

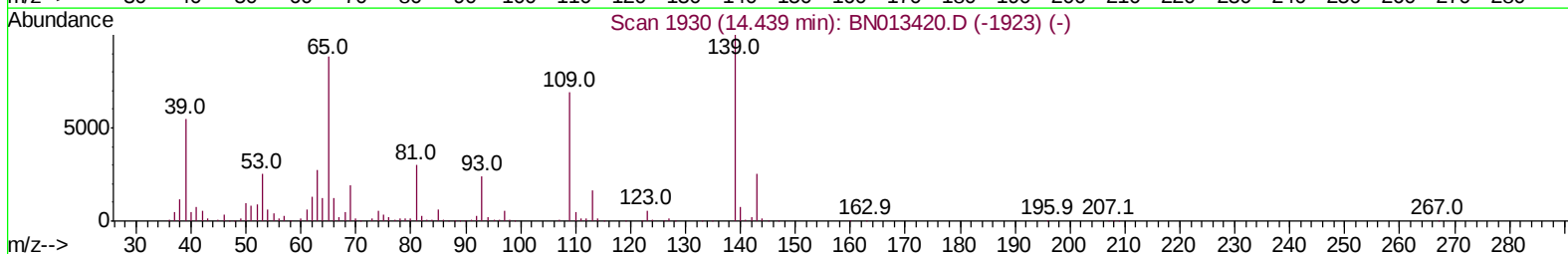
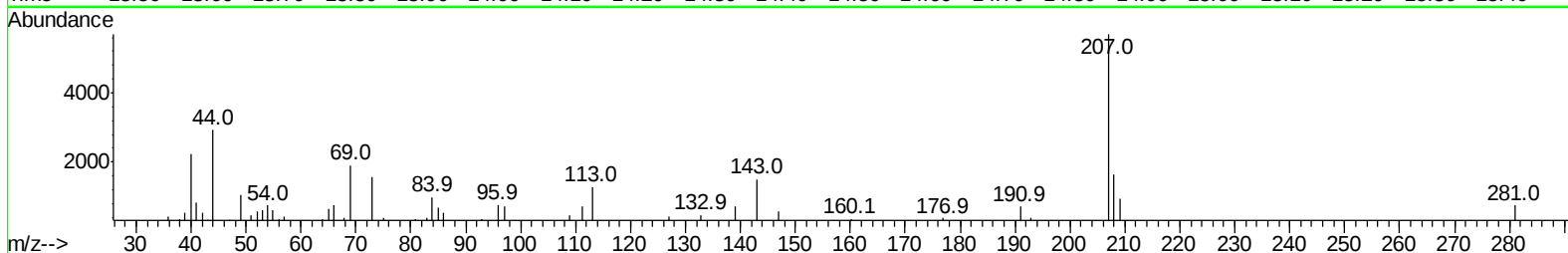
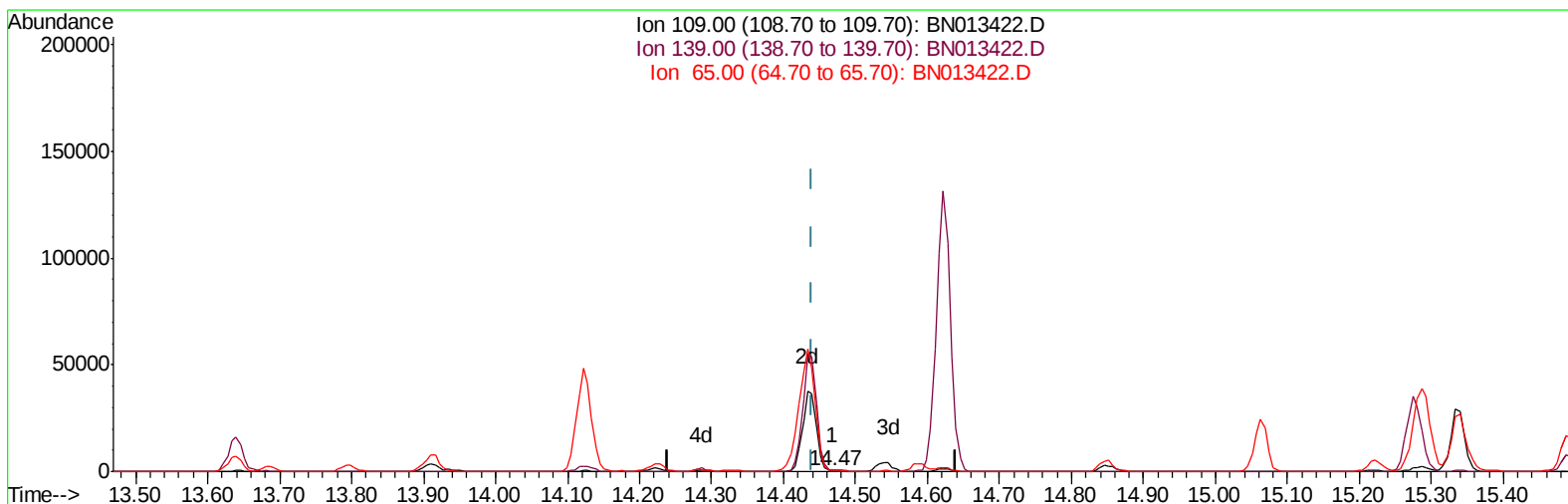
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN011921\  
Data File : BN013422.D  
Acq On : 19 Jan 2021 11:59  
Operator : CG/JU  
Sample : L5214-01  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
MDL-WATER-QT1-2021-01

Quant Time: Jan 19 13:12:23 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SFAM-EPA-BN011521.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Jan 19 10:08:46 2021  
Response via : Initial Calibration

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TIC: BN013422.D

(55) 4-Nitrophenol

14.469min (+0.030) 0.01ng/ul

response 83

| Ion    | Exp%   | Act%   |
|--------|--------|--------|
| 109.00 | 100    | 100    |
| 139.00 | 150.10 | 158.00 |
| 65.00  | 126.70 | 142.44 |
| 0.00   | 0.00   | 0.00   |

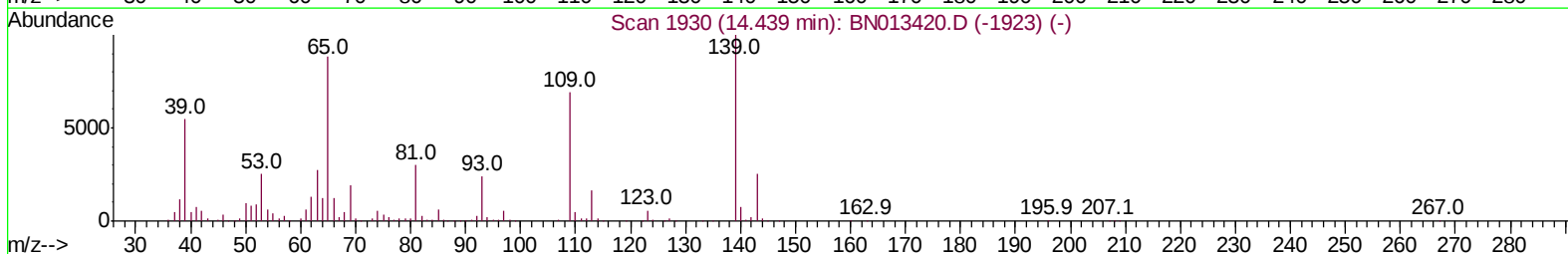
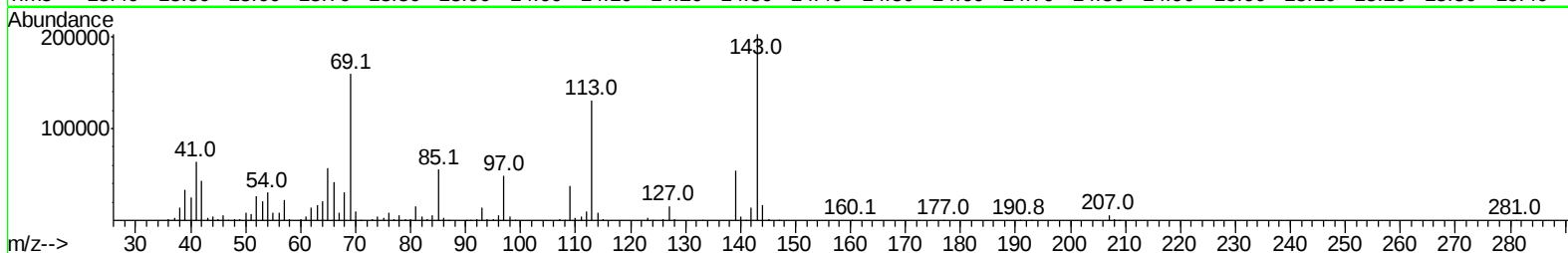
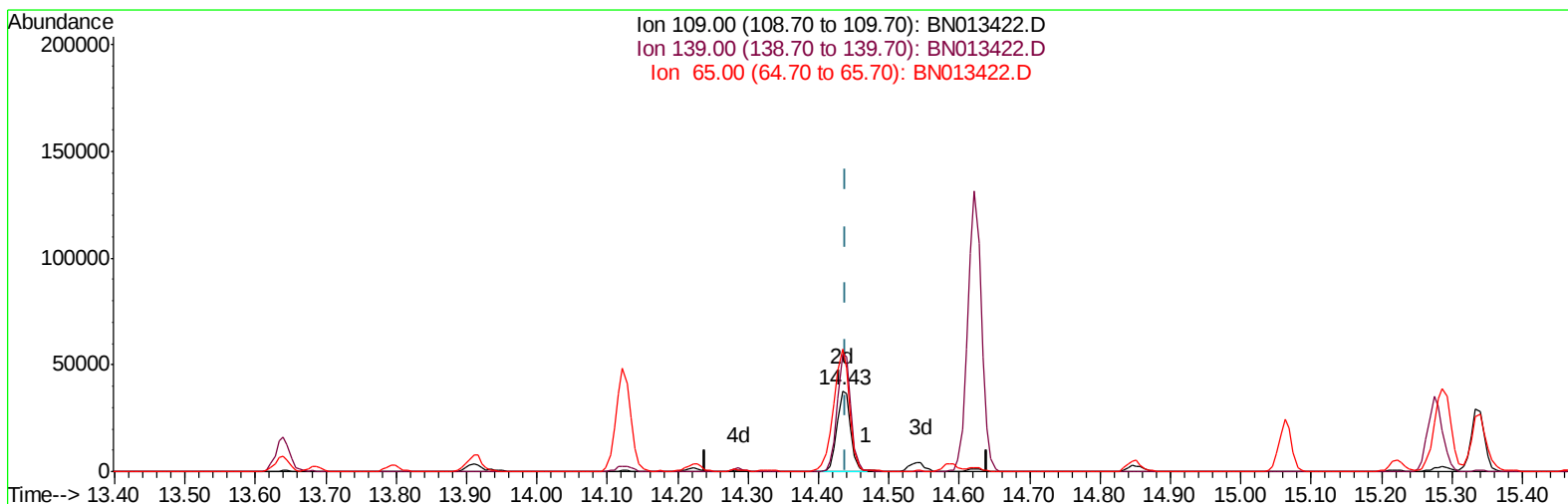
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TIC: BN013422.D

(55) 4-Nitrophenol

14.434min (-0.006) 4.05ng/ul m

response 50278

| Ion    | Exp%   | Act%    |
|--------|--------|---------|
| 109.00 | 100    | 100     |
| 139.00 | 150.10 | 144.85  |
| 65.00  | 126.70 | 152.36# |
| 0.00   | 0.00   | 0.00    |

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 19 10:08:46 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.58  | 152  | 517123   | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.35 | 136  | 2108353  | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.22 | 164  | 1200457  | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 16.97 | 188  | 2336774  | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.17 | 240  | 2039755  | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.36 | 264  | 1695061  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.18  | 96  | 46309   | 3.57  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.56  | 84  | 650117  | 19.31 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 6.76  | 99  | 1517239 | 36.48 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 6.93  | 67  | 941729  | 38.57 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.12  | 132 | 1336581 | 37.41 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 8.28  | 113 | 1245878 | 37.29 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 8.73  | 128 | 623930  | 38.21 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 9.44  | 143 | 646395  | 37.96 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 9.97  | 165 | 1171862 | 35.64 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 10.49 | 131 | 1730369 | 36.44 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 13.64 | 166 | 3300190 | 36.53 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 13.91 | 160 | 4452859 | 39.32 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 14.42 | 143 | 580771  | 33.59 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.22 | 176 | 2871844 | 36.20 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 15.34 | 200 | 445111  | 32.27 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.07 | 188 | 4200941 | 37.00 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 19.37 | 212 | 4334687 | 37.11 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 23.23 | 264 | 3332423 | 37.09 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue    |
|--------------------------------|-------|-----|--------|-------|-----------|
| 2) 1,4-Dioxane                 | 3.22  | 88  | 16975  | 1.266 | ng/uL 91  |
| 5) Pyridine                    | 3.59  | 79  | 136895 | 4.025 | ng/ul# 72 |
| 6) Benzaldehyde                | 6.73  | 77  | 85758  | 5.464 | ng/ul 97  |
| 8) Phenol                      | 6.79  | 94  | 180919 | 4.220 | ng/ul 100 |
| 10) Bis(2-Chloroethyl)ether    | 7.02  | 93  | 153192 | 4.360 | ng/ul 98  |
| 12) 2-Chlorophenol             | 7.15  | 128 | 157207 | 4.258 | ng/ul 98  |
| 13) 2-Methylphenol             | 8.02  | 108 | 122118 | 3.738 | ng/ul 99  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.11  | 45  | 213239 | 4.348 | ng/ul 98  |
| 16) Acetophenone               | 8.39  | 105 | 212271 | 4.188 | ng/ul 99  |
| 17) N-Nitroso-di-n-propylamine | 8.39  | 70  | 89864  | 3.925 | ng/ul 98  |
| 18) 4-Methylphenol             | 8.34  | 108 | 145328 | 4.132 | ng/ul 99  |
| 19) Hexachloroethane           | 8.65  | 117 | 60212  | 3.974 | ng/ul 99  |
| 22) Nitrobenzene               | 8.76  | 77  | 162113 | 4.376 | ng/ul 97  |
| 23) Isophorone                 | 9.29  | 82  | 230230 | 3.507 | ng/ul 98  |
| 25) 2-Nitrophenol              | 9.47  | 139 | 82263  | 4.307 | ng/ul 99  |
| 26) 2,4-Dimethylphenol         | 9.53  | 107 | 150377 | 4.043 | ng/ul 99  |
| 27) Bis(2-Chloroethoxy)methane | 9.77  | 93  | 182359 | 4.035 | ng/ul 99  |
| 29) 2,4-Dichlorophenol         | 10.00 | 162 | 137443 | 4.178 | ng/ul 96  |
| 30) Naphthalene                | 10.40 | 128 | 506326 | 4.230 | ng/ul 99  |
| 32) 4-Chloroaniline            | 10.50 | 127 | 195753 | 4.274 | ng/ul 98  |
| 33) Hexachlorobutadiene        | 10.69 | 225 | 89586  | 4.168 | ng/ul 93  |
| 34) Caprolactam                | 11.29 | 113 | 25403  | 2.659 | ng/ul 96  |

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Manual Integrations  
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Quant Time: Jan 19 13:14:01 2021  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 19 10:08:46 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 4-Chloro-3-methylphenol    | 11.63 | 107  | 117893   | 3.594 | ng/ul  | 97       |
| 36) 2-Methylnaphthalene        | 12.02 | 142  | 337161   | 4.185 | ng/ul  | 96       |
| 37) 1-Methylnaphthalene        | 12.24 | 142  | 328344   | 4.229 | ng/ul  | 96       |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.39 | 216  | 161505   | 4.194 | ng/ul  | 94       |
| 40) Hexachlorocyclopentadiene  | 12.37 | 237  | 88876    | 3.739 | ng/ul  | 95       |
| 41) 2,4,6-Trichlorophenol      | 12.64 | 196  | 81648    | 3.493 | ng/ul  | 100      |
| 42) 2,4,5-Trichlorophenol      | 12.70 | 196  | 95234    | 3.743 | ng/ul  | 98       |
| 43) 1,1'-Biphenyl              | 13.05 | 154  | 427739   | 4.232 | ng/ul  | 99       |
| 44) 2-Chloronaphthalene        | 13.08 | 162  | 332137   | 4.168 | ng/ul  | 99       |
| 45) 2-Nitroaniline             | 13.29 | 65   | 59077    | 3.299 | ng/ul  | 99       |
| 47) Dimethylphthalate          | 13.69 | 163  | 383264   | 4.114 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene         | 13.80 | 165  | 59133    | 3.313 | ng/ul  | 96       |
| 50) Acenaphthylene             | 13.94 | 152  | 492846   | 4.212 | ng/ul  | 99       |
| 51) 3-Nitroaniline             | 14.12 | 138  | 54449    | 3.287 | ng/ul  | 95       |
| 52) Acenaphthene               | 14.29 | 153  | 348416   | 4.174 | ng/ul  | 100      |
| 53) 2,4-Dinitrophenol          | 14.33 | 184  | 20622    | 2.064 | ng/ul  | 93       |
| 55) 4-Nitrophenol              | 14.43 | 109  | 50278m   | 4.054 | ng/ul  |          |
| 56) Dibenzofuran               | 14.62 | 168  | 485198   | 4.289 | ng/ul  | 98       |
| 57) 2,4-Dinitrotoluene         | 14.59 | 165  | 91697    | 3.614 | ng/ul  | 93       |
| 58) 2,3,4,6-Tetrachlorophenol  | 14.85 | 232  | 69360    | 3.378 | ng/ul  | 100      |
| 59) Diethylphthalate           | 15.06 | 149  | 346696   | 3.817 | ng/ul  | 100      |
| 61) Fluorene                   | 15.27 | 166  | 391494   | 4.342 | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenylether | 15.27 | 204  | 189685   | 4.309 | ng/ul  | 98       |
| 63) 4-Nitroaniline             | 15.29 | 138  | 53131    | 3.605 | ng/ul  | 94       |
| 66) 4,6-Dinitro-2-methylphenol | 15.35 | 198  | 55227    | 3.694 | ng/ul# | 75       |
| 67) N-Nitrosodiphenylamine     | 15.49 | 169  | 314535   | 4.158 | ng/ul  | 96       |
| 68) 4-Bromophenyl-phenylether  | 16.17 | 248  | 103078   | 3.906 | ng/ul  | 97       |
| 69) Hexachlorobenzene          | 16.28 | 284  | 123379   | 4.135 | ng/ul  | 97       |
| 70) Atrazine                   | 16.45 | 200  | 85942    | 3.403 | ng/ul  | 97       |
| 71) Pentachlorophenol          | 16.62 | 266  | 43371    | 2.570 | ng/ul  | 96       |
| 72) Phenanthrene               | 17.02 | 178  | 599798   | 4.269 | ng/ul  | 99       |
| 74) Anthracene                 | 17.10 | 178  | 608388   | 4.295 | ng/ul  | 98       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.00 | 216  | 158804   | 4.210 | ng/uL  | 95       |
| 76) Pentachlorobenzene         | 14.54 | 250  | 153336   | 4.176 | ng/uL  | 97       |
| 77) Carbazole                  | 17.37 | 167  | 478126   | 4.106 | ng/ul  | 99       |
| 78) Di-n-butylphthalate        | 17.96 | 149  | 478651   | 3.441 | ng/ul  | 100      |
| 80) Fluoranthene               | 19.04 | 202  | 613662   | 4.042 | ng/ul  | 98       |
| 82) Pyrene                     | 19.40 | 202  | 681147   | 4.353 | ng/ul  | 100      |
| 83) Butylbenzylphthalate       | 20.33 | 149  | 154278   | 2.808 | ng/ul  | 94       |
| 84) 3,3'-Dichlorobenzidine     | 21.10 | 252  | 98409    | 2.574 | ng/ul  | 97       |
| 85) Benzo(a)anthracene         | 21.16 | 228  | 554871   | 4.086 | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.12 | 149  | 219655   | 2.774 | ng/ul  | 100      |
| 87) Chrysene                   | 21.22 | 228  | 573540   | 4.286 | ng/ul  | 97       |
| 89) Di-n-octyl phthalate       | 21.99 | 149  | 299640   | 2.708 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 22.71 | 252  | 498631   | 4.359 | ng/ul  | 100      |
| 91) Benzo(k)fluoranthene       | 22.75 | 252  | 473995   | 4.212 | ng/ul  | 100      |
| 93) Benzo(a)pyrene             | 23.27 | 252  | 432315   | 4.228 | ng/ul  | 96       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.54 | 276  | 459117   | 3.977 | ng/ul  | 98       |
| 95) Dibenzo(a,h)anthracene     | 25.55 | 278  | 385014   | 3.990 | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene       | 26.20 | 276  | 376778   | 3.890 | ng/ul  | 98       |

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| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

|       |  |  |  |  |  |  |
|-------|--|--|--|--|--|--|
| ----- |  |  |  |  |  |  |
| ----- |  |  |  |  |  |  |

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

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