

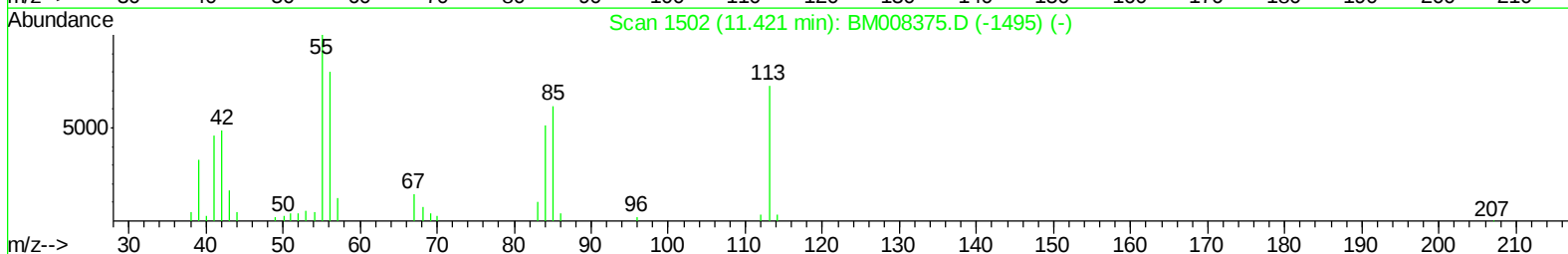
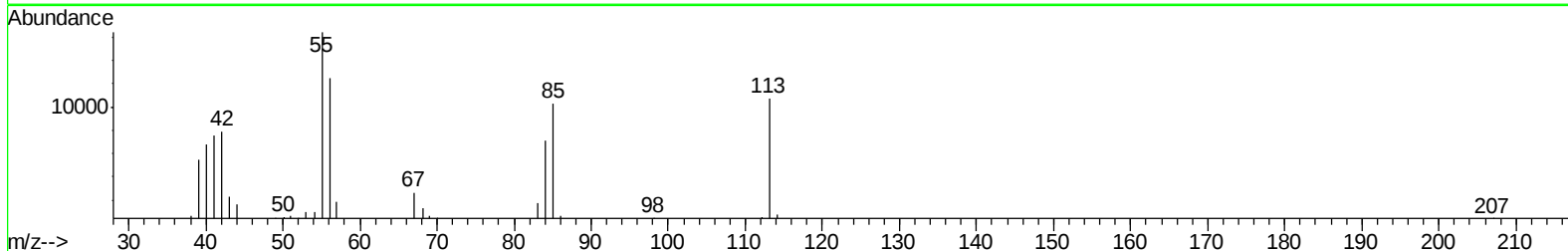
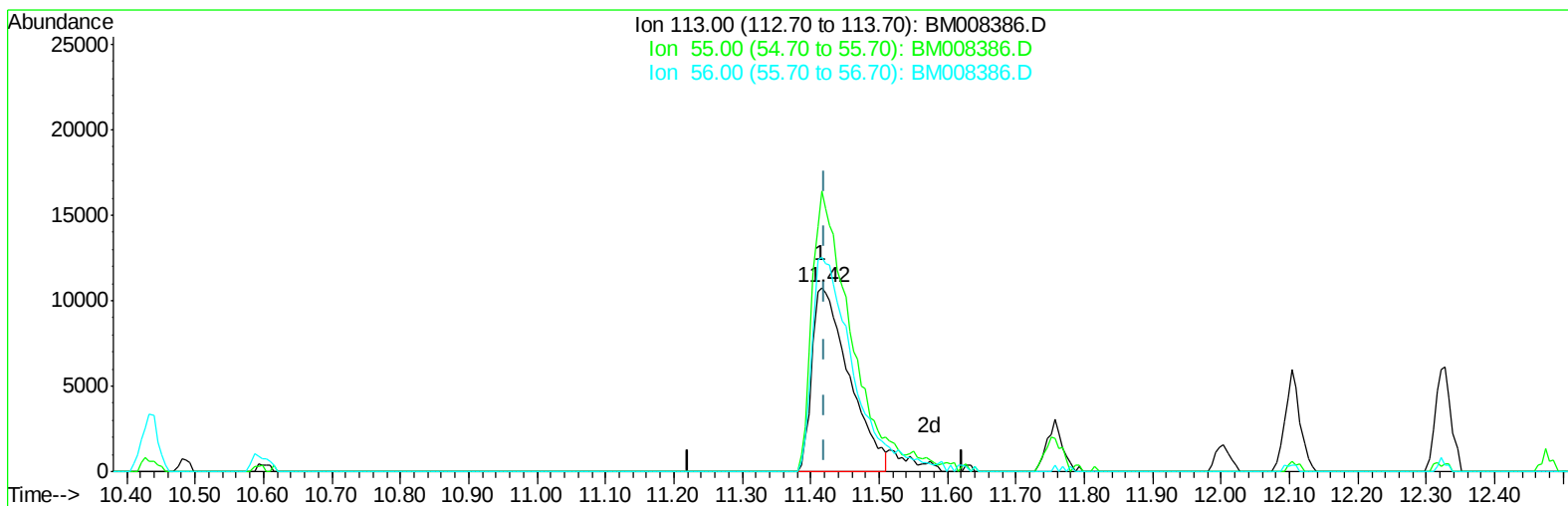
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121716\  
Data File : BM008386.D  
Acq On : 16 Dec 2016 22:50  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
LabSampleId :  
SSTD02060

Manual Integrations  
APPROVED

umangi  
12/20/2016 10:43:10 AM

Quant Time: Dec 17 02:20:09 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM12  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Dec 17 02:12:55 2016  
Response via : Initial Calibration



TIC: BM008386.D

(32) Caprolactam

11.416min (-0.006) 18.88ng/ul

response 40323

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 157.60 | 152.52 |
| 56.00  | 114.10 | 116.36 |
| 0.00   | 0.00   | 0.00   |

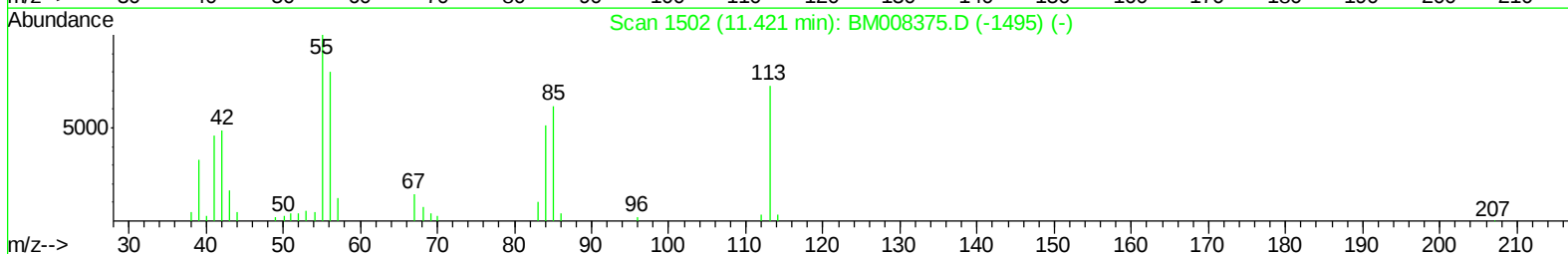
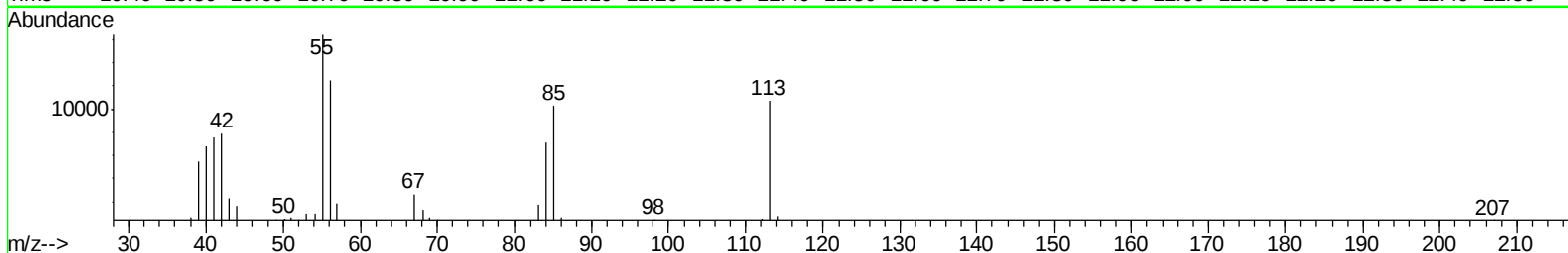
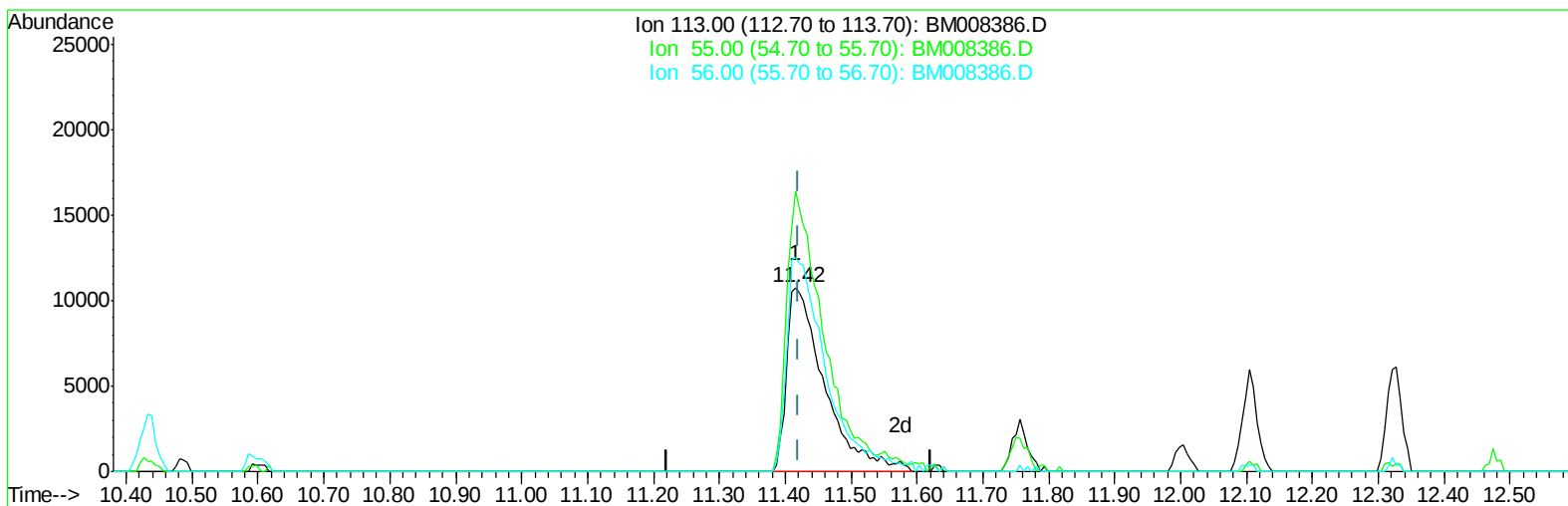
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121716\  
Data File : BM008386.D  
Acq On : 16 Dec 2016 22:50  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_M  
LabSampleId :  
SSTD02060

Manual Integrations  
APPROVED

umangi  
12/20/2016 10:43:10 AM

Quant Time: Dec 17 02:20:09 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM12  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Dec 17 02:12:55 2016  
Response via : Initial Calibration



TIC: BM008386.D

(32) Caprolactam

11.416min (-0.006) 20.33ng/ul m

response 43420

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 157.60 | 152.52 |
| 56.00  | 114.10 | 116.36 |
| 0.00   | 0.00   | 0.00   |

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121716\  
 Data File : BM008386.D  
 Acq On : 16 Dec 2016 22:50  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02060

Manual Integrations  
 APPROVED

umangi  
 12/20/2016 10:43:10 AM

Quant Time: Dec 17 03:19:42 2016

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121

Quant Title : SVOA CALIBRATION

QLast Update : Sat Dec 17 02:12:55 2016

Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.66  | 152  | 118690   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 460259   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.30 | 164  | 302871   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.06 | 188  | 619970   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.26 | 240  | 796854   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.49 | 264  | 808846   | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.23  | 96  | 21505  | 8.53  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.85  | 99  | 187375 | 20.35 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.01  | 67  | 108405 | 20.48 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.20  | 132 | 147843 | 21.34 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.38  | 113 | 144754 | 20.30 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.82  | 128 | 71497  | 21.50 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.54  | 143 | 78757  | 21.37 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.07 | 165 | 148786 | 22.03 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.59 | 131 | 171032 | 22.09 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.72 | 166 | 430425 | 21.38 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.99 | 160 | 523092 | 21.98 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.56 | 143 | 48358  | 16.30 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.30 | 176 | 375354 | 21.45 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.46 | 200 | 67840  | 18.38 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.16 | 188 | 587935 | 21.77 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.46 | 212 | 683558 | 22.21 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.34 | 264 | 745616 | 21.95 | ng/ul | 0.00 |

#### Target Compounds

|                                |       |     |        |       |       | Qvalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.26  | 88  | 23678  | 8.03  | ng/uL | 96     |
| 4) Benzaldehyde                | 6.83  | 77  | 136400 | 22.38 | ng/ul | 91     |
| 6) Phenol                      | 6.88  | 94  | 196374 | 20.54 | ng/ul | 96     |
| 8) Bis(2-Chloroethyl)ether     | 7.10  | 93  | 148890 | 20.86 | ng/ul | 94     |
| 10) 2-Chlorophenol             | 7.23  | 128 | 148191 | 21.18 | ng/ul | 99     |
| 11) 2-Methylphenol             | 8.11  | 108 | 141782 | 20.26 | ng/ul | 97     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.19  | 45  | 171103 | 21.08 | ng/ul | 98     |
| 14) Acetophenone               | 8.49  | 105 | 237385 | 20.36 | ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.47  | 70  | 123700 | 20.64 | ng/ul | 99     |
| 16) 4-Methylphenol             | 8.45  | 108 | 155580 | 20.50 | ng/ul | 99     |
| 17) Hexachloroethane           | 8.72  | 117 | 66869  | 20.96 | ng/ul | 96     |
| 20) Nitrobenzene               | 8.87  | 77  | 182911 | 21.33 | ng/ul | 99     |
| 21) Isophorone                 | 9.38  | 82  | 332867 | 21.59 | ng/ul | 98     |
| 23) 2-Nitrophenol              | 9.57  | 139 | 82856  | 21.98 | ng/ul | 91     |
| 24) 2,4-Dimethylphenol         | 9.63  | 107 | 182248 | 21.64 | ng/ul | 95     |
| 25) Bis(2-Chloroethoxy)methane | 9.86  | 93  | 187525 | 21.42 | ng/ul | 97     |
| 27) 2,4-Dichlorophenol         | 10.10 | 162 | 147507 | 21.67 | ng/ul | 98     |
| 28) Naphthalene                | 10.49 | 128 | 467391 | 21.45 | ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.62 | 127 | 173678 | 22.26 | ng/ul | 98     |
| 31) Hexachlorobutadiene        | 10.76 | 225 | 118299 | 21.38 | ng/ul | 96     |
| 32) Caprolactam                | 11.42 | 113 | 43420m | 20.33 | ng/ul |        |
| 33) 4-Chloro-3-methylphenol    | 11.76 | 107 | 157844 | 21.43 | ng/ul | 99     |
| 34) 2-Methylnaphthalene        | 12.10 | 142 | 335308 | 21.30 | ng/ul | 100    |

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM121716\  
 Data File : BM008386.D  
 Acq On : 16 Dec 2016 22:50  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02060

Manual Integrations  
 APPROVED

umangi  
 12/20/2016 10:43:10 AM

Quant Time: Dec 17 03:19:42 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Dec 17 02:12:55 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.48 | 216  | 208622   | 22.00 | ng/ul  | 99       |
| 37) Hexachlorocyclopentadiene  | 12.45 | 237  | 71769    | 18.74 | ng/ul  | 98       |
| 38) 2,4,6-Trichlorophenol      | 12.74 | 196  | 118541   | 22.02 | ng/ul  | 94       |
| 39) 2,4,5-Trichlorophenol      | 12.82 | 196  | 129024   | 21.97 | ng/ul  | 97       |
| 40) 1,1'-Biphenyl              | 13.13 | 154  | 439343   | 21.74 | ng/ul  | 98       |
| 41) 2-Chloronaphthalene        | 13.17 | 162  | 339452   | 21.91 | ng/ul  | 98       |
| 42) 2-Nitroaniline             | 13.40 | 65   | 101648   | 22.18 | ng/ul  | 96       |
| 44) Dimethylphthalate          | 13.77 | 163  | 422533   | 21.41 | ng/ul  | 99       |
| 45) 2,6-Dinitrotoluene         | 13.90 | 165  | 84348    | 21.94 | ng/ul  | 99       |
| 47) Acenaphthylene             | 14.02 | 152  | 529087   | 21.85 | ng/ul  | 99       |
| 48) 3-Nitroaniline             | 14.24 | 138  | 81165    | 21.19 | ng/ul  | 99       |
| 49) Acenaphthene               | 14.37 | 153  | 364184   | 21.88 | ng/ul  | 99       |
| 50) 2,4-Dinitrophenol          | 14.48 | 184  | 47527    | 21.32 | ng/ul  | 94       |
| 52) 4-Nitrophenol              | 14.57 | 109  | 47976    | 16.86 | ng/ul  | 99       |
| 53) Dibenzofuran               | 14.70 | 168  | 525300   | 21.66 | ng/ul  | 98       |
| 54) 2,4-Dinitrotoluene         | 14.71 | 165  | 125003   | 21.71 | ng/ul  | 96       |
| 55) 2,3,4,6-Tetrachlorophenol  | 14.94 | 232  | 118738   | 21.90 | ng/ul# | 96       |
| 56) Diethylphthalate           | 15.13 | 149  | 414914   | 21.14 | ng/ul  | 99       |
| 58) Fluorene                   | 15.36 | 166  | 429694   | 21.57 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.35 | 204  | 230659   | 21.78 | ng/ul  | 97       |
| 60) 4-Nitroaniline             | 15.42 | 138  | 78867    | 21.40 | ng/ul  | 98       |
| 63) 4,6-Dinitro-2-methylphenol | 15.47 | 198  | 68771    | 18.18 | ng/ul  | 97       |
| 64) N-Nitrosodiphenylamine     | 15.57 | 169  | 367873   | 21.80 | ng/ul  | 95       |
| 65) 4-Bromophenyl-phenylether  | 16.25 | 248  | 149032   | 21.90 | ng/ul  | 96       |
| 66) Hexachlorobenzene          | 16.36 | 284  | 164066   | 21.27 | ng/ul  | 98       |
| 67) Atrazine                   | 16.53 | 200  | 141570   | 20.47 | ng/ul  | 98       |
| 68) Pentachlorophenol          | 16.72 | 266  | 84239    | 20.19 | ng/ul  | 96       |
| 69) Phenanthrene               | 17.10 | 178  | 681755   | 21.46 | ng/ul  | 99       |
| 71) Anthracene                 | 17.19 | 178  | 703430   | 21.81 | ng/ul  | 99       |
| 72) Carbazole                  | 17.47 | 167  | 599627   | 21.11 | ng/ul  | 100      |
| 73) Di-n-butylphthalate        | 18.02 | 149  | 682601   | 21.49 | ng/ul  | 99       |
| 74) Fluoranthene               | 19.13 | 202  | 829250   | 21.25 | ng/ul  | 100      |
| 77) Pyrene                     | 19.49 | 202  | 864672   | 22.10 | ng/ul  | 99       |
| 78) Butylbenzylphthalate       | 20.39 | 149  | 317584   | 22.63 | ng/ul  | 97       |
| 79) 3,3'-Dichlorobenzidine     | 21.18 | 252  | 299196   | 21.42 | ng/ul  | 95       |
| 80) Benzo(a)anthracene         | 21.24 | 228  | 899775   | 21.85 | ng/ul  | 100      |
| 81) Bis(2-ethylhexyl)phthalate | 21.16 | 149  | 460320   | 22.40 | ng/ul  | 99       |
| 82) Chrysene                   | 21.29 | 228  | 860404   | 21.72 | ng/ul  | 98       |
| 84) Di-n-octyl phthalate       | 22.03 | 149  | 826403   | 21.29 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 22.81 | 252  | 940791   | 21.55 | ng/ul  | 99       |
| 86) Benzo(k)fluoranthene       | 22.86 | 252  | 913853   | 21.87 | ng/ul  | 99       |
| 88) Benzo(a)pyrene             | 23.39 | 252  | 910015   | 21.63 | ng/ul  | 98       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.71 | 276  | 1103413  | 21.60 | ng/ul  | 99       |
| 90) Dibenzo(a,h)anthracene     | 25.73 | 278  | 926741   | 21.61 | ng/ul  | 98       |
| 91) Benzo(g,h,i)perylene       | 26.40 | 276  | 918888   | 21.51 | ng/ul  | 98       |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

