

**Instrument :**  
BNA\_M  
**LabSampleId :**  
SSTD02066

mohammad  
11/27/2018 11:52:41 AM

[illegible]

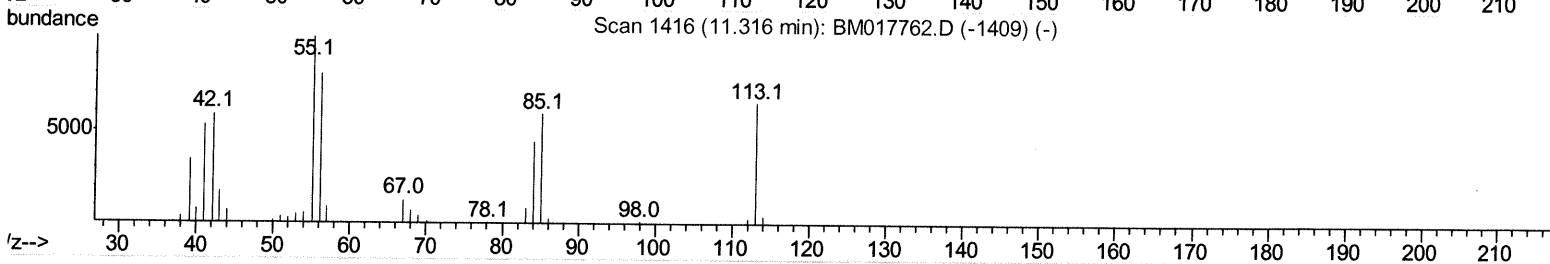
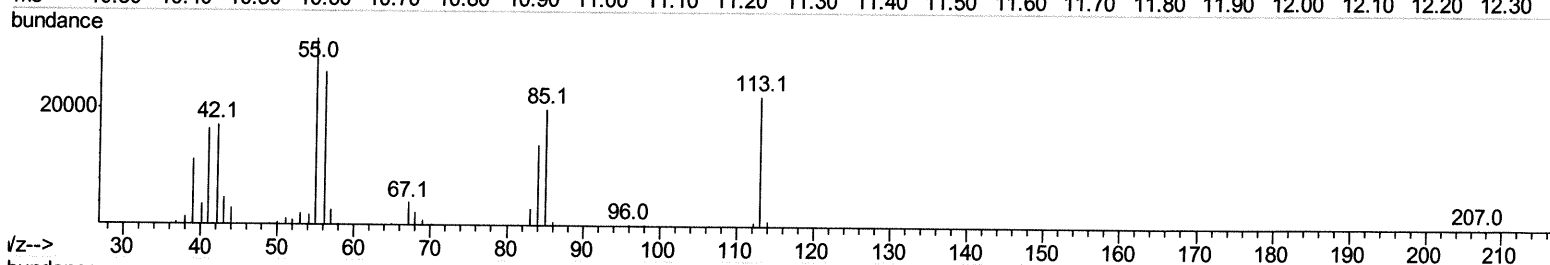
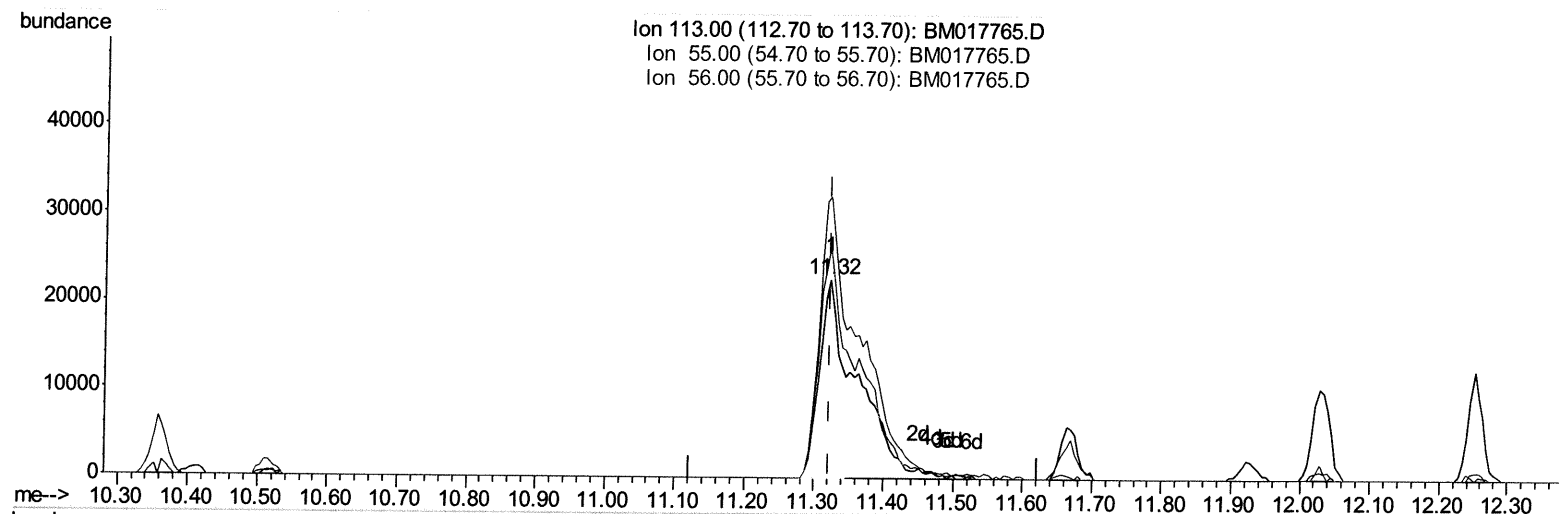
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112618\  
 Data File : BM017765.D  
 Acq On : 26 Nov 2018 12:30  
 Operator : MJ/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

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Manual Integrations  
 APPROVED

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Quant Time: Nov 26 14:50:25 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM111918MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 23 21:57:24 2018  
 Response via : Initial Calibration



TIC: BM017765.D

(32) Caprolactam

11.322min (-0.000) 10.59ng/ul

response 47395

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 138.30 | 141.84 |
| 56.00  | 110.30 | 116.52 |
| 0.00   | 0.00   | 0.00   |

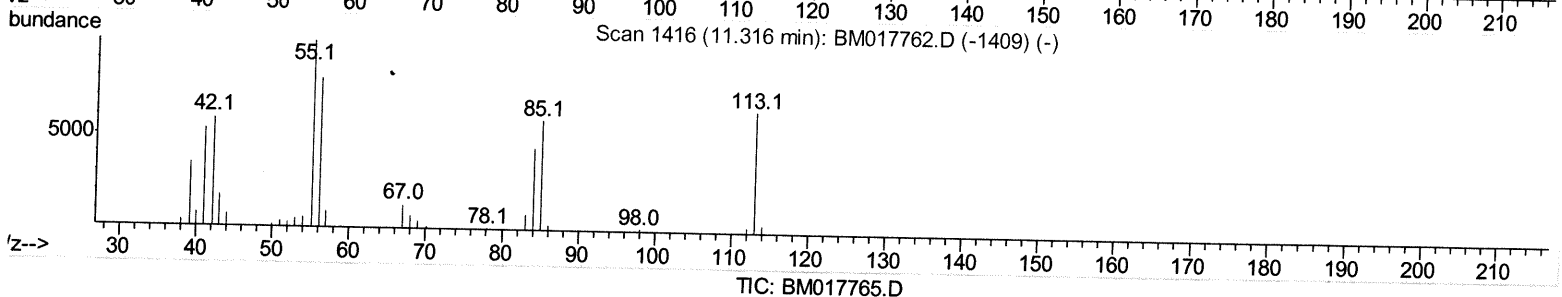
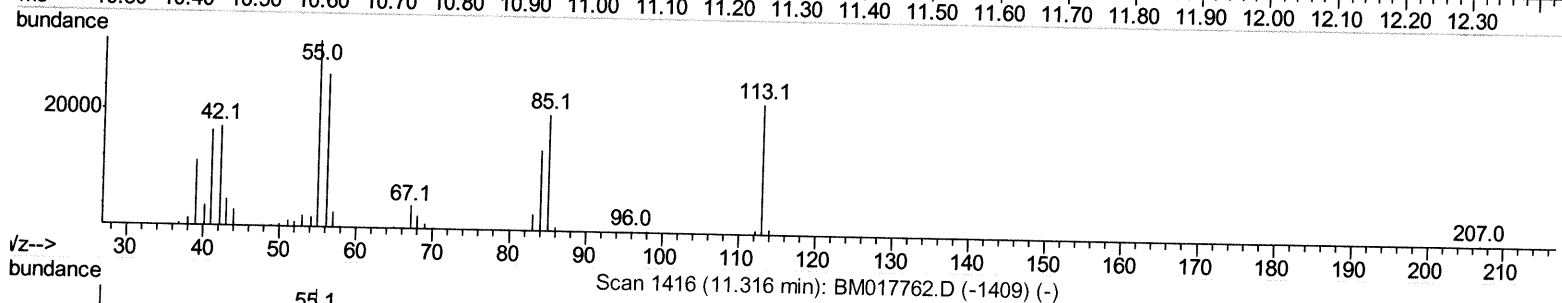
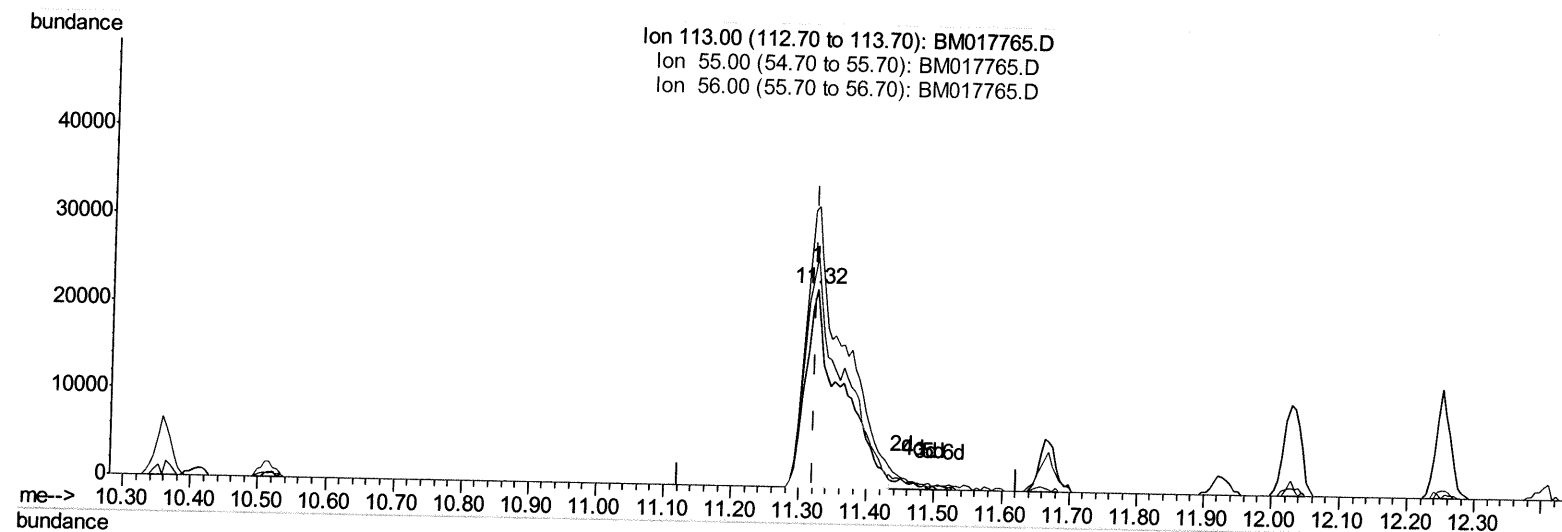
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(32) Caprolactam

11.322min (-0.000) 18.61ng/ul

response 83313

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 113.00 | 100    | 100    |
| 55.00  | 138.30 | 141.84 |
| 56.00  | 110.30 | 116.52 |
| 0.00   | 0.00   | 0.00   |

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Manual Integrations  
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Quant Time: Nov 26 15:41:20 2018  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Nov 23 21:57:24 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.59  | 152  | 186943   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.36 | 136  | 855313   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.24 | 164  | 553851   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.99 | 188  | 1298233  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.20 | 240  | 1538001  | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.38 | 264  | 1522349  | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.18  | 96  | 31396   | 7.69  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.77  | 99  | 335033  | 19.62 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.94  | 67  | 200583  | 19.48 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.13  | 132 | 265135  | 19.85 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.30  | 113 | 277565  | 19.90 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.75  | 128 | 133773  | 20.36 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.46  | 143 | 152795  | 20.30 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.99  | 165 | 289357  | 20.19 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.52 | 131 | 254537  | 20.70 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.66 | 166 | 928001  | 19.31 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.93 | 160 | 1144158 | 19.67 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.46 | 143 | 156003  | 17.85 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.24 | 176 | 816339  | 19.67 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.37 | 200 | 171361  | 18.46 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.09 | 188 | 1287863 | 19.93 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.39 | 212 | 1447299 | 19.45 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.24 | 264 | 1625604 | 19.69 | ng/ul | 0.00 |

#### Target Compounds

|                                |       |     |        | Ovalue |           |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.22  | 88  | 33299  | 7.641  | ng/uL 96  |
| 4) Benzaldehyde                | 6.76  | 77  | 64957  | 20.403 | ng/ul 96  |
| 6) Phenol                      | 6.80  | 94  | 336422 | 19.741 | ng/ul 99  |
| 8) Bis(2-Chloroethyl)ether     | 7.03  | 93  | 255244 | 19.604 | ng/ul 99  |
| 10) 2-Chlorophenol             | 7.16  | 128 | 267892 | 20.071 | ng/ul 98  |
| 11) 2-Methylphenol             | 8.04  | 108 | 257989 | 19.842 | ng/ul 100 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.12  | 45  | 297157 | 19.345 | ng/ul 100 |
| 14) Acetophenone               | 8.42  | 105 | 438989 | 20.189 | ng/ul 99  |
| 15) N-Nitroso-di-n-propylamine | 8.40  | 70  | 229730 | 20.257 | ng/ul 98  |
| 16) 4-Methylphenol             | 8.36  | 108 | 279918 | 19.922 | ng/ul 97  |
| 17) Hexachloroethane           | 8.65  | 117 | 108211 | 19.996 | ng/ul 96  |
| 20) Nitrobenzene               | 8.79  | 77  | 337893 | 20.491 | ng/ul 99  |
| 21) Isophorone                 | 9.31  | 82  | 619158 | 20.184 | ng/ul 99  |
| 23) 2-Nitrophenol              | 9.49  | 139 | 160356 | 20.469 | ng/ul 95  |
| 24) 2,4-Dimethylphenol         | 9.56  | 107 | 339949 | 20.773 | ng/ul 95  |
| 25) Bis(2-Chloroethoxy)methane | 9.79  | 93  | 374995 | 19.977 | ng/ul 99  |
| 27) 2,4-Dichlorophenol         | 10.02 | 162 | 281504 | 20.805 | ng/ul 99  |
| 28) Naphthalene                | 10.41 | 128 | 885591 | 19.873 | ng/ul 100 |
| 30) 4-Chloroaniline            | 10.54 | 127 | 258372 | 20.593 | ng/ul 97  |
| 31) Hexachlorobutadiene        | 10.69 | 225 | 181206 | 20.165 | ng/ul 98  |
| 32) Caprolactam                | 11.32 | 113 | 83313m | 18.611 | ng/ul 98  |
| 33) 4-Chloro-3-methylphenol    | 11.67 | 107 | 316622 | 20.604 | ng/ul 98  |
| 34) 2-Methylnaphthalene        | 12.03 | 142 | 675835 | 20.315 | ng/ul 99  |

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| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.40 | 216  | 359991   | 19.480 | ng/ul | 97       |
| 37) Hexachlorocyclopentadiene  | 12.37 | 237  | 213400   | 17.906 | ng/ul | 98       |
| 38) 2,4,6-Trichlorophenol      | 12.66 | 196  | 234638   | 19.536 | ng/ul | 99       |
| 39) 2,4,5-Trichlorophenol      | 12.73 | 196  | 252501   | 19.657 | ng/ul | 99       |
| 40) 1,1'-Biphenyl              | 13.06 | 154  | 878330   | 19.265 | ng/ul | 100      |
| 41) 2-Chloronaphthalene        | 13.10 | 162  | 697016   | 19.544 | ng/ul | 98       |
| 42) 2-Nitroaniline             | 13.33 | 65   | 209589   | 20.181 | ng/ul | 98       |
| 44) Dimethylphthalate          | 13.70 | 163  | 909537   | 19.535 | ng/ul | 100      |
| 45) 2,6-Dinitrotoluene         | 13.83 | 165  | 185743   | 20.303 | ng/ul | 96       |
| 47) Acenaphthylene             | 13.96 | 152  | 1069379  | 19.688 | ng/ul | 99       |
| 48) 3-Nitroaniline             | 14.16 | 138  | 168310   | 19.285 | ng/ul | 98       |
| 49) Acenaphthene               | 14.30 | 153  | 770065   | 19.672 | ng/ul | 100      |
| 50) 2,4-Dinitrophenol          | 14.37 | 184  | 105742   | 17.047 | ng/ul | 97       |
| 52) 4-Nitrophenol              | 14.48 | 109  | 137453   | 19.127 | ng/ul | 95       |
| 53) Dibenzofuran               | 14.64 | 168  | 1096245  | 19.790 | ng/ul | 100      |
| 54) 2,4-Dinitrotoluene         | 14.63 | 165  | 279623   | 20.186 | ng/ul | 96       |
| 55) 2,3,4,6-Tetrachlorophenol  | 14.87 | 232  | 239294   | 20.204 | ng/ul | 98       |
| 56) Diethylphthalate           | 15.08 | 149  | 946374   | 19.732 | ng/ul | 100      |
| 58) Fluorene                   | 15.29 | 166  | 915794   | 19.825 | ng/ul | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.29 | 204  | 466602   | 19.664 | ng/ul | 99       |
| 60) 4-Nitroaniline             | 15.34 | 138  | 177141   | 18.257 | ng/ul | 99       |
| 63) 4,6-Dinitro-2-methylphenol | 15.39 | 198  | 177895   | 18.599 | ng/ul | 97       |
| 64) N-Nitrosodiphenylamine     | 15.51 | 169  | 797405   | 19.950 | ng/ul | 98       |
| 65) 4-Bromophenyl-phenylether  | 16.19 | 248  | 293118   | 19.759 | ng/ul | 97       |
| 66) Hexachlorobenzene          | 16.29 | 284  | 322908   | 19.583 | ng/ul | 96       |
| 67) Atrazine                   | 16.47 | 200  | 289018   | 19.494 | ng/ul | 99       |
| 68) Pentachlorophenol          | 16.64 | 266  | 188737   | 18.398 | ng/ul | 98       |
| 69) Phenanthrene               | 17.03 | 178  | 1474571  | 19.862 | ng/ul | 99       |
| 71) Anthracene                 | 17.13 | 178  | 1520351  | 20.024 | ng/ul | 99       |
| 72) 1,2,3,4-Tetrachlorobenzene | 13.02 | 216  | 355855   | 19.342 | ng/uL | 100      |
| 73) Pentachlorobenzene         | 14.56 | 250  | 367852   | 19.549 | ng/uL | 99       |
| 74) Carbazole                  | 17.41 | 167  | 1326106  | 19.822 | ng/ul | 100      |
| 75) Di-n-butylphthalate        | 17.97 | 149  | 1663309  | 19.673 | ng/ul | 100      |
| 76) Fluoranthene               | 19.06 | 202  | 1770248  | 20.272 | ng/ul | 99       |
| 79) Pyrene                     | 19.43 | 202  | 1832600  | 19.629 | ng/ul | 100      |
| 80) Butylbenzylphthalate       | 20.34 | 149  | 787863   | 19.735 | ng/ul | 99       |
| 81) 3,3'-Dichlorobenzidine     | 21.13 | 252  | 607815   | 19.198 | ng/ul | 98       |
| 82) Benzo(a)anthracene         | 21.18 | 228  | 1892960  | 19.843 | ng/ul | 100      |
| 83) Bis(2-ethylhexyl)phthalate | 21.12 | 149  | 1193484  | 20.189 | ng/ul | 100      |
| 84) Chrysene                   | 21.23 | 228  | 1792635  | 20.093 | ng/ul | 99       |
| 86) Di-n-octyl phthalate       | 21.99 | 149  | 2024842  | 18.990 | ng/ul | 100      |
| 87) Benzo(b)fluoranthene       | 22.73 | 252  | 1850112  | 19.551 | ng/ul | 99       |
| 88) Benzo(k)fluoranthene       | 22.77 | 252  | 1807824  | 19.657 | ng/ul | 99       |
| 90) Benzo(a)pyrene             | 23.29 | 252  | 1756587  | 19.575 | ng/ul | 99       |
| 91) Indeno(1,2,3-cd)pyrene     | 25.56 | 276  | 1837211  | 19.322 | ng/ul | 100      |
| 92) Dibenzo(a,h)anthracene     | 25.57 | 278  | 1524433  | 18.989 | ng/ul | 99       |
| 93) Benzo(a,h,i)perylene       | 26.22 | 276  | 1495333  | 19.459 | ng/ul | 99       |

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