373-17.573 | 17.468 | 0.004 |
| 78 Indeno (1,2,3-cd) pyrene | 19.818 | 19.802 | 19.802 | 19.799 | 19.818 | 19.833 | 19.833 | 19.733-19.933 | 19.812 | 0.013 |
| \$ 191 Dibenzo (a,h) anthracene | 19.745 | 19.729 | 19.733 | 19.729 | 19.739 | 19.751 | 19.751 | 19.651-19.851 | 19.738 | 0.009 |
| 79 Dibenzo (a,h) anthracene | 19.827 | 19.818 | 19.821 | 19.821 | 19.830 | 19.840 | 19.840 | 19.740-19.940 | 19.826 | 0.008 |
| 80 Benzo (g,h,i) perylene | 20.670 | 20.648 | 20.653 | 20.654 | 20.656 | 20.676 | 20.676 | 20.576-20.776 | 20.663 | 0.010 |
| \$ 55 2,4,6-Tribromophenol | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | 14.264 | 14.164-14.364 | +++++ | +++++ |
| 99 Perylene | 17.536 | 17.536 | 17.536 | 17.530 | 17.533 | 17.536 | 17.536 | 17.436-17.636 | 17.535 | 0.003 |
| 58 Pentachlorophenol | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | 10.007 | 9.907-10.107 | +++++ | +++++ |
| 187 Total Benzofluoranthene | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | 17.639 | 17.539-17.739 | +++++ | +++++ |
| 152 Benzo (e) pyrene | +++++ | +++++ | +++++ | +++++ | +++++ | +++++ | 30.943 | 30.843-31.043 | +++++ | +++++ |

Analytical Resources Inc.: Organics Instrument Log

NT-12 Serial No.: GC=US00032558, MS= US01180091

Date: 3/31/11 Analysis: SIMPAA Analyst: JB
 GC Program: SIMPAAZF Column No: 167992 Column Type: ZB35
 Instrument Tune (.U or .CT.): 110331 EM Voltage: 2660
 Calibration File: 0331102 Curve Date: 3/31/11 Injection Vol.: 1µl

| IS/SS | Ical/Ccal | LCS/ICV |
|--------|-----------|---------|
| 1754-1 | 1818-1 | |
| | 1788-3 | |
| | | |
| | | |
| | | |
| | | |
| | | |

Document All Maintenance Tasks In StarLIMS

| INTERNAL STANDARD SUMMARY FOR DATABATCH - /chem1/nt12.i/20110331.b | | | | | | | | | | | | | | | |
|--|----------|------------|-----------|-----------|--|------------------|--------|------|--------|------|--------|-------|--------|-------|--------|
| Time | Filename | LabID | ClientId | DF | | | | | | | | | | | |
| 1 | 2014 | 03311101.d | DFTPP0331 | DFTPP0331 | 1 | [NO ISTDs FOUND] | | | | | | | | | |
| 2 | 2027 | 03311102.d | IC0250331 | IC0250331 | 1 | 4.92 | 494112 | 7.16 | 280105 | 9.10 | 461353 | 13.80 | 503160 | 17.47 | 442215 |
| 3 | 2055 | 03311103.d | IC010331 | IC010331 | 1 | 4.92 | 382584 | 7.16 | 224662 | 9.10 | 370050 | 13.81 | 404335 | 17.46 | 354686 |
| 4 | 2123 | 03311104.d | IC050331 | IC050331 | 1 | 4.92 | 399629 | 7.16 | 230116 | 9.10 | 380150 | 13.80 | 423330 | 17.46 | 369841 |
| 5 | 2151 | 03311105.d | IC10331 | IC10331 | 1 | 4.92 | 447156 | 7.16 | 253822 | 9.10 | 425121 | 13.80 | 461408 | 17.47 | 420159 |
| 6 | 2219 | 03311106.d | IC50331 | IC50331 | 1 | 4.92 | 424290 | 7.16 | 233090 | 9.10 | 394000 | 13.81 | 423013 | 17.47 | 385980 |
| 7 | 2247 | 03311107.d | IC100331 | IC10331 | 1 | 4.92 | 479714 | 7.16 | 252063 | 9.10 | 420234 | 13.81 | 454758 | 17.47 | 406204 |
| 8 | 2315 | 03311108.d | IC200331 | IC200331 | 1 | 4.92 | 431209 | 7.16 | 238299 | 9.10 | 401066 | 13.81 | 433254 | 17.47 | 394342 |
| | | | | | <div style="font-size: 2em; font-family: cursive;"> JB 04/01/11 </div> | | | | | | | | | | |
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| | | | | | | | | | | | | | | | |

Every line must contain information or be lined out. Make all entries legible.
 Start a new page for each QC period. Document All Maintenance Tasks In StarLIMS

MANUAL INTEGRATION SUMMARY FOR DATABATCH - /chem1/nt12.i/20110331.b
 ARI Job No.: IC02 Method: SIMPNA0331.m Instrument: nt12.i Date: 31-MAR-2011

SE 09/01/11

| Time | Filename | LabID | ClientID | DF | Manually Integrated Compounds |
|------|------------|-----------|-----------|----|---|
| 2027 | 03311102.d | IC0250331 | IC0250331 | 1 | NO MANUAL INTEGRATION |
| 2055 | 03311103.d | IC010331 | IC010331 | 1 | NO MANUAL INTEGRATION |
| 2123 | 03311104.d | IC050331 | IC050331 | 1 | Benzo (a) anthracene, Benzo (a) pyrene, Dibenzo (a,h) anthracene, |
| 2151 | 03311105.d | IC10331 | IC10331 | 1 | NO MANUAL INTEGRATION |
| 2219 | 03311106.d | IC50331 | IC50331 | 1 | NO MANUAL INTEGRATION |
| 2247 | 03311107.d | IC100331 | IC10331 | 1 | NO MANUAL INTEGRATION |

Date : 31-MAR-2011 20:14

Client ID: DFTPP0331

Instrument: nt12.i

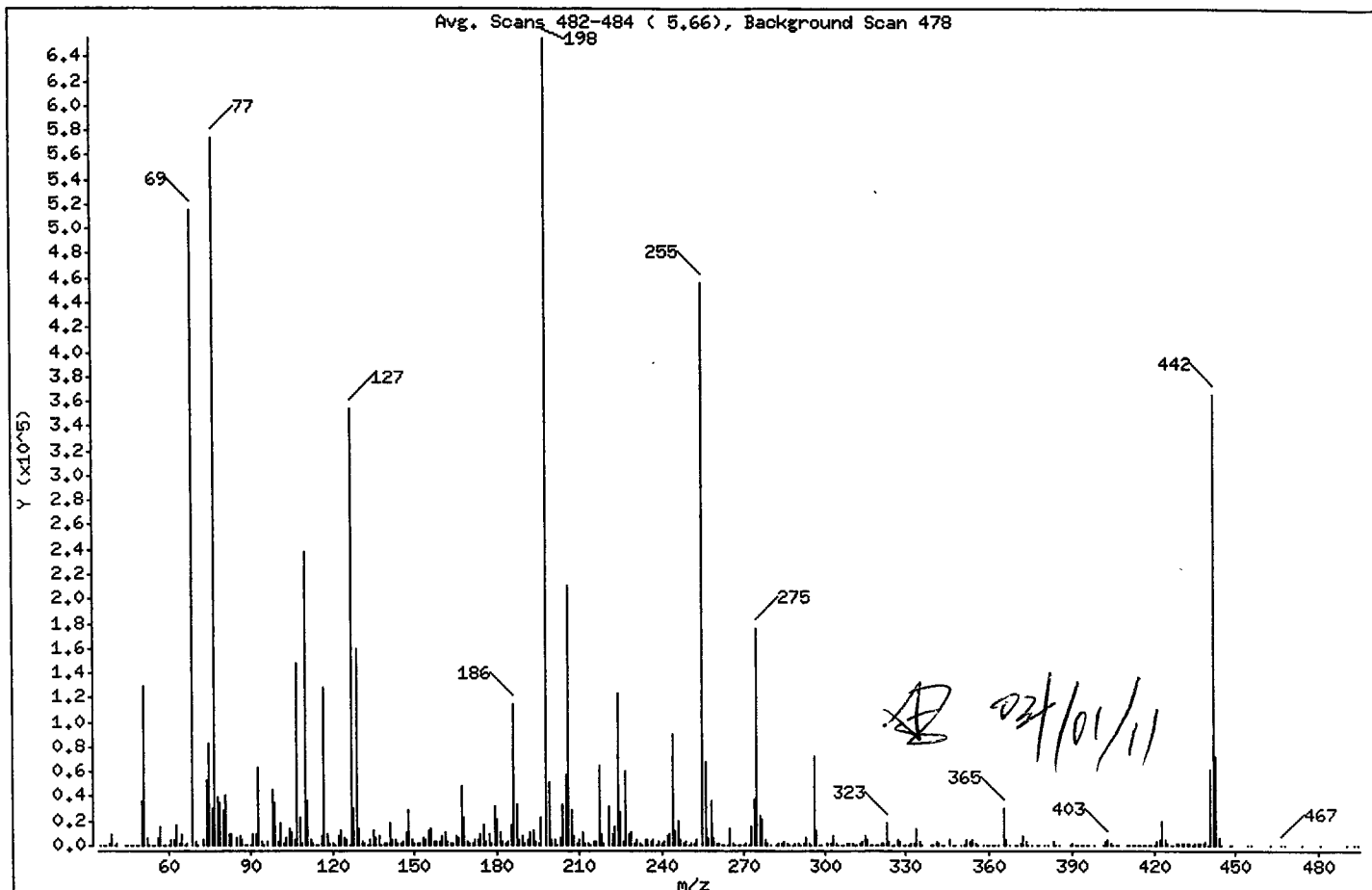
Sample Info: DFTPP0331

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0.32

1 dftpp



| m/e | ION ABUNDANCE CRITERIA | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 198 | Base Peak, 100% relative abundance | 100.00 |
| 51 | 10.00 - 80.00% of mass 198 | 19.69 |
| 68 | Less than 2.00% of mass 69 | 0.00 (0.00) |
| 69 | Mass 69 relative abundance | 78.66 |
| 70 | Less than 2.00% of mass 69 | 0.40 (0.50) |
| 127 | 10.00 - 80.00% of mass 198 | 53.97 |
| 197 | Less than 2.00% of mass 198 | 0.00 |
| 199 | 5.00 - 9.00% of mass 198 | 7.79 |
| 275 | 10.00 - 60.00% of mass 198 | 26.89 |
| 365 | Greater than 1.00% of mass 198 | 4.71 |
| 441 | 0.01 - 24.00% of mass 442 | 9.40 (16.84) |
| 442 | 50.00 - 200.00% of mass 198 | 55.79 |
| 443 | 15.00 - 24.00% of mass 442 | 11.11 (19.92) |

Date : 31-MAR-2011 20:14

Client ID: DFTPP0331

Instrument: nt12.i

Sample Info: DFTPP0331

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0.32

Data File: 03311101.d

Spectrum: Avg. Scans 482-484 (5.66), Background Scan 478

Location of Maximum: 198.00

Number of points: 394

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|-------|--------|--------|-------|--------|--------|--------|-------|
| 35.00 | 612 | 140.00 | 1666 | 241.00 | 3403 | 346.00 | 3950 |
| 36.00 | 181 | 141.00 | 18336 | 242.00 | 7630 | 347.00 | 691 |
| 37.00 | 518 | 142.00 | 5189 | 243.00 | 8406 | 348.00 | 56 |
| 38.00 | 1675 | 143.00 | 4787 | 244.00 | 89752 | 350.00 | 556 |
| 39.00 | 9265 | 144.00 | 1323 | 245.00 | 12268 | 351.00 | 210 |
| 40.00 | 574 | 145.00 | 1048 | 246.00 | 20272 | 352.00 | 5114 |
| 41.00 | 761 | 146.00 | 2945 | 247.00 | 3780 | 353.00 | 3262 |
| 44.00 | 43 | 147.00 | 10618 | 248.00 | 1407 | 354.00 | 4524 |
| 45.00 | 78 | 148.00 | 28280 | 249.00 | 3356 | 355.00 | 1609 |
| 46.00 | 60 | 149.00 | 5053 | 250.00 | 508 | 356.00 | 277 |
| 47.00 | 107 | 150.00 | 1458 | 251.00 | 1606 | 358.00 | 205 |
| 48.00 | 214 | 151.00 | 2111 | 252.00 | 1728 | 359.00 | 461 |
| 49.00 | 523 | 152.00 | 1434 | 253.00 | 4807 | 360.00 | 303 |
| 50.00 | 36520 | 153.00 | 6118 | 255.00 | 456320 | 361.00 | 550 |
| 51.00 | 129064 | 154.00 | 5165 | 256.00 | 68120 | 362.00 | 219 |
| 52.00 | 6002 | 155.00 | 12120 | 257.00 | 5397 | 363.00 | 193 |
| 53.00 | 733 | 156.00 | 13350 | 258.00 | 36192 | 365.00 | 30872 |
| 54.00 | 62 | 157.00 | 3318 | 259.00 | 5278 | 366.00 | 4506 |
| 55.00 | 729 | 158.00 | 3317 | 260.00 | 1306 | 367.00 | 402 |
| 56.00 | 6448 | 159.00 | 3009 | 261.00 | 973 | 369.00 | 293 |
| 57.00 | 15302 | 160.00 | 7531 | 262.00 | 164 | 370.00 | 602 |
| 58.00 | 658 | 161.00 | 10379 | 263.00 | 281 | 371.00 | 947 |
| 59.00 | 164 | 162.00 | 2879 | 264.00 | 535 | 372.00 | 7911 |
| 60.00 | 227 | 163.00 | 1429 | 265.00 | 13069 | 373.00 | 2464 |
| 61.00 | 4122 | 164.00 | 977 | 266.00 | 1736 | 374.00 | 187 |
| 62.00 | 5135 | 165.00 | 7676 | 267.00 | 555 | 375.00 | 156 |
| 63.00 | 16760 | 166.00 | 6736 | 268.00 | 329 | 377.00 | 466 |
| 64.00 | 2178 | 167.00 | 48624 | 269.00 | 493 | 378.00 | 72 |
| 65.00 | 8420 | 168.00 | 23120 | 270.00 | 706 | 380.00 | 72 |
| 66.00 | 430 | 169.00 | 3088 | 271.00 | 1482 | 381.00 | 174 |
| 67.00 | 970 | 170.00 | 1245 | 272.00 | 720 | 383.00 | 2322 |
| 69.00 | 515584 | 171.00 | 1484 | 273.00 | 15391 | 384.00 | 571 |
| 70.00 | 2598 | 172.00 | 4473 | 274.00 | 36928 | 385.00 | 257 |
| 71.00 | 278 | 173.00 | 5014 | 275.00 | 176256 | 389.00 | 134 |
| 73.00 | 4478 | 174.00 | 8950 | 276.00 | 24792 | 390.00 | 1063 |

Date : 31-MAR-2011 20:14

Client ID: DFTPP0331

Instrument: nt12.i

Sample Info: DFTPP0331

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0,32

Data File: 03311101.d

Spectrum: Avg. Scans 482-484 (5.66), Background Scan 478

Location of Maximum: 198.00

Number of points: 394

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|--------|--------|--------|--------|--------|-------|--------|-------|
| 74.00 | 52336 | 175.00 | 17000 | 277.00 | 20608 | 391.00 | 613 |
| 75.00 | 82696 | 176.00 | 3561 | 278.00 | 3839 | 392.00 | 675 |
| 76.00 | 29520 | 177.00 | 8429 | 279.00 | 903 | 393.00 | 155 |
| 77.00 | 574208 | 178.00 | 1888 | 280.00 | 405 | 394.00 | 53 |
| 78.00 | 39336 | 179.00 | 32192 | 282.00 | 611 | 395.00 | 78 |
| 79.00 | 35048 | 180.00 | 21832 | 283.00 | 1994 | 396.00 | 156 |
| 80.00 | 28976 | 181.00 | 11130 | 284.00 | 1528 | 398.00 | 61 |
| 81.00 | 40616 | 182.00 | 2550 | 285.00 | 3123 | 401.00 | 473 |
| 82.00 | 9469 | 183.00 | 920 | 286.00 | 1215 | 402.00 | 3429 |
| 83.00 | 9491 | 184.00 | 2918 | 287.00 | 98 | 403.00 | 4444 |
| 85.00 | 6025 | 185.00 | 16113 | 288.00 | 204 | 404.00 | 1400 |
| 86.00 | 7864 | 186.00 | 114832 | 289.00 | 1493 | 405.00 | 246 |
| 87.00 | 4530 | 187.00 | 33608 | 290.00 | 1150 | 406.00 | 85 |
| 88.00 | 456 | 188.00 | 4155 | 291.00 | 424 | 407.00 | 171 |
| 89.00 | 741 | 189.00 | 7729 | 292.00 | 1282 | 410.00 | 338 |
| 90.00 | 359 | 190.00 | 1679 | 293.00 | 5422 | 411.00 | 59 |
| 91.00 | 8697 | 191.00 | 3925 | 294.00 | 1367 | 412.00 | 288 |
| 92.00 | 9722 | 192.00 | 10797 | 295.00 | 570 | 413.00 | 62 |
| 93.00 | 62768 | 193.00 | 11404 | 296.00 | 72376 | 414.00 | 116 |
| 94.00 | 3354 | 194.00 | 3574 | 297.00 | 11570 | 415.00 | 319 |
| 95.00 | 414 | 195.00 | 2250 | 298.00 | 710 | 416.00 | 229 |
| 96.00 | 2725 | 196.00 | 23304 | 299.00 | 504 | 417.00 | 241 |
| 98.00 | 45184 | 198.00 | 655488 | 301.00 | 960 | 418.00 | 258 |
| 99.00 | 34768 | 199.00 | 51056 | 302.00 | 1550 | 419.00 | 24 |
| 100.00 | 2441 | 200.00 | 3768 | 303.00 | 7568 | 420.00 | 98 |
| 101.00 | 18736 | 201.00 | 3895 | 304.00 | 1736 | 421.00 | 3247 |
| 102.00 | 1238 | 202.00 | 358 | 305.00 | 141 | 422.00 | 3938 |
| 103.00 | 6585 | 203.00 | 6471 | 306.00 | 199 | 423.00 | 19888 |
| 104.00 | 13969 | 204.00 | 33520 | 307.00 | 330 | 424.00 | 3955 |
| 105.00 | 10643 | 205.00 | 57048 | 308.00 | 1058 | 425.00 | 780 |
| 106.00 | 3843 | 206.00 | 210944 | 309.00 | 1110 | 426.00 | 619 |
| 107.00 | 148416 | 207.00 | 28512 | 310.00 | 1005 | 427.00 | 477 |
| 108.00 | 22168 | 208.00 | 7802 | 311.00 | 191 | 428.00 | 1065 |
| 109.00 | 832 | 209.00 | 2100 | 312.00 | 489 | 429.00 | 922 |
| 110.00 | 237440 | 210.00 | 4845 | 313.00 | 1078 | 430.00 | 1034 |

Date : 31-MAR-2011 20:14

Client ID: DFTPP0331

Instrument: nt12.i

Sample Info: DFTPP0331

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0.32

Data File: 03311101.d

Spectrum: Avg. Scans 482-484 (5.66), Background Scan 478

Location of Maximum: 198.00

Number of points: 394

| m/z | Y | m/z | Y | m/z | Y | m/z | Y |
|--------|--------|--------|--------|--------|-------|--------|--------|
| 111.00 | 35864 | 211.00 | 10197 | 314.00 | 2957 | 431.00 | 864 |
| 112.00 | 4264 | 212.00 | 1707 | 315.00 | 7566 | 432.00 | 776 |
| 113.00 | 1063 | 213.00 | 1107 | 316.00 | 4170 | 433.00 | 1166 |
| 114.00 | 505 | 214.00 | 424 | 317.00 | 532 | 434.00 | 520 |
| 115.00 | 444 | 215.00 | 3179 | 318.00 | 603 | 435.00 | 2007 |
| 116.00 | 7606 | 216.00 | 3532 | 319.00 | 289 | 436.00 | 1581 |
| 117.00 | 127472 | 217.00 | 64840 | 320.00 | 625 | 437.00 | 1718 |
| 118.00 | 8568 | 218.00 | 9360 | 321.00 | 2037 | 438.00 | 1285 |
| 119.00 | 992 | 219.00 | 1436 | 322.00 | 1011 | 439.00 | 2279 |
| 120.00 | 1205 | 221.00 | 31184 | 323.00 | 17768 | 440.00 | 404 |
| 121.00 | 534 | 222.00 | 9561 | 324.00 | 2786 | 441.00 | 61584 |
| 122.00 | 7775 | 223.00 | 14903 | 325.00 | 372 | 442.00 | 365696 |
| 123.00 | 12511 | 224.00 | 122816 | 326.00 | 344 | 443.00 | 72840 |
| 124.00 | 6616 | 225.00 | 27784 | 327.00 | 3904 | 444.00 | 6371 |
| 125.00 | 4767 | 226.00 | 3099 | 328.00 | 2526 | 445.00 | 677 |
| 126.00 | 473 | 227.00 | 60992 | 329.00 | 418 | 448.00 | 85 |
| 127.00 | 353792 | 228.00 | 8680 | 330.00 | 240 | 449.00 | 114 |
| 128.00 | 29984 | 229.00 | 9960 | 331.00 | 297 | 455.00 | 74 |
| 129.00 | 159488 | 230.00 | 1173 | 332.00 | 1662 | 456.00 | 69 |
| 130.00 | 13978 | 231.00 | 4680 | 333.00 | 1729 | 463.00 | 93 |
| 131.00 | 2437 | 232.00 | 831 | 334.00 | 13645 | 467.00 | 220 |
| 132.00 | 1247 | 233.00 | 265 | 335.00 | 2611 | 468.00 | 93 |
| 133.00 | 289 | 234.00 | 4005 | 336.00 | 404 | 474.00 | 72 |
| 134.00 | 4813 | 235.00 | 4166 | 339.00 | 357 | 481.00 | 103 |
| 135.00 | 12545 | 236.00 | 3099 | 340.00 | 429 | 490.00 | 116 |
| 136.00 | 5676 | 237.00 | 4095 | 341.00 | 2293 | 493.00 | 55 |
| 137.00 | 6839 | 238.00 | 280 | 342.00 | 839 | 494.00 | 62 |
| 138.00 | 686 | 239.00 | 2416 | 343.00 | 81 | | |
| 139.00 | 1502 | 240.00 | 1585 | 344.00 | 118 | | |

Date : 31-MAR-2011 20:14

Client ID: DFTPP0331

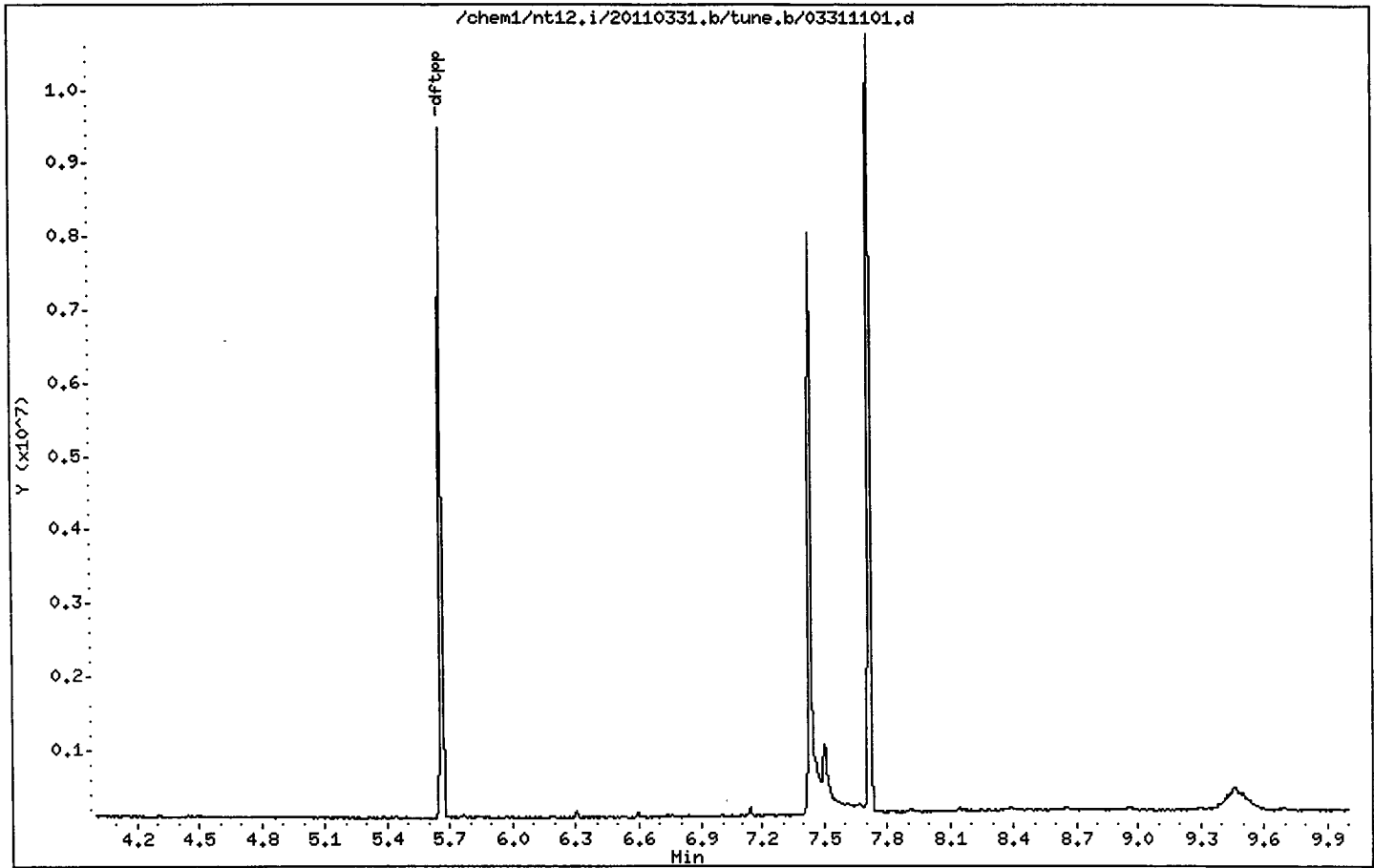
Instrument: nt12.i

Sample Info: DFTPP0331

Operator: JZ

Column phase: ZB-5msi

Column diameter: 0.32



Analytical Resources Inc.
ABN by sw846 8270C
DDT Breakdown Report

Data file: /chem1/nt12.i/20110331.b/ddt.b/03311101.d ARI ID: DDT0331
Method: /chem1/nt12.i/20110331.b/ddt.b/sw846ddt.m Misc: 11-
Analysis Date: 31-MAR-2011 13:17 Instrument: nt12.i

| COMPOUND | RT | AREA |
|-------------------|-------|--------|
| Pentachlorophenol | 5.774 | 52518 |
| Benzidine | 7.452 | 816154 |
| 4,4'-DDE | ---- | ---- |
| 4,4'-DDD | 7.527 | 26312 |
| 4,4'-DDT | 7.751 | 412197 |

$$\text{DDT Percent Breakdown} = \frac{(\text{DDE Area} + \text{DDD Area}) * 100}{(\text{DDE Area} + \text{DDD Area} + \text{DDT Area})}$$

$$\text{DDT Percent Breakdown} = \frac{(0 + 26312) * 100}{(0 + 26312 + 412197)}$$

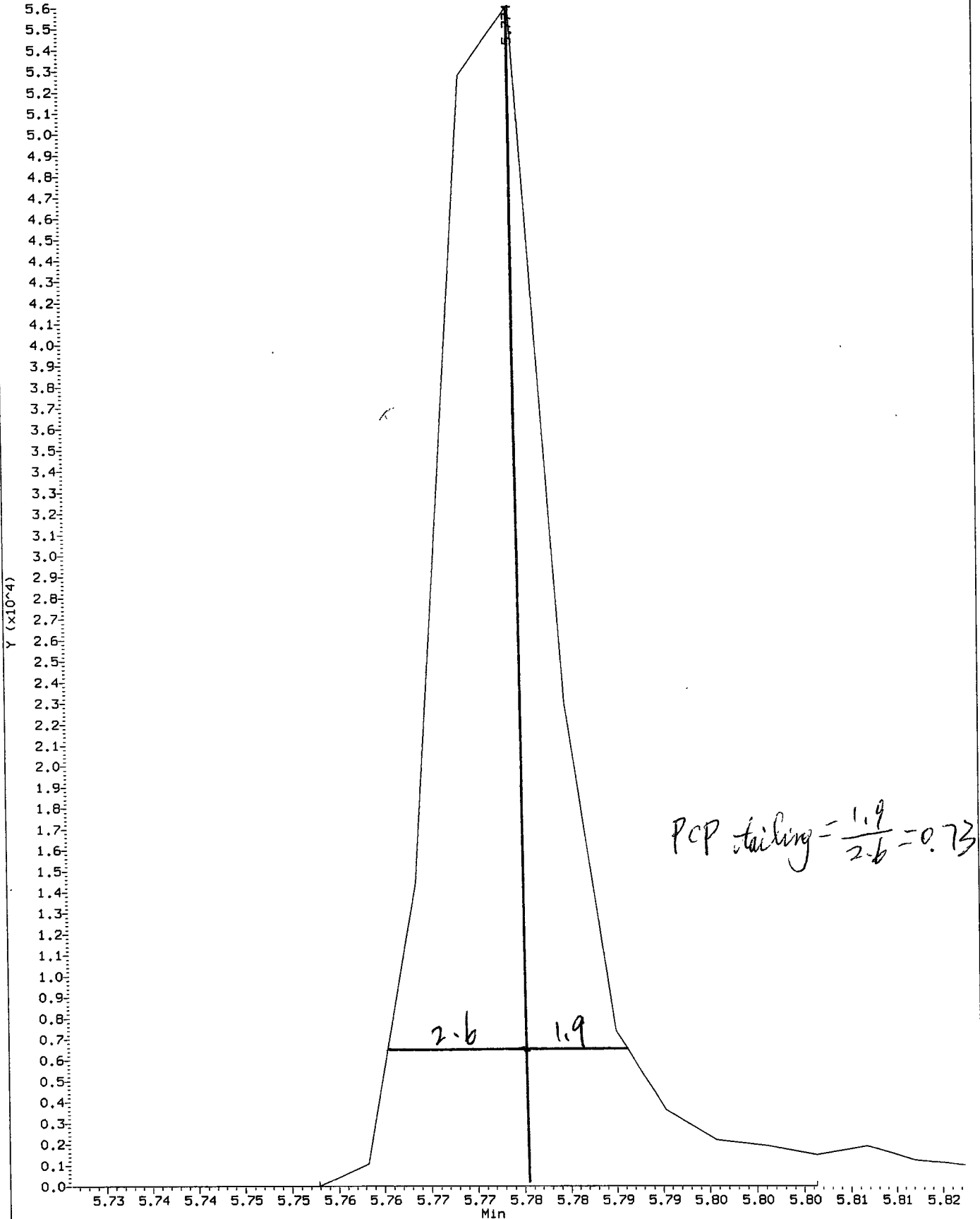
DDT Percent Breakdown = 6.0 %

ok *to* *04/01/11*

Data File: /chem1/nt12.1/20110331.b/ddt.b/03311101.d
Injection Date: 31-MAR-2011 13:17
Instrument: nt12.i
Client Sample ID: DDT0331

Compound: Pentachlorophenol
CAS Number: 87-86-5

Ion 266.00: Area: 52518 Height: 56104

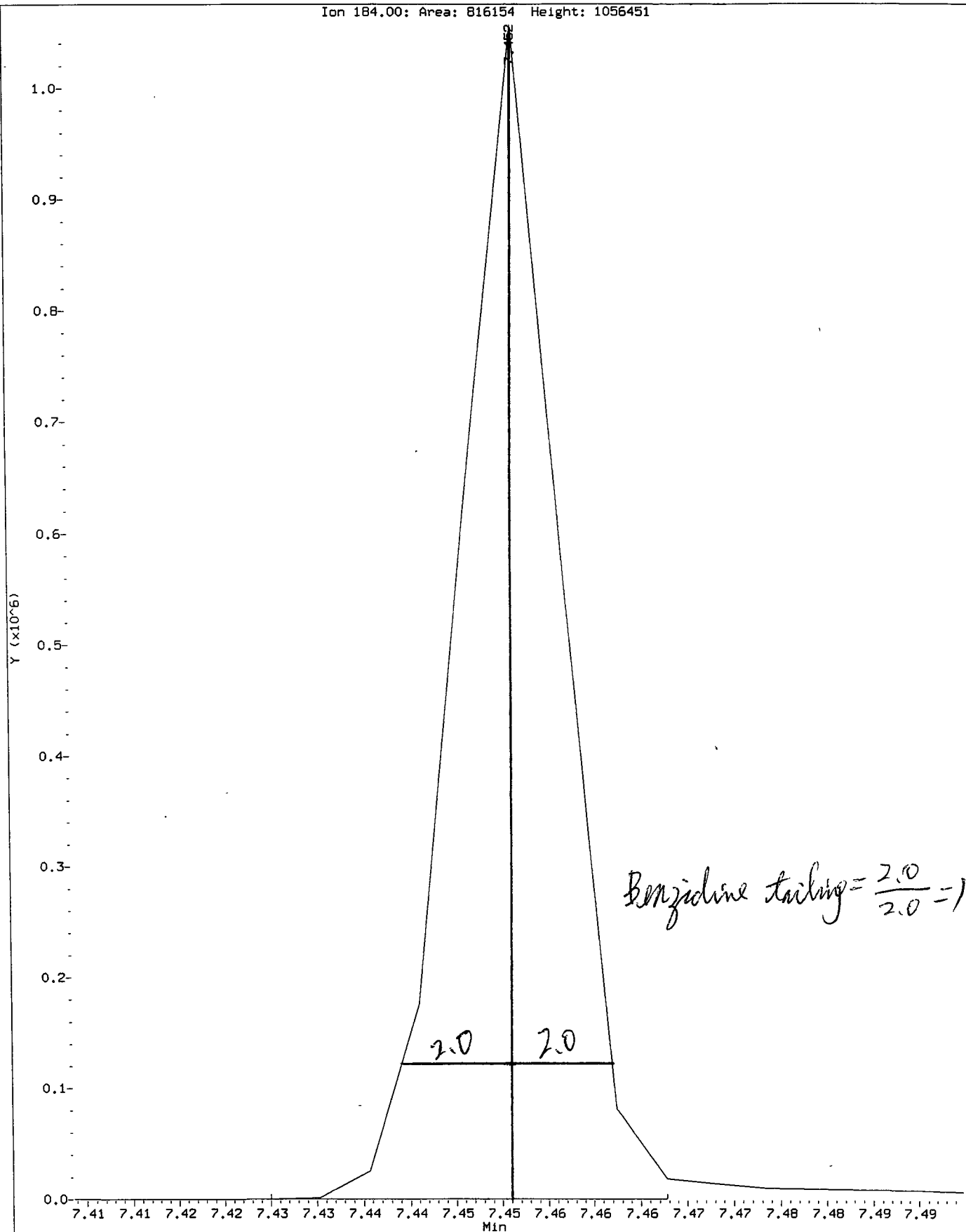


SN54: 00147

Data File: /chem1/nt12.1/20110331.b/ddt.b/03311101.d
Injection Date: 31-MAR-2011 13:17
Instrument: nt12.1
Client Sample ID: DDT0331

Compound: Benzidine
CAS Number:

Ion 184.00: Area: 816154 Height: 1056451



SN54:00148

Analytical Resources, Inc.

Semivolatle Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311103.d
 Lab Smp Id: IC010331 Client Smp ID: IC010331
 Inj Date : 31-MAR-2011 20:55
 Operator : JZ Inst ID: nt12.i
 Smp Info : IC010331,
 Misc Info : 11-
 Comment : 1ul Injection
 Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 3 Calibration Sample, Level: 1
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: pnax.sub
 Target Version: 3.50

Handwritten: 04/01/11

| Compounds | QUANT SIG | MASS | RT | EXP RT | REL RT | RESPONSE | AMOUNTS | |
|-------------------------------------|-----------|--------|--------|---------|--------|----------|-----------------|----------------|
| | | | | | | | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.917 | 4.917 | (1.000) | 382584 | 2.00000 | | |
| 28 Naphthalene | 128 | 4.945 | 4.945 | (1.006) | 19700 | 0.10000 | 0.1066 | |
| \$ 190 2-Methylnaphthalene-d10 | 152 | 5.655 | 5.655 | (1.150) | 13809 | 0.10000 | 0.1181 | |
| 32 2-Methylnaphthalene | 141 | 5.700 | 5.700 | (1.159) | 12073 | 0.10000 | 0.1176 | |
| 105 1-methylnaphthalene | 141 | 5.889 | 5.892 | (1.198) | 13974 | 0.10000 | 0.1279 | |
| 40 Acenaphthylene | 152 | 7.025 | 7.025 | (0.981) | 22074 | 0.10000 | 0.1123 | |
| * 42 Acenaphthene-d10 | 164 | 7.161 | 7.164 | (1.000) | 224662 | 2.00000 | | |
| 44 Acenaphthene | 153 | 7.211 | 7.208 | (1.007) | 14906 | 0.10000 | 0.1217 | |
| 46 Dibenzofuran | 168 | 7.350 | 7.353 | (1.026) | 19734 | 0.10000 | 0.1185 | |
| 49 Fluorene | 166 | 7.804 | 7.808 | (1.090) | 15973 | 0.10000 | 0.1165 | |
| * 59 Phenanthrene-d10 | 188 | 9.098 | 9.098 | (1.000) | 370050 | 2.00000 | | |
| 60 Phenanthrene | 178 | 9.130 | 9.133 | (1.003) | 23257 | 0.10000 | 0.1191 | |
| 61 Anthracene | 178 | 9.164 | 9.164 | (1.007) | 22926 | 0.10000 | 0.1169 | |
| 64 Fluoranthene | 202 | 10.821 | 10.827 | (1.189) | 24809 | 0.10000 | 0.1168 | |
| 65 Pyrene | 202 | 11.298 | 11.301 | (0.818) | 25492 | 0.10000 | 0.1171 | |
| 68 Benzo (a) anthracene | 228 | 13.686 | 13.693 | (0.991) | 22501 | 0.10000 | 0.1158 | |
| * 69 Chrysene-d12 | 240 | 13.806 | 13.806 | (1.000) | 404335 | 2.00000 | | |
| 71 Chrysene | 228 | 13.873 | 13.873 | (1.005) | 21911 | 0.10000 | 0.1153 | |
| 74 Benzo (b) fluoranthene | 252 | 16.302 | 16.312 | (0.934) | 21792 | 0.10000 | 0.1090 | |
| 75 Benzo (k) fluoranthene | 252 | 16.359 | 16.369 | (0.937) | 25137 | 0.10000 | 0.1198 | |
| 188 Benzo (j) fluoranthene | 252 | 16.425 | 16.444 | (0.941) | 24420 | 0.10000 | 0.1218 | |
| 76 Benzo (a) pyrene | 252 | 17.255 | 17.268 | (0.988) | 22219 | 0.10000 | 0.1210 | |
| * 77 Perylene-d12 | 264 | 17.464 | 17.473 | (1.000) | 354686 | 2.00000 | | |
| 78 Indeno (1,2,3-cd) pyrene | 276 | 19.802 | 19.833 | (1.134) | 24381 | 0.10000 | 0.1092 | |
| \$ 191 Dibenzo (a,h) anthracene-d14 | 292 | 19.729 | 19.751 | (1.130) | 19561 | 0.10000 | 0.1175 | |
| 79 Dibenzo (a,h) anthracene | 278 | 19.818 | 19.840 | (1.135) | 22181 | 0.10000 | 0.1215 | |
| 80 Benzo (g,h,i) perylene | 276 | 20.648 | 20.676 | (1.182) | 24243 | 0.10000 | 0.1224 | |

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-------------|-----------|--------|---------|---------|----------|--------------------|-------------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| ===== | ==== | == | ===== | ===== | ===== | ===== | ===== |
| 99 Perylene | 252 | 17.536 | 17.536 | (1.004) | 18544 | 0.10000 | 0.1194 |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 03311103.d
 Lab Smp Id: IC010331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 31-MAR-2011
 Calibration Time: 20:27
 Client Smp ID: IC010331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|--------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 382584 | -22.57 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 224662 | -19.79 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 370050 | -19.79 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 404335 | -19.64 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 354686 | -19.79 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | -0.06 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | -0.04 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.00 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.81 | 0.02 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.46 | -0.04 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/nt12.i/20110331.b/03311103.d

Date : 31-MAR-2011 20:55

Client ID: IC010331

Sample Info: IC010331,

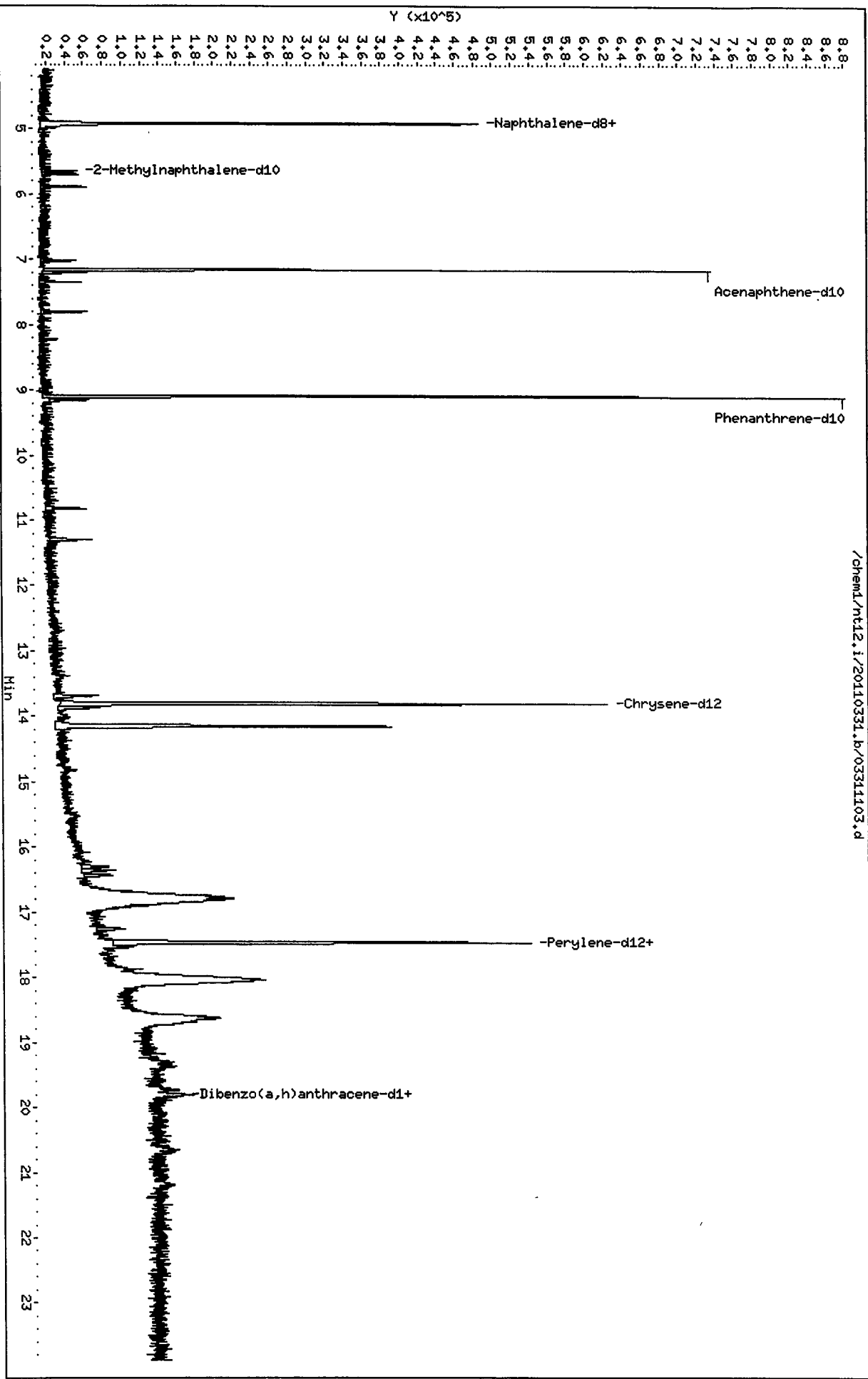
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110331.b/03311103.d



CO-ELUTION SUMMARY FOR FILE - 03311103.d

Lab ID: IC010331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00153

Analytical Resources, Inc.

Semivolatile Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311104.d
 Lab Smp Id: IC050331 Client Smp ID: IC050331
 Inj Date : 31-MAR-2011 21:23
 Operator : JZ Inst ID: nt12.i
 Smp Info : IC050331,
 Misc Info : 11-
 Comment : 1ul Injection
 Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 4 Calibration Sample, Level: 2
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: pmax.sub
 Target Version: 3.50

Handwritten signature and date: 04/01/11

| Compounds | QUANT | SIG | MASS | RT | EXP RT | REL RT | RESPONSE | AMOUNTS | |
|-------------------------------------|-------|-----|--------|--------|---------|--------|----------|-----------------|----------------|
| | | | | | | | | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | | 4.917 | 4.917 | (1.000) | 399629 | 2.00000 | | |
| 28 Naphthalene | 128 | | 4.949 | 4.945 | (1.006) | 104132 | 0.50000 | 0.5396 | |
| \$ 190 2-Methylnaphthalene-d10 | 152 | | 5.655 | 5.655 | (1.150) | 65979 | 0.50000 | 0.5400 | |
| 32 2-Methylnaphthalene | 141 | | 5.700 | 5.700 | (1.159) | 57513 | 0.50000 | 0.5363 | |
| 105 1-methylnaphthalene | 141 | | 5.892 | 5.892 | (1.198) | 60865 | 0.50000 | 0.5334 | |
| 40 Acenaphthylene | 152 | | 7.025 | 7.025 | (0.981) | 106123 | 0.50000 | 0.5270 | |
| * 42 Acenaphthene-d10 | 164 | | 7.161 | 7.164 | (1.000) | 230116 | 2.00000 | | |
| 44 Acenaphthene | 153 | | 7.208 | 7.208 | (1.007) | 63191 | 0.50000 | 0.5038 | |
| 46 Dibenzofuran | 168 | | 7.353 | 7.353 | (1.027) | 89653 | 0.50000 | 0.5257 | |
| 49 Fluorene | 166 | | 7.808 | 7.808 | (1.090) | 71498 | 0.50000 | 0.5092 | |
| * 59 Phenanthrene-d10 | 188 | | 9.098 | 9.098 | (1.000) | 380150 | 2.00000 | | |
| 60 Phenanthrene | 178 | | 9.130 | 9.133 | (1.003) | 104524 | 0.50000 | 0.5210 | |
| 61 Anthracene | 178 | | 9.164 | 9.164 | (1.007) | 104184 | 0.50000 | 0.5172 | |
| 64 Fluoranthene | 202 | | 10.827 | 10.827 | (1.190) | 115111 | 0.50000 | 0.5276 | |
| 65 Pyrene | 202 | | 11.294 | 11.301 | (0.818) | 116889 | 0.50000 | 0.5128 | |
| 68 Benzo (a) anthracene | 228 | | 13.683 | 13.693 | (0.991) | 100173 | 0.50000 | 0.4922 (M) | |
| * 69 Chrysene-d12 | 240 | | 13.803 | 13.806 | (1.000) | 423330 | 2.00000 | | |
| 71 Chrysene | 228 | | 13.866 | 13.873 | (1.005) | 101086 | 0.50000 | 0.5082 | |
| 74 Benzo (b) fluoranthene | 252 | | 16.299 | 16.312 | (0.933) | 105384 | 0.50000 | 0.5054 | |
| 75 Benzo (k) fluoranthene | 252 | | 16.359 | 16.369 | (0.937) | 111647 | 0.50000 | 0.5105 | |
| 188 Benzo (j) fluoranthene | 252 | | 16.429 | 16.444 | (0.941) | 105384 | 0.50000 | 0.5043 | |
| 76 Benzo (a) pyrene | 252 | | 17.255 | 17.268 | (0.988) | 93384 | 0.50000 | 0.4878 (M) | |
| * 77 Perylene-d12 | 264 | | 17.464 | 17.473 | (1.000) | 369841 | 2.00000 | | |
| 78 Indeno (1,2,3-cd) pyrene | 276 | | 19.802 | 19.833 | (1.134) | 121347 | 0.50000 | 0.5213 | |
| \$ 191 Dibenzo (a,h) anthracene-d14 | 292 | | 19.733 | 19.751 | (1.130) | 87297 | 0.50000 | 0.5030 | |
| 79 Dibenzo (a,h) anthracene | 278 | | 19.821 | 19.840 | (1.135) | 86728 | 0.50000 | 0.4556 (M) | |
| 80 Benzo (g,h,i) perylene | 276 | | 20.663 | 20.676 | (1.183) | 104467 | 0.50000 | 0.5060 | |

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-------------|-----------|--------|---------|---------|----------|--------------------|-------------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| ----- | ---- | == | ===== | ===== | ===== | ===== | ===== |
| 99 Perylene | 252 | 17.536 | 17.536 | (1.004) | 81801 | 0.50000 | 0.5053 |

QC Flag Legend

M - Compound response manually integrated.

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 03311104.d
 Lab Smp Id: IC050331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 31-MAR-2011
 Calibration Time: 20:27
 Client Smp ID: IC050331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|--------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 399629 | -19.12 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 230116 | -17.85 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 380150 | -17.60 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 423330 | -15.87 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 369841 | -16.37 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | -0.06 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | -0.04 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.00 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.80 | 0.00 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.46 | -0.04 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/rt12.i/20110331.b/03311104.d

Date: 31-MAR-2011 21:23

Client ID: IC050331

Sample Info: IC050331,

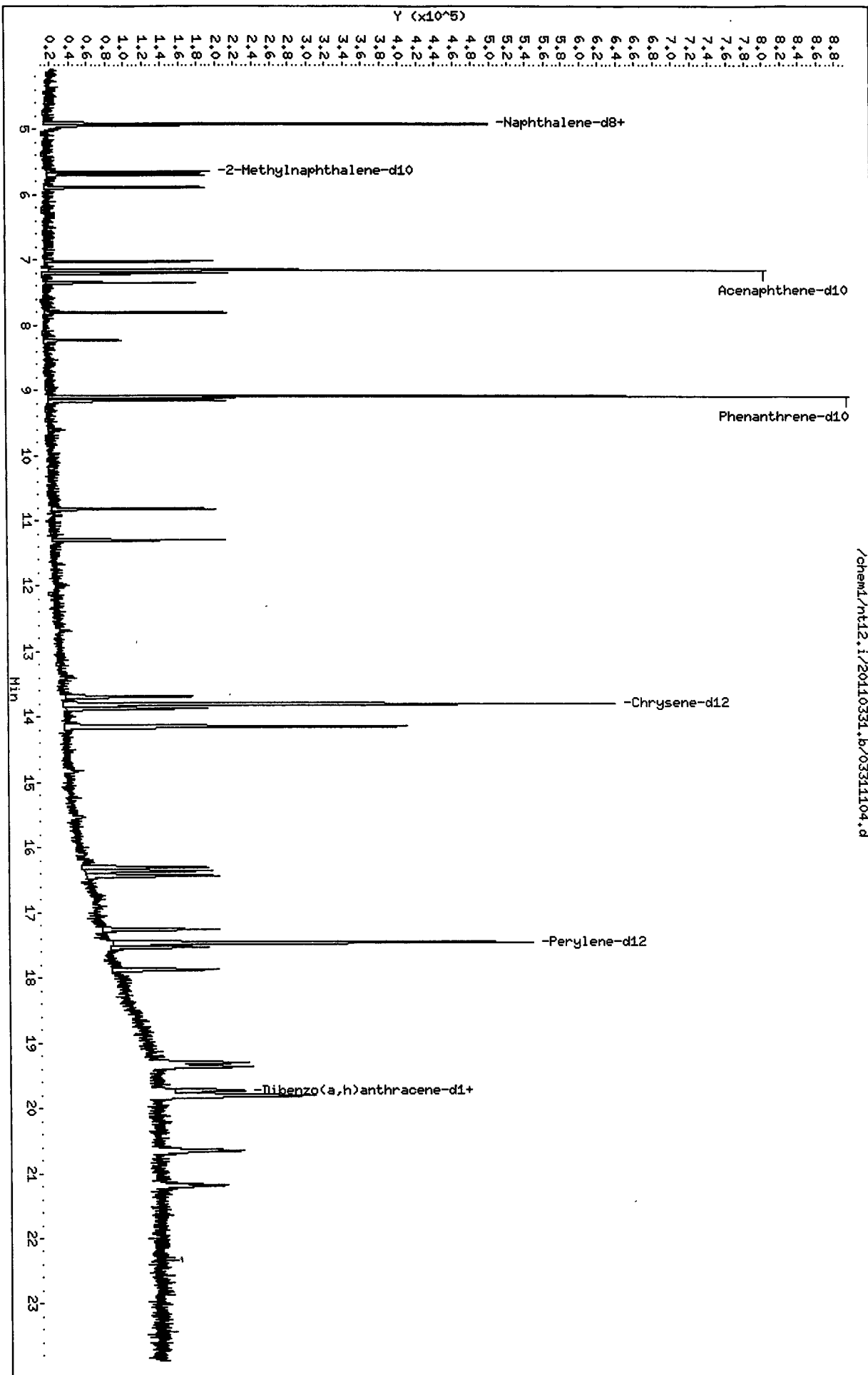
Column phase: ZB35

Instrument: rt12.i

Operator: JZ

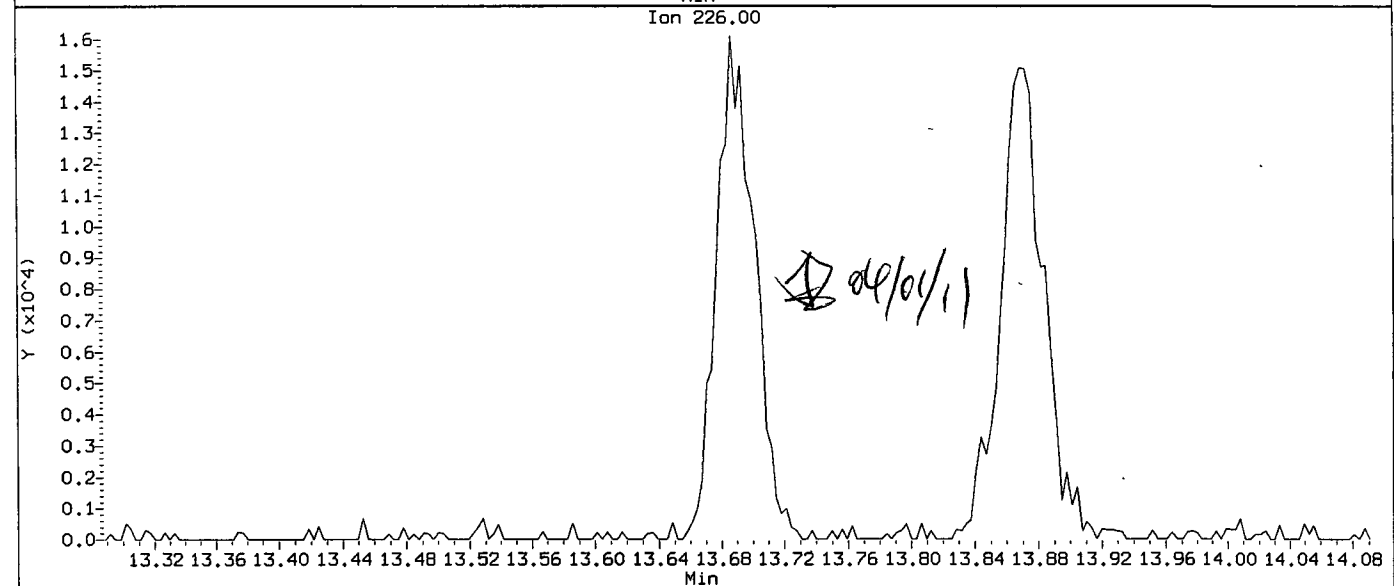
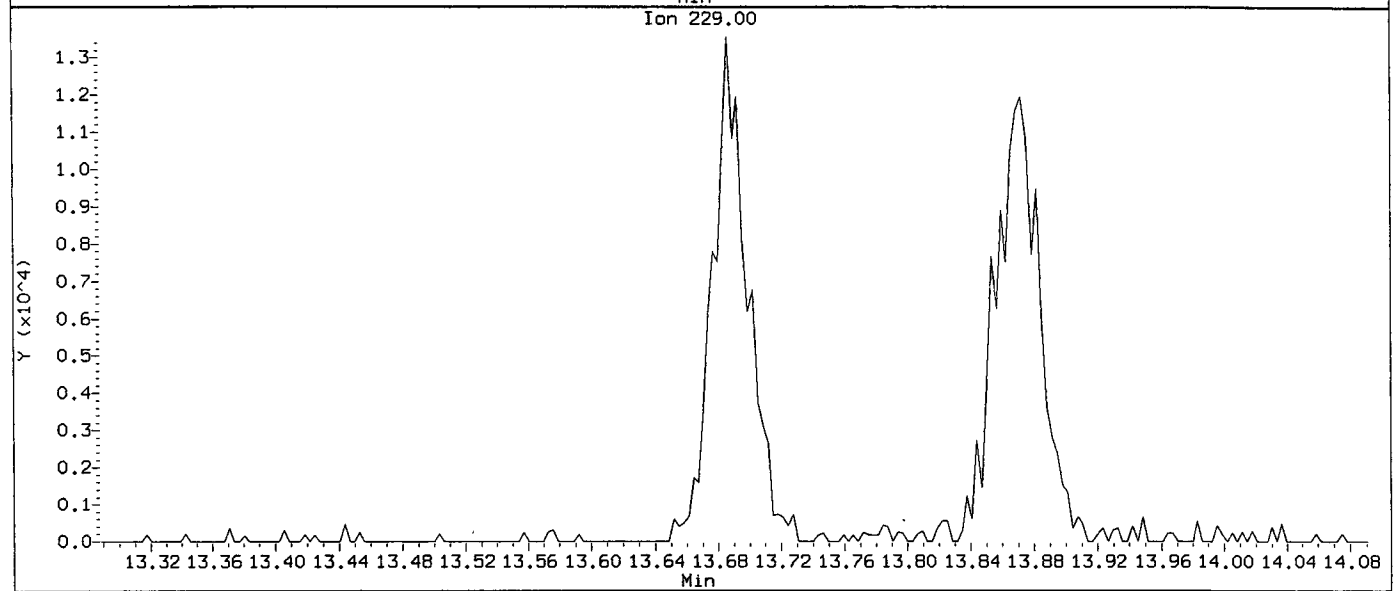
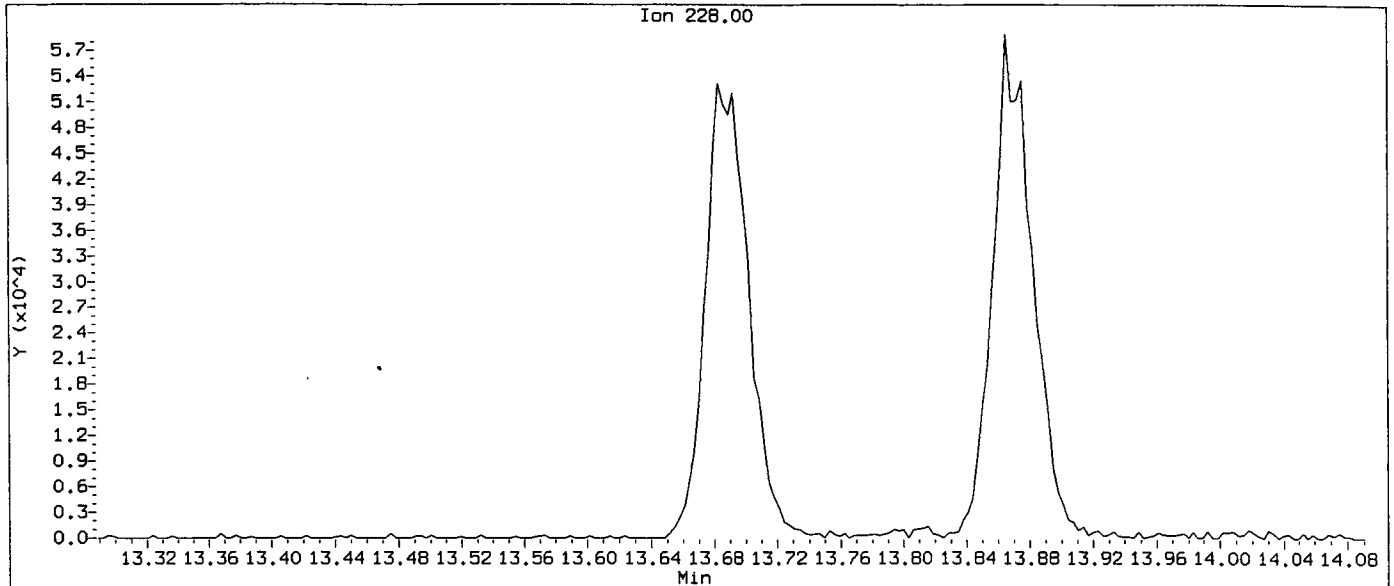
Column diameter: 0.32

/chem1/rt12.i/20110331.b/03311104.d



Data File: /chem1/nt12.1/20110331.b/03311104.d
Injection Date: 31-MAR-2011 21:23
Instrument: nt12.1
Client Sample ID: IC050331

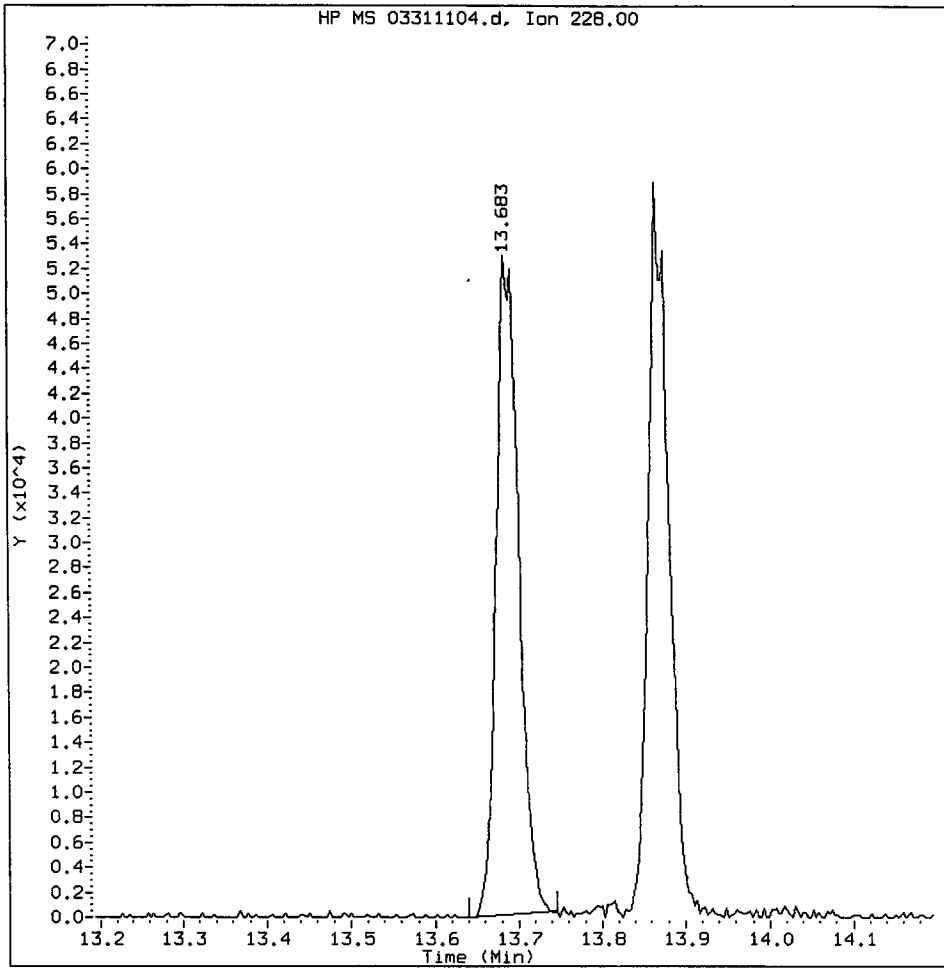
Compound: Benzo(a)anthracene
CAS Number: 56-55-3



SN54:00158

IC050331, /chem1/nt12.i/20110331.b/03311104.d

Benzo(a)anthracene Amount: 0.49 Area: 100173



MANUAL INTEGRATION for Benzo(a)anthracene

1. Baseline correction
2. Poor chromatography
3. Peak not found
4. Totals calculation

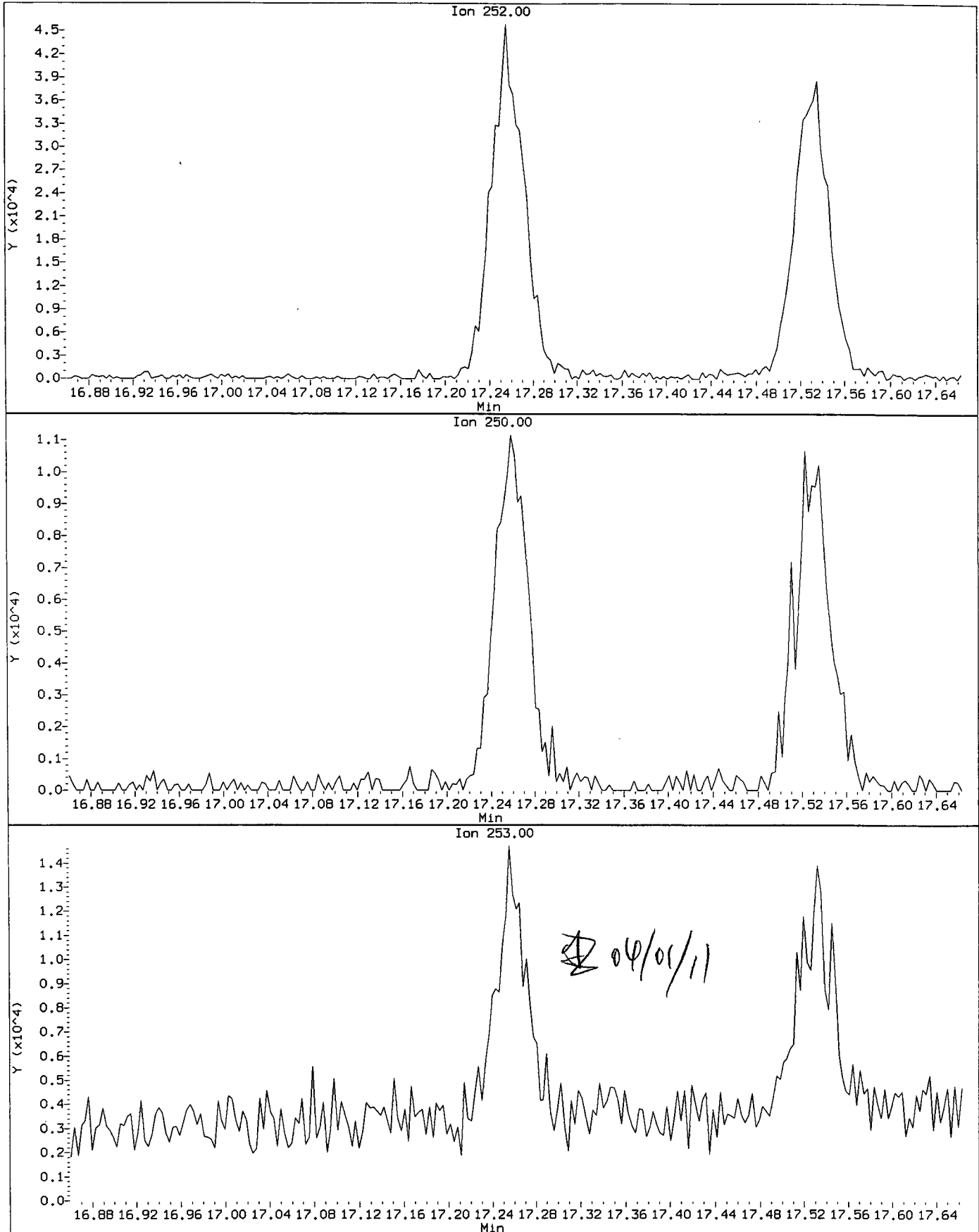
5. Other _____

Analyst: AD

Date: 04/01/11

Data File: /chem1/nt12.1/20110331.b/03311104.d
Injection Date: 31-MAR-2011 21:23
Instrument: nt12.1
Client Sample ID: IC050331

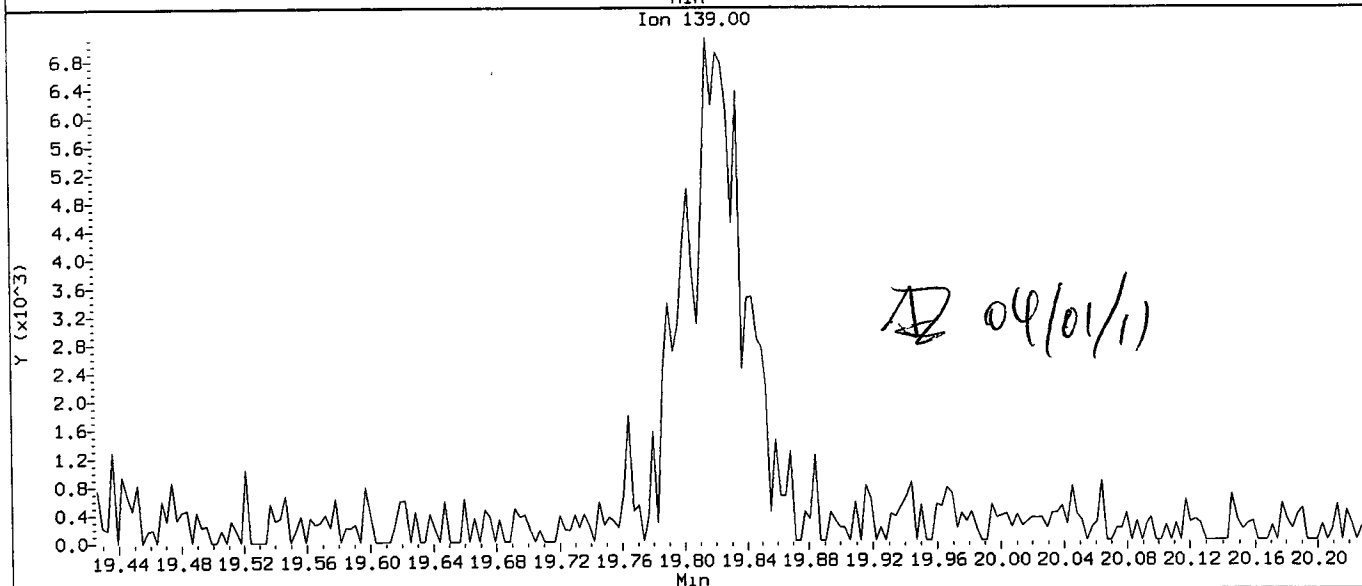
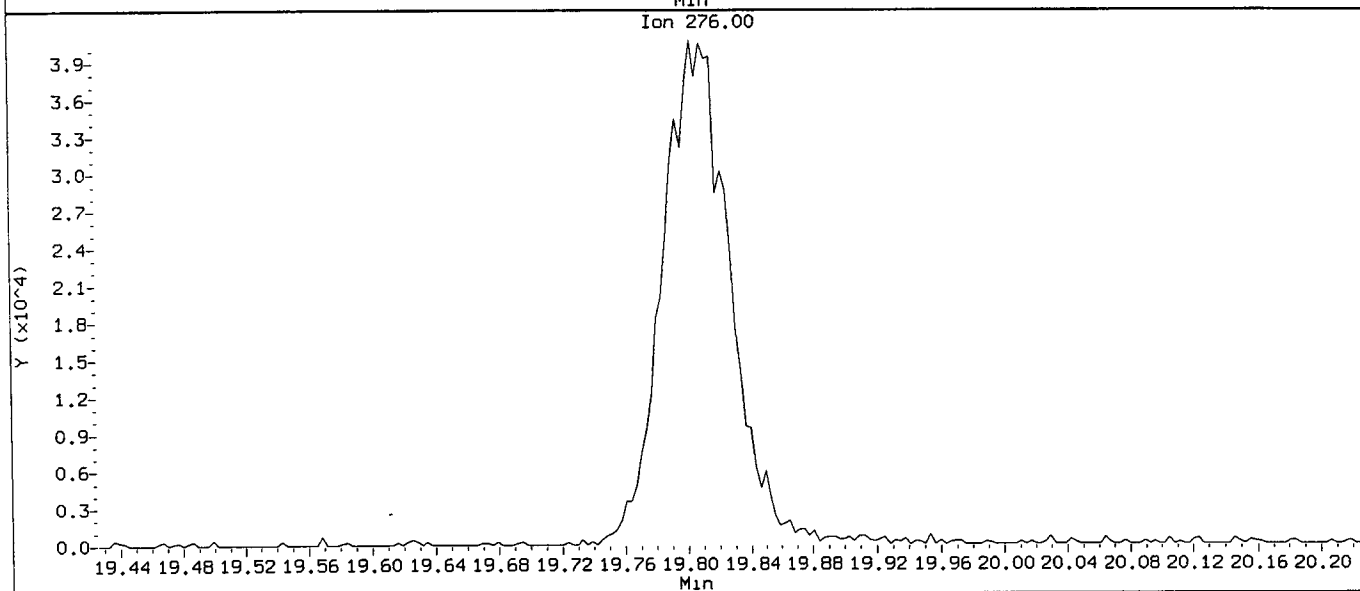
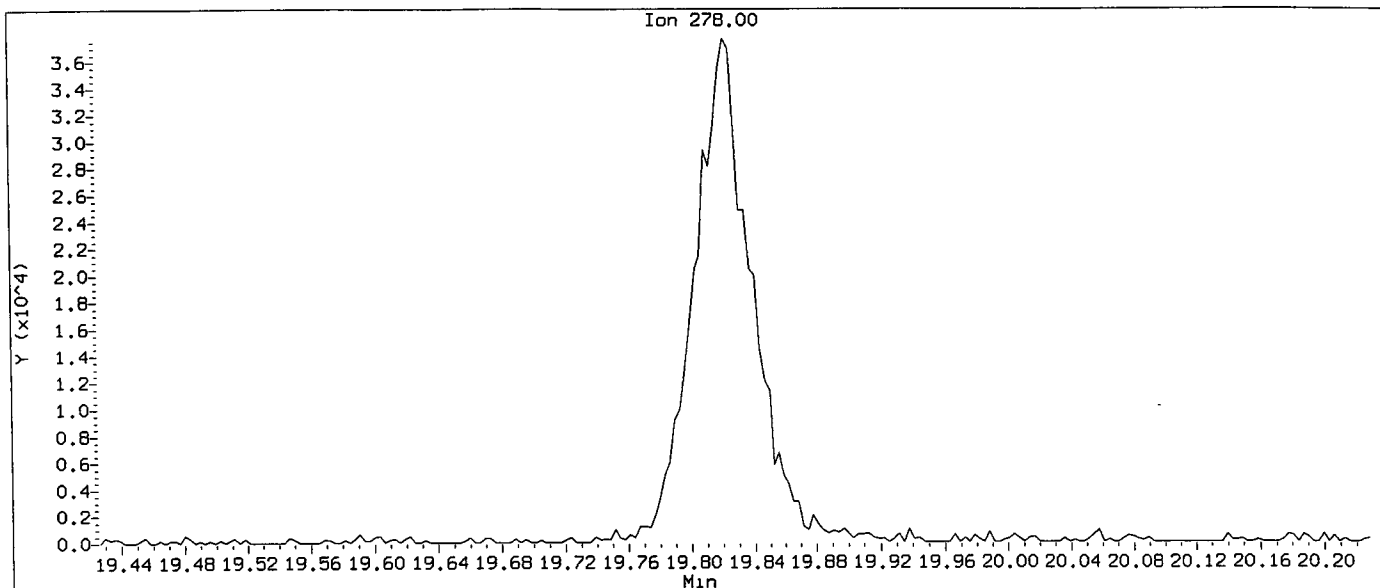
Compound: Benzo(a)pyrene
CAS Number: 50-32-8



SN54 : 00160

Data File: /chem1/nt12.1/20110331.b/03311104.d
Injection Date: 31-MAR-2011 21:23
Instrument: nt12.1
Client Sample ID: IC050331

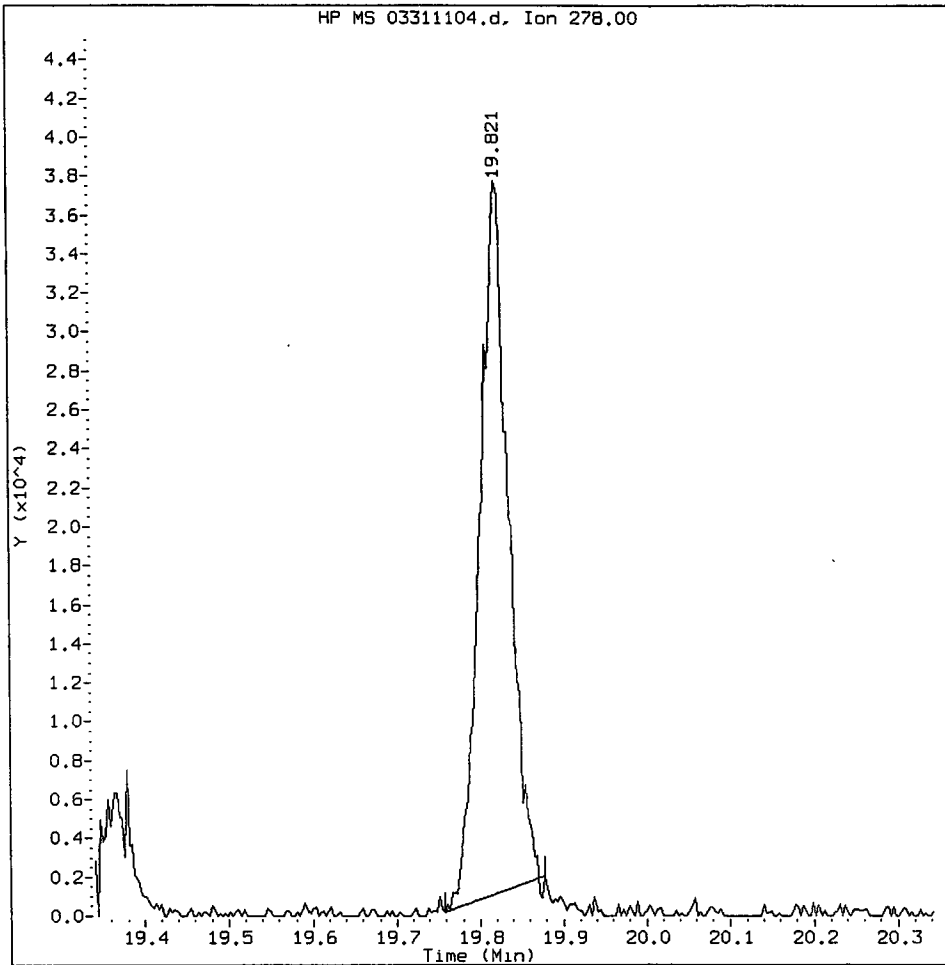
Compound: Dibenzo(a,h)anthracene
CAS Number: 53-70-3



SN54:00162

IC050331, /chem1/nt12.i/20110331.b/03311104.d

Dibenzo(a,h)anthracene Amount: 0.46 Area: 86728



MANUAL INTEGRATION for Dibenzo(a,h)anthracene

1. Baseline correction
2. Poor chromatography
3. Peak not found
4. Totals calculation

5. Other _____

Analyst: AD

Date: 06/01/11

SN54: 00163

CO-ELUTION SUMMARY FOR FILE - 03311104.d

Lab ID: IC050331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54:00164

Analytical Resources, Inc.

Semivolatiles Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311105.d
Lab Smp Id: IC10331 Client Smp ID: IC10331
Inj Date : 31-MAR-2011 21:51
Operator : JZ Inst ID: nt12.i
Smp Info : IC10331,
Misc Info : 11-
Comment : 1ul Injection
Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
Als bottle: 5 Calibration Sample, Level: 3
Dil Factor: 1.00000
Integrator: HP RTE Compound Sublist: pnax.sub
Target Version: 3.50

Handwritten signature and date: 04/01/11
AMOUNTS

| Compounds | QUANT | SIG | | | | AMOUNTS | |
|-----------------------------------|-------|--------|--------|---------|----------|-----------------|----------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.920 | 4.917 | (1.000) | 447156 | 2.00000 | |
| 28 Naphthalene | 128 | 4.945 | 4.945 | (1.005) | 215855 | 1.00000 | 0.9997 |
| \$ 190 2-Methylnaphthalene-d10 | 152 | 5.655 | 5.655 | (1.149) | 138509 | 1.00000 | 1.013 |
| 32 2-Methylnaphthalene | 141 | 5.696 | 5.700 | (1.158) | 121414 | 1.00000 | 1.012 |
| 105 1-methylnaphthalene | 141 | 5.892 | 5.892 | (1.198) | 126734 | 1.00000 | 0.9925 |
| 40 Acenaphthylene | 152 | 7.025 | 7.025 | (0.981) | 226770 | 1.00000 | 1.021 |
| * 42 Acenaphthene-d10 | 164 | 7.164 | 7.164 | (1.000) | 253822 | 2.00000 | |
| 44 Acenaphthene | 153 | 7.208 | 7.208 | (1.006) | 139508 | 1.00000 | 1.008 |
| 46 Dibenzofuran | 168 | 7.353 | 7.353 | (1.026) | 188297 | 1.00000 | 1.001 |
| 49 Fluorene | 166 | 7.807 | 7.808 | (1.090) | 155567 | 1.00000 | 1.004 |
| * 59 Phenanthrene-d10 | 188 | 9.098 | 9.098 | (1.000) | 425121 | 2.00000 | |
| 60 Phenanthrene | 178 | 9.133 | 9.133 | (1.004) | 225236 | 1.00000 | 1.004 |
| 61 Anthracene | 178 | 9.164 | 9.164 | (1.007) | 225493 | 1.00000 | 1.001 |
| 64 Fluoranthene | 202 | 10.824 | 10.827 | (1.190) | 242321 | 1.00000 | 0.9932 |
| 65 Pyrene | 202 | 11.298 | 11.301 | (0.818) | 248242 | 1.00000 | 0.9992 |
| 68 Benzo(a)anthracene | 228 | 13.689 | 13.693 | (0.992) | 217524 | 1.00000 | 0.9806 |
| * 69 Chrysene-d12 | 240 | 13.803 | 13.806 | (1.000) | 461408 | 2.00000 | |
| 71 Chrysene | 228 | 13.872 | 13.873 | (1.005) | 215600 | 1.00000 | 0.9945 |
| 74 Benzo(b)fluoranthene | 252 | 16.302 | 16.312 | (0.933) | 240884 | 1.00000 | 1.017 |
| 75 Benzo(k)fluoranthene | 252 | 16.359 | 16.369 | (0.937) | 241536 | 1.00000 | 0.9721 |
| 188 Benzo(j)fluoranthene | 252 | 16.432 | 16.444 | (0.941) | 228216 | 1.00000 | 0.9612 |
| 76 Benzo(a)pyrene | 252 | 17.252 | 17.268 | (0.988) | 209214 | 1.00000 | 0.9620 |
| * 77 Perylene-d12 | 264 | 17.467 | 17.473 | (1.000) | 420159 | 2.00000 | |
| 78 Indeno(1,2,3-cd)pyrene | 276 | 19.799 | 19.833 | (1.133) | 256435 | 1.00000 | 0.9697 |
| \$ 191 Dibenzo(a,h)anthracene-d14 | 292 | 19.729 | 19.751 | (1.130) | 186714 | 1.00000 | 0.9470 |
| 79 Dibenzo(a,h)anthracene | 278 | 19.821 | 19.840 | (1.135) | 208586 | 1.00000 | 0.9645 |
| 80 Benzo(g,h,i)perylene | 276 | 20.654 | 20.676 | (1.182) | 219349 | 1.00000 | 0.9352 |

| Compounds | QUANT SIG | | | | | | AMOUNTS | |
|-------------|-----------|--------|--------|---------|----------|--------------------|-------------------|--|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) | |
| ----- | ---- | == | ===== | ===== | ===== | ===== | ===== | |
| 99 Perylene | 252 | 17.530 | 17.536 | (1.004) | 179605 | 1.00000 | 0.9765 | |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 03311105.d
 Lab Smp Id: IC10331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 31-MAR-2011
 Calibration Time: 20:27
 Client Smp ID: IC10331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 447156 | -9.50 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 253822 | -9.38 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 425121 | -7.85 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 461408 | -8.30 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 420159 | -4.99 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | 0.00 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | 0.00 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.00 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.80 | 0.00 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.47 | -0.02 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/nt12.i/20110331.b/03311105.d

Date: 31-MAR-2011 21:51

Client ID: IC10331

Sample Info: IC10331,

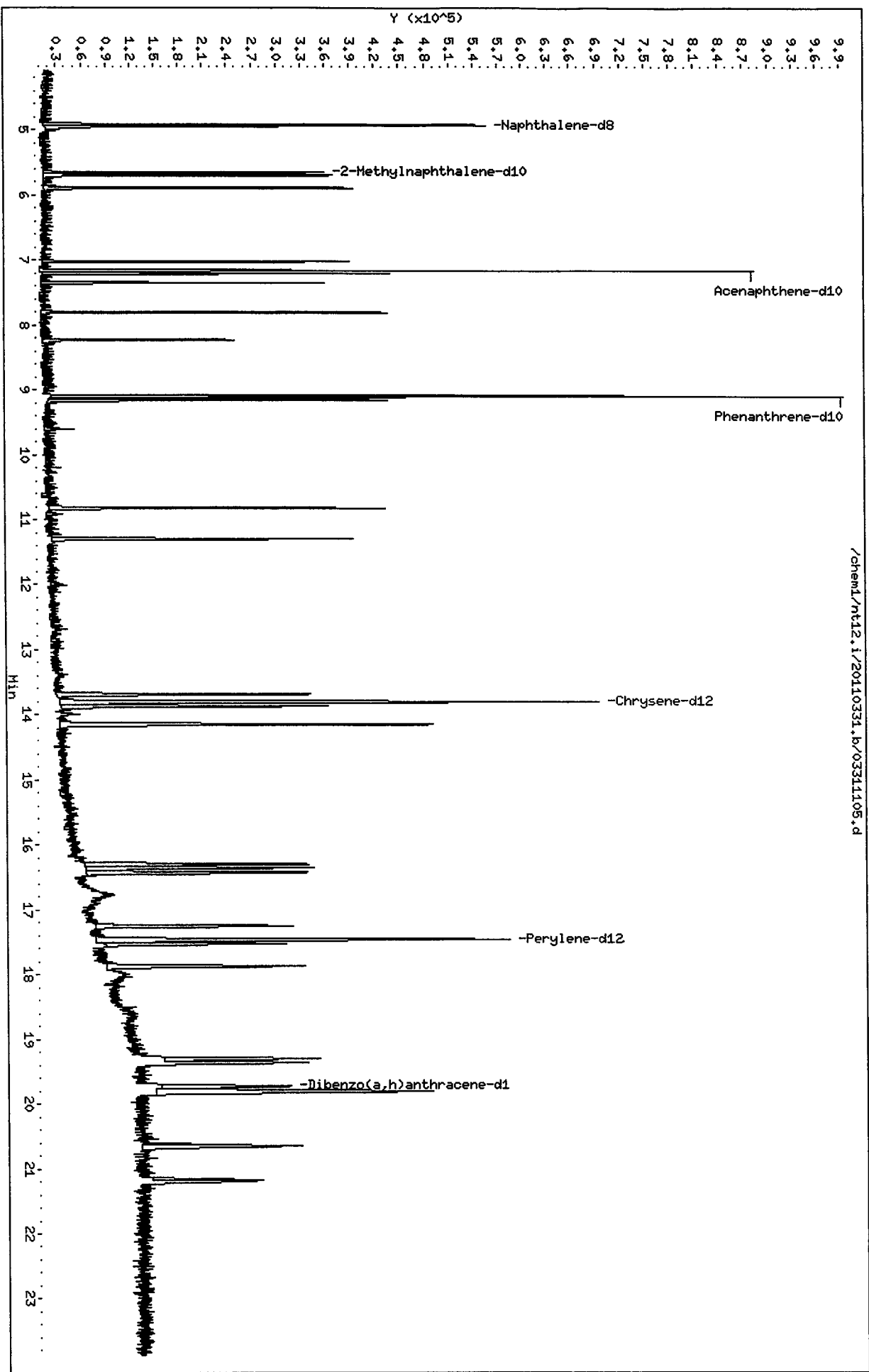
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110331.b/03311105.d



CO-ELUTION SUMMARY FOR FILE - 03311105.d

Lab ID: IC10331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00169

Analytical Resources, Inc.

Semivolatile Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311102.d
 Lab Smp Id: IC0250331 Client Smp ID: IC0250331
 Inj Date : 31-MAR-2011 20:27
 Operator : JZ Inst ID: nt12.i
 Smp Info : IC0250331
 Misc Info : 11-
 Comment : lul Injection
 Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 2 Calibration Sample, Level: 4
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: pnax.sub
 Target Version: 3.50

B 04/01/11

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-----------------------------------|-----------|--------|---------|---------|----------|-----------------|----------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.920 | 4.917 | (1.000) | 494112 | 2.00000 | . |
| 28 Naphthalene | 128 | 4.945 | 4.945 | (1.005) | 609116 | 2.50000 | 2.553 |
| \$ 190 2-Methylnaphthalene-d10 | 152 | 5.655 | 5.655 | (1.149) | 364043 | 2.50000 | 2.410 |
| 32 2-Methylnaphthalene | 141 | 5.700 | 5.700 | (1.158) | 323517 | 2.50000 | 2.440 |
| 105 1-methylnaphthalene | 141 | 5.892 | 5.892 | (1.198) | 335266 | 2.50000 | 2.376 |
| 40 Acenaphthylene | 152 | 7.025 | 7.025 | (0.981) | 592188 | 2.50000 | 2.416 |
| * 42 Acenaphthene-d10 | 164 | 7.164 | 7.164 | (1.000) | 280105 | 2.00000 | |
| 44 Acenaphthene | 153 | 7.208 | 7.208 | (1.006) | 360113 | 2.50000 | 2.359 |
| 46 Dibenzofuran | 168 | 7.353 | 7.353 | (1.026) | 494475 | 2.50000 | 2.382 |
| 49 Fluorene | 166 | 7.808 | 7.808 | (1.090) | 409605 | 2.50000 | 2.397 |
| * 59 Phenanthrene-d10 | 188 | 9.098 | 9.098 | (1.000) | 461353 | 2.00000 | |
| 60 Phenanthrene | 178 | 9.133 | 9.133 | (1.004) | 577350 | 2.50000 | 2.371 |
| 61 Anthracene | 178 | 9.164 | 9.164 | (1.007) | 611931 | 2.50000 | 2.503 |
| 64 Fluoranthene | 202 | 10.827 | 10.827 | (1.190) | 638608 | 2.50000 | 2.412 |
| 65 Pyrene | 202 | 11.301 | 11.301 | (0.819) | 648018 | 2.50000 | 2.392 |
| 68 Benzo(a)anthracene | 228 | 13.690 | 13.693 | (0.992) | 588425 | 2.50000 | 2.433 |
| * 69 Chrysene-d12 | 240 | 13.803 | 13.806 | (1.000) | 503160 | 2.00000 | |
| 71 Chrysene | 228 | 13.873 | 13.873 | (1.005) | 567061 | 2.50000 | 2.399 |
| 74 Benzo(b)fluoranthene | 252 | 16.306 | 16.312 | (0.933) | 603897 | 2.50000 | 2.422 |
| 75 Benzo(k)fluoranthene | 252 | 16.362 | 16.369 | (0.937) | 626194 | 2.50000 | 2.395 |
| 188 Benzo(j)fluoranthene | 252 | 16.435 | 16.444 | (0.941) | 598489 | 2.50000 | 2.395 |
| 76 Benzo(a)pyrene | 252 | 17.262 | 17.268 | (0.988) | 549027 | 2.50000 | 2.399 |
| * 77 Perylene-d12 | 264 | 17.470 | 17.473 | (1.000) | 442215 | 2.00000 | |
| 78 Indeno(1,2,3-cd)pyrene | 276 | 19.818 | 19.833 | (1.134) | 666802 | 2.50000 | 2.396 |
| \$ 191 Dibenzo(a,h)anthracene-d14 | 292 | 19.745 | 19.751 | (1.130) | 498072 | 2.50000 | 2.400 |
| 79 Dibenzo(a,h)anthracene | 278 | 19.827 | 19.840 | (1.135) | 555306 | 2.50000 | 2.440 |
| 80 Benzo(g,h,i)perylene | 276 | 20.670 | 20.676 | (1.183) | 582882 | 2.50000 | 2.361 |

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-------------|-----------|--------|---------|---------|----------|--------------------|-------------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| ===== | ==== | == | ===== | ===== | ===== | ===== | ===== |
| 99 Perylene | 252 | 17.536 | 17.536 | (1.004) | 462859 | 2.50000 | 2.391 |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 03311102.d
 Lab Smp Id: IC0250331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 31-MAR-2011
 Calibration Time: 20:27
 Client Smp ID: IC0250331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 494112 | 0.00 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 280105 | 0.00 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 461353 | 0.00 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 503160 | 0.00 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 442215 | 0.00 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | 0.00 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | 0.00 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.00 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.80 | 0.00 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.47 | 0.00 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/nt12.i/20110331.b/03311102.d

Date: 31-MAR-2011 20:27

Client ID: IC0250331

Sample Info: IC0250331

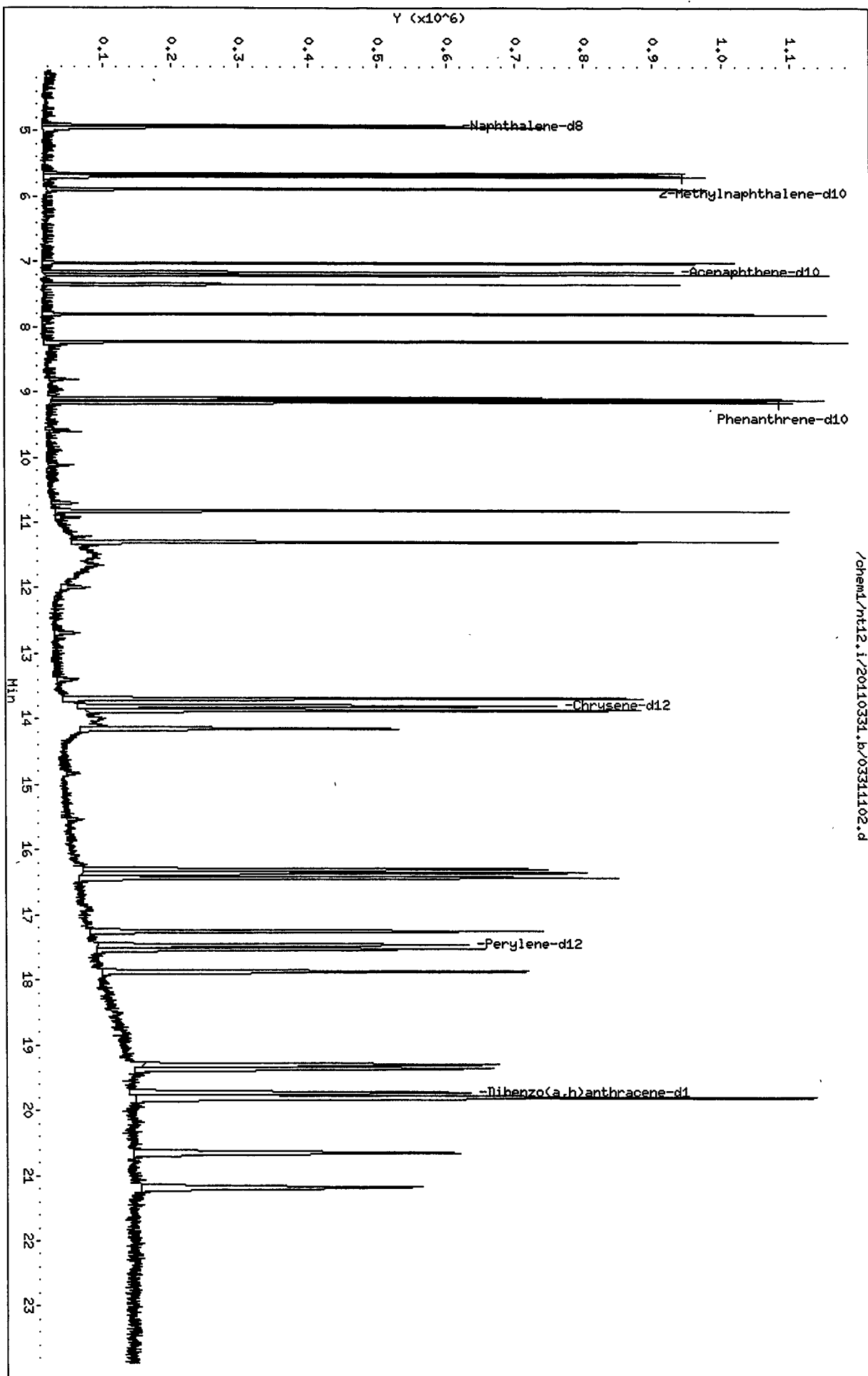
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110331.b/03311102.d



CO-ELUTION SUMMARY FOR FILE - 03311102.d

Lab ID: IC0250331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-201

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00174

Analytical Resources, Inc.

Semivolatile Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311106.d
 Lab Smp Id: IC50331 Client Smp ID: IC50331
 Inj Date : 31-MAR-2011 22:19
 Operator : JZ Inst ID: nt12.i
 Smp Info : IC50331,
 Misc Info : 11-
 Comment : lul Injection
 Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 6 Calibration Sample, Level: 5
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: pnax.sub
 Target Version: 3.50

J 04/01/11

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-----------------------------------|-----------|--------|---------|---------|----------|-----------------|----------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.920 | 4.917 | (1.000) | 424290 | 2.00000 | |
| 28 Naphthalene | 128 | 4.948 | 4.945 | (1.006) | 995611 | 5.00000 | 4.860 |
| \$ 190 2-Methylnaphthalene-d10 | 152 | 5.655 | 5.655 | (1.149) | 610176 | 5.00000 | 4.704 |
| 32 2-Methylnaphthalene | 141 | 5.699 | 5.700 | (1.158) | 536889 | 5.00000 | 4.716 |
| 105 1-methylnaphthalene | 141 | 5.892 | 5.892 | (1.198) | 554359 | 5.00000 | 4.576 |
| 40 Acenaphthylene | 152 | 7.022 | 7.025 | (0.980) | 989235 | 5.00000 | 4.850 |
| * 42 Acenaphthene-d10 | 164 | 7.164 | 7.164 | (1.000) | 233090 | 2.00000 | |
| 44 Acenaphthene | 153 | 7.208 | 7.208 | (1.006) | 606119 | 5.00000 | 4.771 |
| 46 Dibenzofuran | 168 | 7.353 | 7.353 | (1.026) | 823076 | 5.00000 | 4.765 |
| 49 Fluorene | 166 | 7.807 | 7.808 | (1.090) | 688644 | 5.00000 | 4.842 |
| * 59 Phenanthrene-d10 | 188 | 9.101 | 9.098 | (1.000) | 394000 | 2.00000 | |
| 60 Phenanthrene | 178 | 9.133 | 9.133 | (1.003) | 970112 | 5.00000 | 4.665 |
| 61 Anthracene | 178 | 9.164 | 9.164 | (1.007) | 1000694 | 5.00000 | 4.793 |
| 64 Fluoranthene | 202 | 10.824 | 10.827 | (1.189) | 1074309 | 5.00000 | 4.751 |
| 65 Pyrene | 202 | 11.297 | 11.301 | (0.818) | 1100784 | 5.00000 | 4.833 |
| 68 Benzo(a)anthracene | 228 | 13.693 | 13.693 | (0.992) | 1011134 | 5.00000 | 4.972 |
| * 69 Chrysene-d12 | 240 | 13.806 | 13.806 | (1.000) | 423013 | 2.00000 | |
| 71 Chrysene | 228 | 13.876 | 13.873 | (1.005) | 971908 | 5.00000 | 4.890 |
| 74 Benzo(b)fluoranthene | 252 | 16.305 | 16.312 | (0.933) | 1067461 | 5.00000 | 4.906 |
| 75 Benzo(k)fluoranthene | 252 | 16.359 | 16.369 | (0.936) | 1075252 | 5.00000 | 4.711 |
| 188 Benzo(j)fluoranthene | 252 | 16.435 | 16.444 | (0.941) | 1050965 | 5.00000 | 4.819 |
| 76 Benzo(a)pyrene | 252 | 17.262 | 17.268 | (0.988) | 981267 | 5.00000 | 4.912 |
| * 77 Perylene-d12 | 264 | 17.470 | 17.473 | (1.000) | 385980 | 2.00000 | |
| 78 Indeno(1,2,3-cd)pyrene | 276 | 19.818 | 19.833 | (1.134) | 1209015 | 5.00000 | 4.977 |
| \$ 191 Dibenzo(a,h)anthracene-d14 | 292 | 19.739 | 19.751 | (1.130) | 889305 | 5.00000 | 4.910 |
| 79 Dibenzo(a,h)anthracene | 278 | 19.830 | 19.840 | (1.135) | 978944 | 5.00000 | 4.927 |
| 80 Benzo(g,h,i)perylene | 276 | 20.666 | 20.676 | (1.183) | 1044239 | 5.00000 | 4.846 |

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-------------|-----------|--------|---------|---------|----------|--------------------|-------------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| ----- | ---- | == | ===== | ===== | ===== | ===== | ===== |
| 99 Perylene | 252 | 17.533 | 17.536 | (1.004) | 808709 | 5.00000 | 4.786 |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

| | |
|--|-------------------------------|
| Instrument ID: nt12.i | Calibration Date: 31-MAR-2011 |
| Lab File ID: 03311106.d | Calibration Time: 20:27 |
| Lab Smp Id: IC50331 | Client Smp ID: IC50331 |
| Analysis Type: SV | Level: |
| Quant Type: ISTD | Sample Type: |
| Operator: JZ | |
| Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m | |
| Misc Info: 11- | |

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|--------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 424290 | -14.13 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 233090 | -16.78 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 394000 | -14.60 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 423013 | -15.93 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 385980 | -12.72 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | 0.00 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | 0.00 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.03 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.81 | 0.02 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.47 | 0.00 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/nt12.i/20110331.b/03311106.d

Date : 31-MAR-2011 22:19

Client ID: IC50331

Sample Info: IC50331,

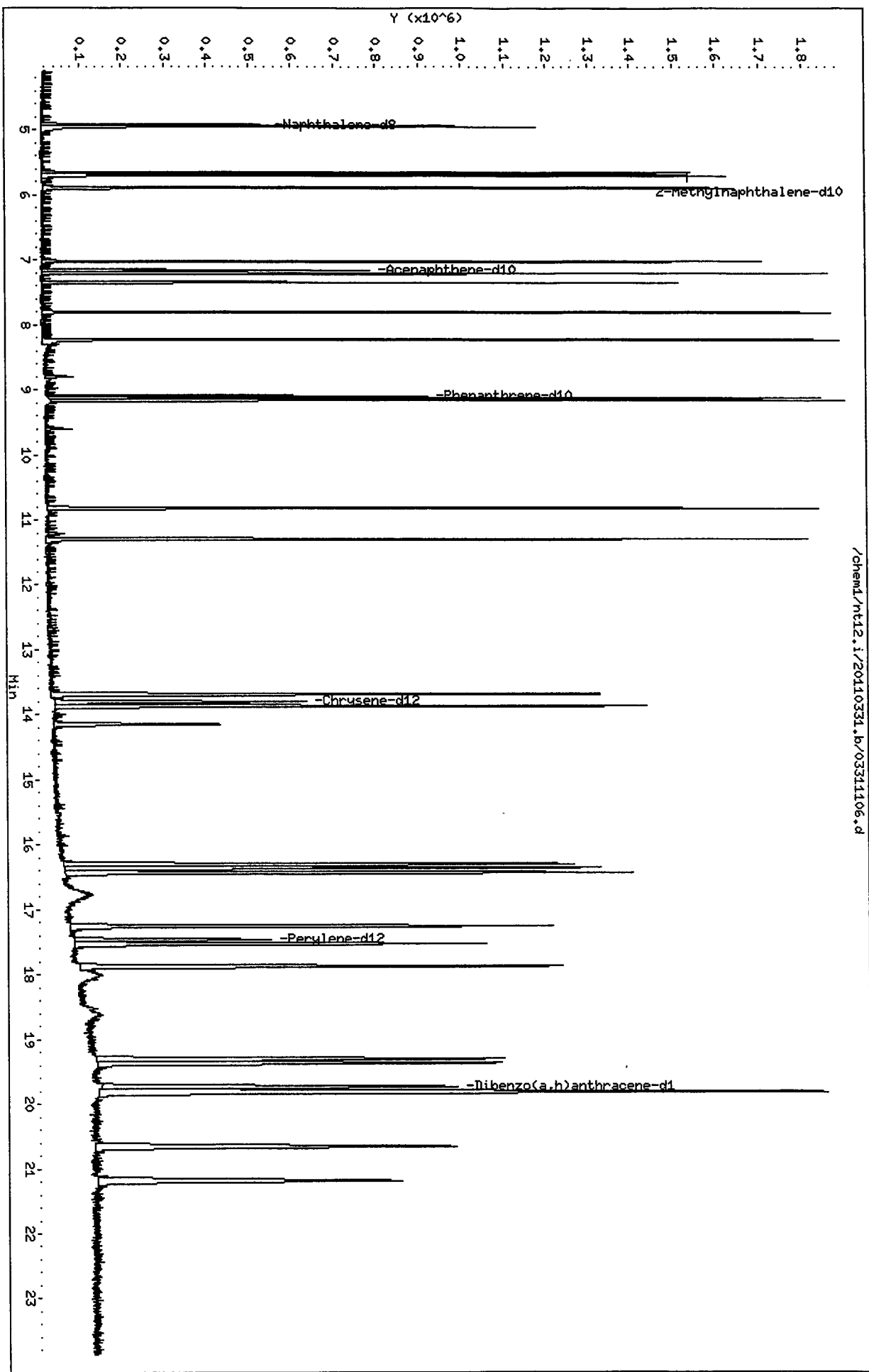
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110331.b/03311106.d



CO-ELUTION SUMMARY FOR FILE - 03311106.d

Lab ID: IC50331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00179

Analytical Resources, Inc.

Semivolatle Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110331.b/03311107.d
 Lab Smp Id: IC100331 Client Smp ID: IC10331
 Inj Date : 31-MAR-2011 22:47
 Operator : JZ Inst ID: nt12.i
 Smp Info : IC100331,
 Misc Info : 11-
 Comment : 1ul Injection
 Method : /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Meth Date : 01-Apr-2011 17:46 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 7 Calibration Sample, Level: 6
 Dil Factor: 1.00000
 Integrator: HP RTE
 Target Version: 3.50
 Compound Sublist: pnax.sub

Q 04/01/11

| Compounds | QUANT SIG | | AMOUNTS | | | ON-COL | |
|-------------------------------------|-----------|--------|---------|---------|----------|----------|-----------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | | CAL-AMT (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.917 | 4.917 | (1.000) | 479714 | 2.00000 | |
| 28 Naphthalene | 128 | 4.945 | 4.945 | (1.006) | 1995389 | 10.00000 | 8.614 |
| \$ 190 2-Methylnaphthalene-d10 | 152 | 5.655 | 5.655 | (1.150) | 1205283 | 10.00000 | 8.218 |
| 32 2-Methylnaphthalene | 141 | 5.700 | 5.700 | (1.159) | 1055784 | 10.00000 | 8.202 |
| 105 1-methylnaphthalene | 141 | 5.892 | 5.892 | (1.198) | 1090496 | 10.00000 | 7.961 |
| 40 Acenaphthylene | 152 | 7.025 | 7.025 | (0.981) | 1909756 | 10.00000 | 8.658 |
| * 42 Acenaphthene-d10 | 164 | 7.164 | 7.164 | (1.000) | 252063 | 2.00000 | |
| 44 Acenaphthene | 153 | 7.208 | 7.208 | (1.006) | 1194013 | 10.00000 | 8.691 |
| 46 Dibenzofuran | 168 | 7.353 | 7.353 | (1.026) | 1599433 | 10.00000 | 8.563 |
| 49 Fluorene | 166 | 7.808 | 7.808 | (1.090) | 1361038 | 10.00000 | 8.849 |
| * 59 Phenanthrene-d10 | 188 | 9.098 | 9.098 | (1.000) | 420234 | 2.00000 | |
| 60 Phenanthrene | 178 | 9.133 | 9.133 | (1.004) | 1955746 | 10.00000 | 8.818 |
| 61 Anthracene | 178 | 9.164 | 9.164 | (1.007) | 1861190 | 10.00000 | 8.358 |
| 64 Fluoranthene | 202 | 10.827 | 10.827 | (1.190) | 2094104 | 10.00000 | 8.683 |
| 65 Pyrene | 202 | 11.301 | 11.301 | (0.819) | 2156305 | 10.00000 | 8.807 |
| 68 Benzo (a) anthracene | 228 | 13.693 | 13.693 | (0.992) | 1989377 | 10.00000 | 9.100 |
| * 69 Chrysene-d12 | 240 | 13.806 | 13.806 | (1.000) | 454758 | 2.00000 | |
| 71 Chrysene | 228 | 13.873 | 13.873 | (1.005) | 1919561 | 10.00000 | 8.984 |
| 74 Benzo (b) fluoranthene | 252 | 16.312 | 16.312 | (0.934) | 2134788 | 10.00000 | 9.322 |
| 75 Benzo (k) fluoranthene | 252 | 16.369 | 16.369 | (0.937) | 2182504 | 10.00000 | 9.086 |
| 188 Benzo (j) fluoranthene | 252 | 16.444 | 16.444 | (0.941) | 2042956 | 10.00000 | 8.900 |
| 76 Benzo (a) pyrene | 252 | 17.268 | 17.268 | (0.988) | 1913767 | 10.00000 | 9.102 |
| * 77 Perylene-d12 | 264 | 17.473 | 17.473 | (1.000) | 406204 | 2.00000 | |
| 78 Indeno (1,2,3-cd) pyrene | 276 | 19.833 | 19.833 | (1.135) | 2407866 | 10.00000 | 9.418 |
| \$ 191 Dibenzo (a,h) anthracene-d14 | 292 | 19.751 | 19.751 | (1.130) | 1771889 | 10.00000 | 9.296 |
| 79 Dibenzo (a,h) anthracene | 278 | 19.840 | 19.840 | (1.135) | 1982350 | 10.00000 | 9.481 |
| 80 Benzo (g,h,i) perylene | 276 | 20.676 | 20.676 | (1.183) | 2073848 | 10.00000 | 9.146 |

| Compounds | QUANT SIG | | AMOUNTS | | | | |
|-------------|-----------|--------|---------|---------|----------|--------------------|-------------------|
| | MASS | RT | EXP RT | REL RT | RESPONSE | CAL-AMT (ug/mL) | ON-COL (ug/mL) |
| ===== | ==== | == | ===== | ===== | ===== | ===== | ===== |
| 99 Perylene | 252 | 17.536 | 17.536 | (1.004) | 1608876 | 10.0000 | 9.048 |

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 03311107.d
 Lab Smp Id: IC100331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110331.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 31-MAR-2011
 Calibration Time: 20:27
 Client Smp ID: IC10331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|--------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 479714 | -2.91 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 252063 | -10.01 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 420234 | -8.91 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 454758 | -9.62 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 406204 | -8.14 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.92 | 4.42 | 5.42 | 4.92 | -0.07 |
| 42 Acenaphthene-d10 | 7.16 | 6.66 | 7.66 | 7.16 | 0.00 |
| 59 Phenanthrene-d10 | 9.10 | 8.60 | 9.60 | 9.10 | 0.00 |
| 69 Chrysene-d12 | 13.80 | 13.30 | 14.30 | 13.81 | 0.02 |
| 77 Perylene-d12 | 17.47 | 16.97 | 17.97 | 17.47 | 0.02 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Data File: /chem1/nt12.i/20110331.b/03311107.d

Date : 31-MAR-2011 22:47

Client ID: IC100331

Sample Info: IC100331,

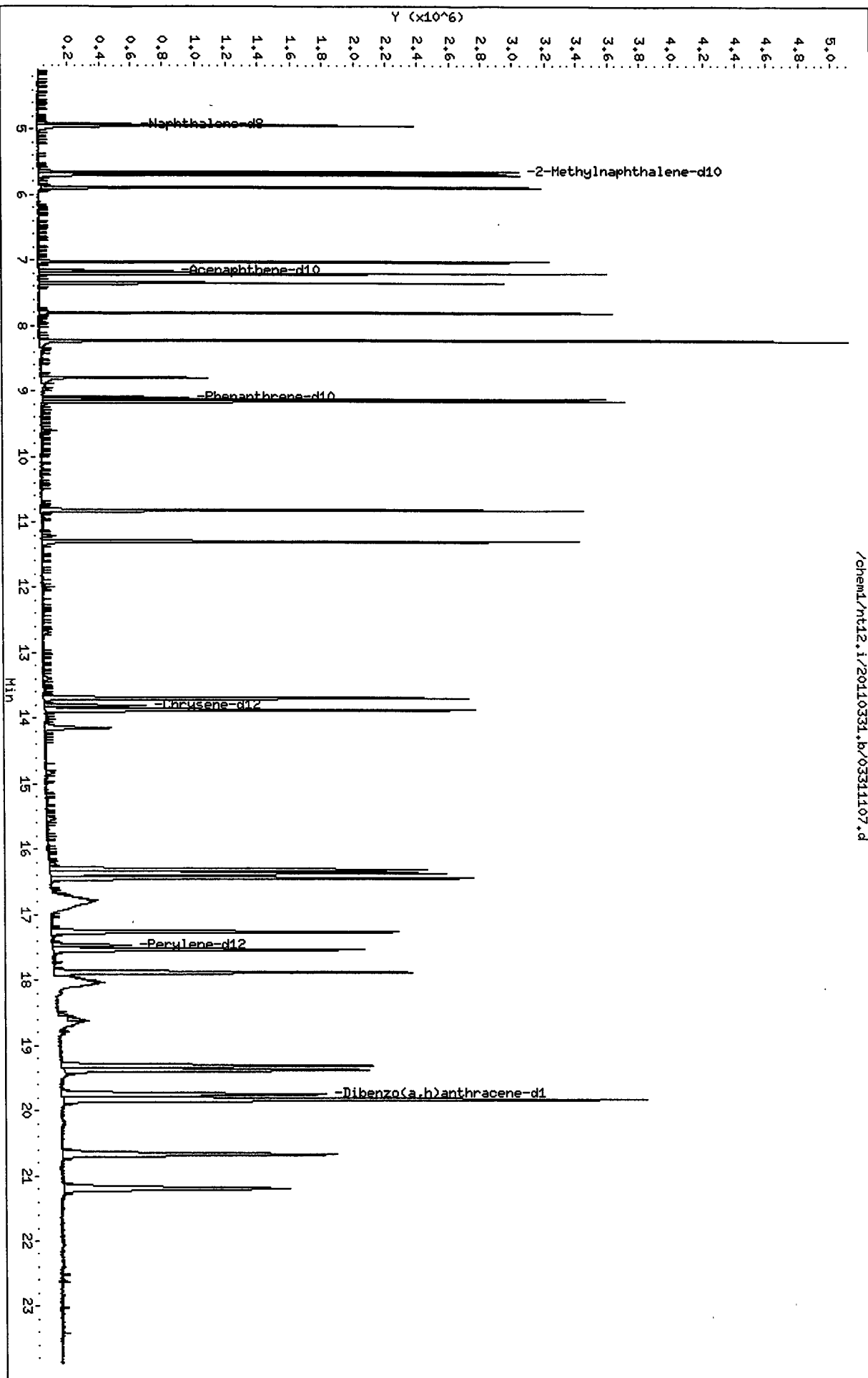
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110331.b/03311107.d



CO-ELUTION SUMMARY FOR FILE - 03311107.d

Lab ID: IC100331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 31-MAR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00184

Analytical Resources, Inc.

Semivolatle Report SW846 Method 8270D

Data file : /chem1/nt12.i/20110401.b/04011103.d
 Lab Smp Id: ICV0331 Client Smp ID: ICV0331
 Inj Date : 01-APR-2011 15:53
 Operator : JZ Inst ID: nt12.i
 Smp Info : ICV0331
 Misc Info : 11-
 Comment : lul Injection
 Method : /chem1/nt12.i/20110401.b/SIMPNA0331.m
 Meth Date : 04-Apr-2011 08:48 jianqing Quant Type: ISTD
 Cal Date : 31-MAR-2011 22:47 Cal File: 03311107.d
 Als bottle: 3 QC Sample: LCS
 Dil Factor: 1.00000
 Integrator: HP RTE Compound Sublist: pnax.sub
 Target Version: 3.50

B *04/04/11*

| Compounds | QUANT SIG | MASS | RT | EXP RT | REL RT | RESPONSE | CONCENTRATIONS | |
|-----------------------------------|-----------|------------------------|--------|---------|--------|----------|-------------------|---------------|
| | | | | | | | ON-COLUMN (ug/mL) | FINAL (ug/mL) |
| * 27 Naphthalene-d8 | 136 | 4.851 | 4.851 | (1.000) | 405752 | 2.00000 | - | |
| 28 Naphthalene | 128 | 4.876 | 4.879 | (1.005) | 524717 | 2.67821 | 2.678 (R) | |
| \$ 190 2-Methylnaphthalene-d10 | 152 | Compound Not Detected. | | | | | | |
| 32 2-Methylnaphthalene | 141 | 5.630 | 5.630 | (1.161) | 294338 | 2.70344 | 2.703 (R) | |
| 105 1-methylnaphthalene | 141 | 5.823 | 5.823 | (1.200) | 265630 | 2.29260 | 2.293 | |
| 40 Acenaphthylene | 152 | 6.952 | 6.952 | (0.980) | 496272 | 2.32632 | 2.326 | |
| * 42 Acenaphthene-d10 | 164 | 7.091 | 7.091 | (1.000) | 243785 | 2.00000 | | |
| 44 Acenaphthene | 153 | 7.139 | 7.139 | (1.007) | 304522 | 2.29176 | 2.292 | |
| 46 Dibenzofuran | 168 | 7.281 | 7.281 | (1.027) | 450265 | 2.49236 | 2.492 | |
| 49 Fluorene | 166 | 7.735 | 7.738 | (1.091) | 353391 | 2.37571 | 2.376 | |
| * 59 Phenanthrene-d10 | 188 | 9.026 | 9.029 | (1.000) | 404374 | 2.00000 | | |
| 60 Phenanthrene | 178 | 9.060 | 9.060 | (1.004) | 500017 | 2.34292 | 2.343 | |
| 61 Anthracene | 178 | 9.092 | 9.092 | (1.007) | 526287 | 2.45603 | 2.456 | |
| 64 Fluoranthene | 202 | 10.733 | 10.736 | (1.189) | 548293 | 2.36264 | 2.363 | |
| 65 Pyrene | 202 | 11.200 | 11.203 | (0.819) | 564584 | 2.46598 | 2.466 | |
| 68 Benzo(a)anthracene | 228 | 13.570 | 13.570 | (0.992) | 499141 | 2.44164 | 2.442 | |
| * 69 Chrysene-d12 | 240 | 13.680 | 13.683 | (1.000) | 425229 | 2.00000 | | |
| 71 Chrysene | 228 | 13.749 | 13.753 | (1.005) | 475813 | 2.38146 | 2.381 | |
| 74 Benzo(b)fluoranthene | 252 | 16.164 | 16.167 | (0.933) | 509580 | 2.48981 | 2.490 | |
| 75 Benzo(k)fluoranthene | 252 | 16.217 | 16.227 | (0.936) | 535096 | 2.49240 | 2.492 | |
| 188 Benzo(j)fluoranthene | 252 | Compound Not Detected. | | | | | | |
| 76 Benzo(a)pyrene | 252 | 17.113 | 17.117 | (0.988) | 427034 | 2.27255 | 2.273 | |
| * 77 Perylene-d12 | 264 | 17.322 | 17.325 | (1.000) | 363042 | 2.00000 | | |
| 78 Indeno(1,2,3-cd)pyrene | 276 | 19.635 | 19.644 | (1.134) | 562724 | 2.46275 | 2.463 | |
| \$ 191 Dibenzo(a,h)anthracene-d14 | 292 | Compound Not Detected. | | | | | | |
| 79 Dibenzo(a,h)anthracene | 278 | 19.654 | 19.660 | (1.135) | 461801 | 2.47126 | 2.471 | |
| 80 Benzo(g,h,i)perylene | 276 | 20.458 | 20.471 | (1.181) | 485824 | 2.39724 | 2.397 | |
| 99 Perylene | 252 | Compound Not Detected. | | | | | | |

| Compounds | QUANT SIG | | | | | | CONCENTRATIONS | |
|-----------|-----------|----|--------|--------|----------|----------------------|------------------|--|
| | MASS | RT | EXP RT | REL RT | RESPONSE | ON-COLUMN (ug/mL) | FINAL (ug/mL) | |
| ===== | ==== | == | ===== | ===== | ===== | ===== | ===== | |

QC Flag Legend

R - Spike/Surrogate failed recovery limits.

Analytical Resources, Inc.

INTERNAL STANDARD COMPOUNDS
 AREA AND RT SUMMARY

Instrument ID: nt12.i
 Lab File ID: 04011103.d
 Lab Smp Id: ICV0331
 Analysis Type: SV
 Quant Type: ISTD
 Operator: JZ
 Method File: /chem1/nt12.i/20110401.b/SIMPNA0331.m
 Misc Info: 11-

Calibration Date: 01-APR-2011
 Calibration Time: 15:16
 Client Smp ID: ICV0331
 Level:
 Sample Type:

Test Mode:
 Use Initial Calibration Level 4.

| COMPOUND | STANDARD | AREA LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|------------|---------|--------|--------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 494112 | 247056 | 988224 | 405752 | -17.88 |
| 42 Acenaphthene-d10 | 280105 | 140052 | 560210 | 243785 | -12.97 |
| 59 Phenanthrene-d10 | 461353 | 230676 | 922706 | 404374 | -12.35 |
| 69 Chrysene-d12 | 503160 | 251580 | 1006320 | 425229 | -15.49 |
| 77 Perylene-d12 | 442215 | 221108 | 884430 | 363042 | -17.90 |

| COMPOUND | STANDARD | RT LIMIT | | SAMPLE | %DIFF |
|---------------------|----------|----------|-------|--------|-------|
| | | LOWER | UPPER | | |
| 27 Naphthalene-d8 | 4.85 | 4.35 | 5.35 | 4.85 | 0.00 |
| 42 Acenaphthene-d10 | 7.09 | 6.59 | 7.59 | 7.09 | 0.00 |
| 59 Phenanthrene-d10 | 9.03 | 8.53 | 9.53 | 9.03 | -0.04 |
| 69 Chrysene-d12 | 13.68 | 13.18 | 14.18 | 13.68 | -0.02 |
| 77 Perylene-d12 | 17.32 | 16.82 | 17.82 | 17.32 | -0.02 |

AREA UPPER LIMIT = +100% of internal standard area.
 AREA LOWER LIMIT = - 50% of internal standard area.
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT.
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT.

Analytical Resources, Inc.

RECOVERY REPORT

Client Name: Client SDG: 20110401
 Sample Matrix: NONE Fraction: SV
 Lab Smp Id: ICV0331 Client Smp ID: ICV0331
 Level: Operator: JZ
 Data Type: MS DATA SampleType: LCS
 SpikeList File: pnalcss.spk Quant Type: ISTD
 Sublist File: pnax.sub
 Method File: /chem1/nt12.i/20110401.b/SIMPNA0331.m
 Misc Info: 11-

| SPIKE COMPOUND | AMOUNT ADDED ug/mL | AMOUNT RECOVERED ug/mL | % RECOVERED | LIMITS |
|-------------------------|--------------------|------------------------|-------------|--------|
| 28 Naphthalene | 2.505 | 2.678 | 106.91* | |
| 32 2-Methylnaphthalen | 2.505 | 2.703 | 107.92* | |
| 105 1-methylnaphthalen | 2.505 | 2.293 | 91.52 | |
| 40 Acenaphthylene | 2.505 | 2.326 | 92.87 | |
| 44 Acenaphthene | 2.505 | 2.292 | 91.49 | |
| 46 Dibenzofuran | 2.505 | 2.492 | 99.50 | |
| 49 Fluorene | 2.505 | 2.376 | 94.84 | |
| 60 Phenanthrene | 2.505 | 2.343 | 93.53 | |
| 61 Anthracene | 2.505 | 2.456 | 98.04 | |
| 64 Fluoranthene | 2.505 | 2.363 | 94.32 | |
| 65 Pyrene | 2.505 | 2.466 | 98.44 | |
| 68 Benzo(a)anthracene | 2.505 | 2.442 | 97.47 | |
| 71 Chrysene | 2.505 | 2.381 | 95.07 | |
| 74 Benzo(b)fluoranthene | 2.505 | 2.490 | 99.39 | |
| 75 Benzo(k)fluoranthene | 2.505 | 2.492 | 99.50 | |
| 76 Benzo(a)pyrene | 2.505 | 2.273 | 90.72 | |
| 78 Indeno(1,2,3-cd)py | 2.505 | 2.463 | 98.31 | |
| 79 Dibenzo(a,h)anthra | 2.505 | 2.471 | 98.65 | |
| 80 Benzo(g,h,i)perylene | 2.505 | 2.397 | 95.70 | |
| 99 Perylene | 2.505 | 0.000 | <i>NR</i> * | |

| SURROGATE COMPOUND | AMOUNT ADDED ug/mL | AMOUNT RECOVERED ug/mL | % RECOVERED | LIMITS |
|----------------------------|--------------------|------------------------|---------------------------|--------|
| \$ 190 2-Methylnaphthalene | 1.875 | 0.000 | * | |
| \$ 191 Dibenzo(a,h)anthr | 1.875 | 0.000 | <i>no second source</i> * | |

\$ 06/06/11

Data File: /chem1/nt12.i/20110401.b/04011103.d

Date: 01-APR-2011 15:53

Client ID: ICV0331

Sample Info: ICV0331

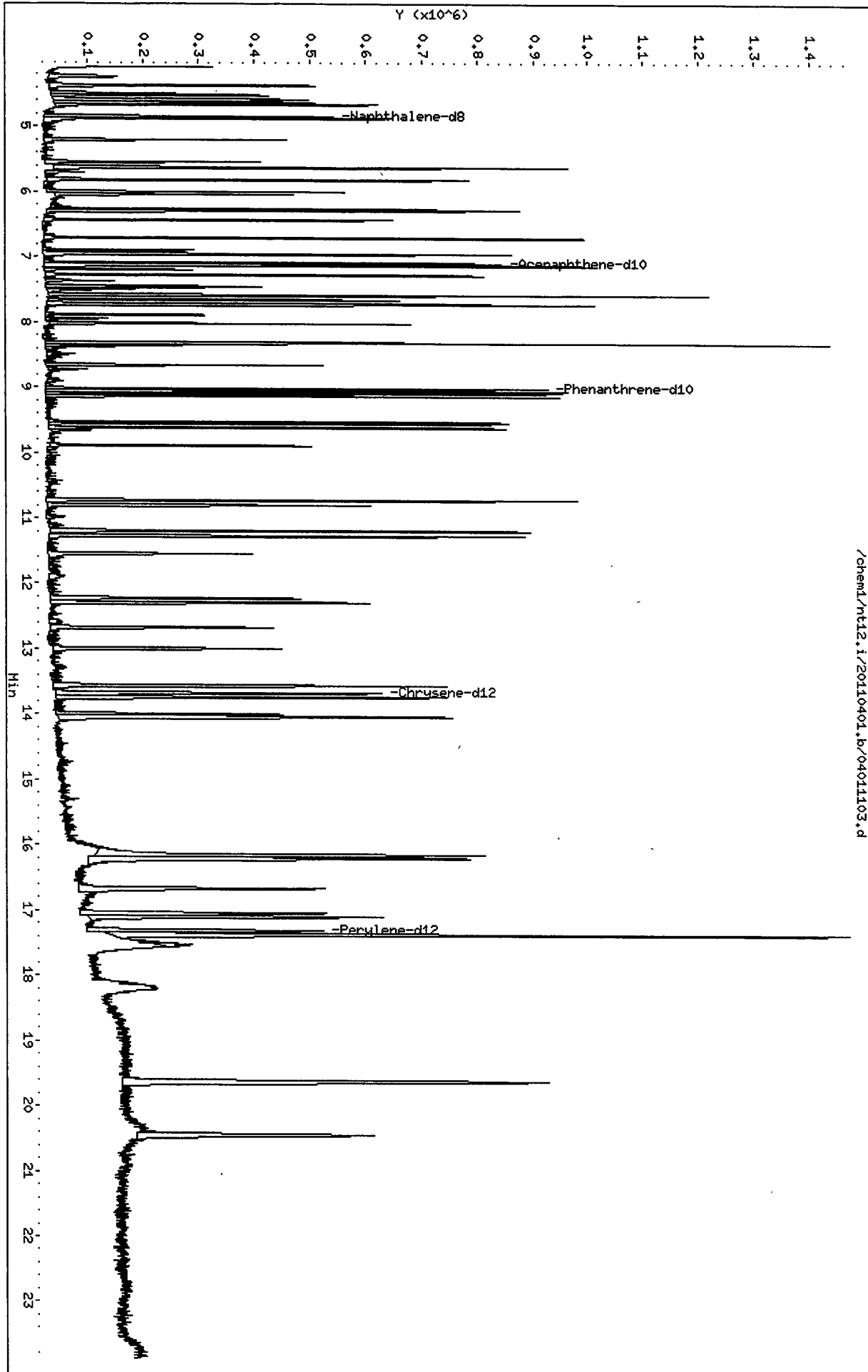
Column phase: ZB35

Instrument: nt12.i

Operator: JZ

Column diameter: 0.32

/chem1/nt12.i/20110401.b/04011103.d



CO-ELUTION SUMMARY FOR FILE - 04011103.d

Lab ID: ICV0331, Method: SIMPNA0331.m, Instrument: nt12.i, Date: 01-APR-2011

RT CO-ELUTION COMPOUNDS

NO CO-ELUTIONS

SN54 : 00190

SIM PAH Raw Data
Run Logs, Continuing Calibrations, and Raw Data

ARI Job ID: SN54

GC/MS SVOA Analyst Notes / Corrective Action Log

ARI Project ID: SN54 Client ID: Floyd Spitzer

ARI SOP: 801S(SIM-PNA) 802S(Butyl Tins) 804S(SVOA-8270D) 805S(op-Pest)

Parameter(s): sim PNA

Instrument: NT-2 NT-4 NT-6 NT-8 NT11, (NT12)

Curve Date: 3/31/11 Analysis Start Date: 4/1; 4/4/11

| | | | |
|--------------------------------|----------------------|-----------------------------------|------------------------|
| DFTPP Tune Meets Criteria? | <u>YES</u> / NO | Internal Standard Meets Criteria? | <u>YES</u> / <u>NO</u> |
| DDT Breakdown <20%? | <u>YES</u> / NO / NA | Method Blank In Control? | <u>YES</u> / NO |
| Peak Tailing Factor ≤2? | <u>YES</u> / NO / NA | LCS / LCSD Recovery In Control? | <u>YES</u> / NO |
| ICal acceptable? | <u>YES</u> / NO | CCal acceptable? | <u>YES</u> / NO |
| Q flag applied? | YES / <u>NO</u> | Q flag applied? | YES / <u>NO</u> |
| Surrogate Recovery in Control? | <u>YES</u> / NO | Special Analysis Criteria Met? | YES / NO / <u>NA</u> |
| Manual Integrations for ICal? | <u>YES</u> / NO | Manual Integrations for Samples? | <u>Yes</u> / NO |

Detail problems, corrective actions and/or other pertinent information below (use reverse side when necessary):

4/1 samples SN54 A-D + MB/LCS/LCSD
 4/4 samples E-H + MS/MSD + Dilutions for samples F-H.
 Due to matrix effect, 15, perylene-d12, response out of DC limit at low bias in all the dilutions runs. PM was informed. no corrective action needed. Forms included.

Additional Details on Reverse: Yes / No

Analyst: [Signature] Date: 04/05/11
 Reviewer: [Signature] Date: 4/6/11

Analytical Resources Inc.: Organics Instrument Log

NT-12 Serial No.: GC=US00032558, MS= US01180091

Date: 4/1/11 Analysis: SIM PAA Analyst: D
 GC Program: SIM PAA 25 Column No: 157992 Column Type: ZB35
 Instrument Tune (.U or .CT.): 40331 EM Voltage: 2682
 Calibration File: 041110V Curve Date: 3/31/11 Injection Vol.: _____

| IS/SS | Ical/Ccal | LCS/ICV |
|---------------|---------------|---------|
| <u>1754-1</u> | <u>1818-1</u> | |
| | <u>788-3</u> | |
| | | |
| | | |
| | | |
| | | |

Document All Maintenance Tasks In StarLIMS

INTERNAL STANDARD SUMMARY FOR DATABATCH - /chem1/nt12.i/20110401.b

| Time | Filename | LabID | ClientId | DF | |
|------|----------|------------|------------|--------------|---|
| 1 | 1457 | 04011101.d | DFTPP0401 | DFTPP0401 | 1 NO ISTDs FOUND |
| 2 | 1516 | 04011102.d | CC0401 | CC0401 | 1 4.85 499536 7.09 294599 9.03 482436 13.68 526425 17.32 439084 |
| 3 | 1553 | 04011103.d | ICV0331 | ICV0331 | 1 4.85 405752 7.09 243785 9.03 404374 13.68 425229 17.32 363042 |
| 4 | 1621 | 04011104.d | SN54MBS2 | SN54MBS2 | 1 4.84 362184 7.08 207575 9.02 332402 13.67 371865 17.32 319297 |
| 5 | 1649 | 04011105.d | SN54LCSS2 | SN54LCSS2 | 1 4.84 366499 7.08 217644 9.02 347080 13.67 381079 17.32 335486 |
| 6 | 1717 | 04011106.d | SN54LCSDS2 | SN54LCSDS2 | 1 4.84 364394 7.08 218342 9.02 356245 13.68 383014 17.32 345990 |
| 7 | 1745 | 04011107.d | SN54QLS | SN54QLS | 1 4.84 354887 7.08 221805 9.02 359036 13.68 388118 17.31 344619 |
| 8 | 1813 | 04011108.d | SN54A | LL-SED3-0-36 | 1 4.84 342514 7.09 202181 9.03 322755 13.73 373966 17.39 298676 |
| 9 | 1841 | 04011109.d | SN54B | LL-SED3-36-1 | 3 4.84 301159 7.09 187766 9.02 301621 13.69 322033 17.35 280959 |
| 10 | 1909 | 04011110.d | SN54C | LL-SED3-141- | 3 4.85 262380 7.09 167023 9.03 271187 13.69 293472 17.36 247444 |
| 11 | 1937 | 04011111.d | SN54D | LL-SED2-0-56 | 1 4.84 249435 7.09 163363 9.03 258445 13.73 339995 17.38 242975 |
| 12 | 2038 | 04011112.d | SN54E | LL-SED2-56-1 | 3 4.85 261670 7.09 166252 9.03 276567 13.70 293916 17.36 214585 |
| 13 | 2106 | 04011113.d | SN54EMS | LL-SED2-56-1 | 3 4.85 231693 7.09 145320 9.03 254871 13.70 270090 17.36 201254 |
| 14 | 2134 | 04011114.d | SN54EMSD | LL-SED2-56-1 | 3 4.85 232782 7.09 152172 9.03 249322 13.70 277173 17.36 190902 |
| 15 | 2202 | 04011115.d | SN54B | LL-SED3-36-1 | 1 4.85 202894 7.09 131352 9.03 218351 13.72 239578 17.40 164864 |
| 16 | 2230 | 04011116.d | SN54C | LL-SED3-141- | 1 4.85 211721 7.09 131101 9.03 230943 13.72 243922 17.41 161259 |

D 04/01/11

Every line must contain information or be lined out. Make all entries legible.
 Start a new page for each QC period. Document All Maintenance Tasks In StarLIMS

MANUAL INTEGRATION SUMMARY FOR DATABATCH - /chem1/nt12.i/20110401.b

ARI Job No.: CC04 Method: SIMPNA0331.m Instrument: nt12.i Date: 01-APR-2011

B. off/05/11

| Time | Filename | LabID | ClientId | DF | Manually Integrated Compounds |
|------|------------|------------|------------|----|---|
| 1516 | 04011102.d | CC0401 | CC0401 | 1 | NO MANUAL INTEGRATION |
| 1621 | 04011104.d | SN54MBS2 | SN54MBS2 | 1 | NO MANUAL INTEGRATION |
| 1649 | 04011105.d | SN54LCSS2 | SN54LCSS2 | 1 | NO MANUAL INTEGRATION |
| 1717 | 04011106.d | SN54LCSDS2 | SN54LCSDS2 | 1 | NO MANUAL INTEGRATION |
| 1813 | 04011108.d | SN54A | LL-SED3-0- | 1 | NO MANUAL INTEGRATION |
| 1841 | 04011109.d | SN54B | LL-SED3-36 | 3 | NO MANUAL INTEGRATION |
| 1909 | 04011110.d | SN54C | LL-SED3-14 | 3 | NO MANUAL INTEGRATION |
| 1937 | 04011111.d | SN54D | LL-SED2-0- | 1 | Dibenzo(a,h)anthracene, Benzo(g,h,i)perylene, |

Q-FLAG SUMMARY FOR DATABATCH - /chem1/nt12.i/20110401.b

Instrument: nt12.i Date: 01-APR-2011 Method: SIMPNA0331.m

INITIAL CAL: 31-MAR-2011

| Compound | %RSD or R ² |
|------------|------------------------|
| ----- | |
| NO Q-FLAGS | |
| ----- | |

CONTINUING CAL: 01-APR-2011

| Compound | %D |
|------------|----|
| ----- | |
| NO Q-FLAGS | |
| ----- | |

JB 04/01/11