

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM070621\  
 Data File : BM030819.D  
 Acq On : 06 Jul 2021 13:19  
 Operator : CG/JU  
 Sample : SSTD08006  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080106

Quant Time: Jul 06 13:48:48 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM070621.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 06 13:25:17 2021  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.76  | 152  | 92145    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 10.55 | 136  | 398477   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.40 | 164  | 253763   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.14 | 188  | 473797   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.31 | 240  | 385349   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 23.56 | 264  | 365028   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |       |
|--------------------------------|-------|-----|---------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.25  | 96  | 74073   | 32.59 | ng/uL | 0.00  |
| 4) Pyridine-d5                 | 3.66  | 84  | 530621  | 86.89 | ng/ul | -0.01 |
| 7) Phenol-d5                   | 6.94  | 99  | 670110  | 84.32 | ng/ul | 0.00  |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.10  | 67  | 419927  | 84.07 | ng/ul | 0.00  |
| 11) 2-Chlorophenol-d4          | 7.30  | 132 | 542132  | 88.38 | ng/ul | 0.00  |
| 15) 4-Methylphenol-d8          | 8.48  | 113 | 541146  | 87.50 | ng/ul | 0.01  |
| 21) Nitrobenzene-d5            | 8.93  | 128 | 265261  | 89.21 | ng/ul | 0.00  |
| 24) 2-Nitrophenol-d4           | 9.65  | 143 | 305205  | 92.36 | ng/ul | 0.00  |
| 28) 2,4-Dichlorophenol-d3      | 10.17 | 165 | 562183  | 88.81 | ng/ul | 0.00  |
| 31) 4-Chloroaniline-d4         | 10.69 | 131 | 733259  | 83.06 | ng/ul | 0.00  |
| 46) Dimethylphthalate-d6       | 13.82 | 166 | 1511482 | 86.26 | ng/ul | 0.00  |
| 49) Acenaphthylene-d8          | 14.09 | 160 | 1981433 | 87.12 | ng/ul | 0.00  |
| 54) 4-Nitrophenol-d4           | 14.61 | 143 | 273109  | 82.27 | ng/ul | 0.01  |
| 60) Fluorene-d10               | 15.39 | 176 | 1326438 | 85.35 | ng/ul | 0.00  |
| 65) 4,6-Dinitro-2-methylphenol | 15.52 | 200 | 274508  | 88.56 | ng/ul | 0.00  |
| 73) Anthracene-d10             | 17.24 | 188 | 1898521 | 85.86 | ng/ul | 0.00  |
| 81) Pyrene-d10                 | 19.53 | 212 | 1816919 | 89.98 | ng/ul | 0.00  |
| 92) Benzo(a)pyrene-d12         | 23.42 | 264 | 1627945 | 85.78 | ng/ul | 0.01  |

## Target Compounds

|                                |       |     |         |        | Ovalue    |
|--------------------------------|-------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.29  | 88  | 76592   | 31.685 | ng/uL 98  |
| 5) Pyridine                    | 3.68  | 79  | 556935  | 84.724 | ng/ul 98  |
| 6) Benzaldehyde                | 6.92  | 77  | 317312  | 82.119 | ng/ul 99  |
| 8) Phenol                      | 6.97  | 94  | 682113  | 84.888 | ng/ul 99  |
| 10) Bis(2-Chloroethyl)ether    | 7.20  | 93  | 524576  | 82.831 | ng/ul 100 |
| 12) 2-Chlorophenol             | 7.33  | 128 | 541118  | 87.538 | ng/ul 99  |
| 13) 2-Methylphenol             | 8.21  | 108 | 503553  | 86.479 | ng/ul 99  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.30  | 45  | 858802  | 85.548 | ng/ul 99  |
| 16) Acetophenone               | 8.59  | 105 | 841652  | 88.129 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 8.59  | 70  | 451883  | 95.659 | ng/ul 97  |
| 18) 4-Methylphenol             | 8.55  | 108 | 549069  | 86.587 | ng/ul 97  |
| 19) Hexachloroethane           | 8.84  | 117 | 228282  | 88.137 | ng/ul 94  |
| 22) Nitrobenzene               | 8.97  | 77  | 655894  | 87.439 | ng/ul 97  |
| 23) Isophorone                 | 9.50  | 82  | 1179362 | 90.792 | ng/ul 99  |
| 25) 2-Nitrophenol              | 9.67  | 139 | 316034  | 90.071 | ng/ul 97  |
| 26) 2,4-Dimethylphenol         | 9.73  | 107 | 620590  | 87.508 | ng/ul 99  |
| 27) Bis(2-Chloroethoxy)methane | 9.97  | 93  | 718881  | 85.063 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 10.20 | 162 | 534917  | 88.206 | ng/ul 99  |
| 30) Naphthalene                | 10.60 | 128 | 1680316 | 83.286 | ng/ul 99  |
| 32) 4-Chloroaniline            | 10.72 | 127 | 731590  | 83.478 | ng/ul 100 |
| 33) Hexachlorobutadiene        | 10.88 | 225 | 353688  | 86.253 | ng/ul 98  |
| 34) Caprolactam                | 11.52 | 113 | 82955   | 47.128 | ng/ul 97  |

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 Sample : SSTD08006  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080106

Quant Time: Jul 06 13:48:48 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM070621.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 06 13:25:17 2021  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|---------|--------|----------|
| 35) 4-Chloro-3-methylphenol    | 11.85 | 107  | 546459   | 91.769  | ng/ul  | 100      |
| 36) 2-Methylnaphthalene        | 12.22 | 142  | 1224335  | 86.962  | ng/ul  | 100      |
| 37) 1-Methylnaphthalene        | 12.43 | 142  | 1226476  | 86.063  | ng/ul  | 100      |
| 39) 1,2,4,5-Tetrachlorobenzene | 12.59 | 216  | 665144   | 84.014  | ng/ul  | 99       |
| 40) Hexachlorocyclopentadiene  | 12.56 | 237  | 414640   | 80.260  | ng/ul  | 98       |
| 41) 2,4,6-Trichlorophenol      | 12.83 | 196  | 440323   | 88.311  | ng/ul  | 97       |
| 42) 2,4,5-Trichlorophenol      | 12.90 | 196  | 468317   | 88.007  | ng/ul  | 99       |
| 43) 1,1'-Biphenyl              | 13.23 | 154  | 1570205  | 83.840  | ng/ul  | 99       |
| 44) 2-Chloronaphthalene        | 13.27 | 162  | 1218211  | 82.737  | ng/ul  | 99       |
| 45) 2-Nitroaniline             | 13.48 | 65   | 382692   | 96.713  | ng/ul  | 99       |
| 47) Dimethylphthalate          | 13.86 | 163  | 1468604  | 85.544  | ng/ul  | 100      |
| 48) 2,6-Dinitrotoluene         | 13.99 | 165  | 314762   | 95.157  | ng/ul  | 95       |
| 50) Acenaphthylene             | 14.12 | 152  | 1800827  | 85.359  | ng/ul  | 99       |
| 51) 3-Nitroaniline             | 14.32 | 138  | 301206   | 85.480  | ng/ul  | 94       |
| 52) Acenaphthene               | 14.46 | 153  | 1291432  | 84.348  | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol          | 14.52 | 184  | 196831   | 90.517  | ng/ul  | 99       |
| 55) 4-Nitrophenol              | 14.63 | 109  | 229385   | 86.375  | ng/ul  | 96       |
| 56) Dibenzofuran               | 14.80 | 168  | 1790052  | 83.865  | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene         | 14.77 | 165  | 428505   | 92.520  | ng/ul  | 98       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.03 | 232  | 385666   | 89.530  | ng/ul  | 99       |
| 59) Diethylphthalate           | 15.22 | 149  | 1471879  | 89.249  | ng/ul  | 99       |
| 61) Fluorene                   | 15.44 | 166  | 1535534  | 87.292  | ng/ul  | 98       |
| 62) 4-Chlorophenyl-phenylether | 15.44 | 204  | 777015   | 88.696  | ng/ul  | 99       |
| 63) 4-Nitroaniline             | 15.48 | 138  | 282566   | 84.210  | ng/ul  | 98       |
| 66) 4,6-Dinitro-2-methylphenol | 15.54 | 198  | 261749   | 86.770  | ng/ul# | 97       |
| 67) N-Nitrosodiphenylamine     | 15.66 | 169  | 1241496  | 87.919  | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether  | 16.33 | 248  | 445941   | 92.242  | ng/ul  | 96       |
| 69) Hexachlorobenzene          | 16.45 | 284  | 496627   | 88.811  | ng/ul  | 96       |
| 70) Atrazine                   | 16.61 | 200  | 420945   | 84.046  | ng/ul  | 100      |
| 71) Pentachlorophenol          | 16.79 | 266  | 304565   | 86.917  | ng/ul  | 98       |
| 72) Phenanthrene               | 17.18 | 178  | 2135569  | 83.933  | ng/ul  | 99       |
| 74) Anthracene                 | 17.27 | 178  | 2223185  | 85.457  | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.20 | 216  | 649894   | 86.000  | ng/uL  | 99       |
| 76) Pentachlorobenzene         | 14.72 | 250  | 632900   | 87.921  | ng/uL  | 99       |
| 77) Carbazole                  | 17.54 | 167  | 1846098  | 79.136  | ng/ul  | 100      |
| 78) Di-n-butylphthalate        | 18.10 | 149  | 2274564  | 95.688  | ng/ul  | 99       |
| 80) Fluoranthene               | 19.19 | 202  | 2254150  | 91.408  | ng/ul  | 99       |
| 82) Pyrene                     | 19.56 | 202  | 2321890  | 89.608  | ng/ul  | 99       |
| 83) Butylbenzylphthalate       | 20.45 | 149  | 856347   | 100.400 | ng/ul  | 94       |
| 84) 3,3'-Dichlorobenzidine     | 21.23 | 252  | 541750   | 71.321  | ng/ul  | 99       |
| 85) Benzo(a)anthracene         | 21.30 | 228  | 2008791  | 84.000  | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.22 | 149  | 1271990  | 103.685 | ng/ul  | 100      |
| 87) Chrysene                   | 21.34 | 228  | 1945344  | 83.444  | ng/ul  | 100      |
| 89) Di-n-octyl phthalate       | 22.10 | 149  | 1967820  | 93.713  | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 22.89 | 252  | 1965728  | 84.821  | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene       | 22.93 | 252  | 1934910  | 86.877  | ng/ul  | 99       |
| 93) Benzo(a)pyrene             | 23.47 | 252  | 1781232  | 85.479  | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene     | 25.84 | 276  | 2312284  | 88.143  | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene     | 25.85 | 278  | 1951980  | 89.758  | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene       | 26.53 | 276  | 1711670  | 78.204  | ng/ul  | 99       |

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Data File : BM030819.D  
Acq On : 06 Jul 2021 13:19  
Operator : CG/JU  
Sample : SSTD08006  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD080106

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM070621.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Jul 06 13:25:17 2021  
Response via : Initial Calibration

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

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

