

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051320\  
 Data File : BM025978.D  
 Acq On : 13 May 2020 13:00  
 Operator : MA/SJ  
 Sample : SSTD04055  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD04055

Quant Time: May 13 13:45:32 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 13 13:00:56 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.50  | 152  | 122379   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.27 | 136  | 490453   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.15 | 164  | 285856   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.90 | 188  | 600994   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.10 | 240  | 645004   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.22 | 264  | 665444   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.08  | 96  | 58278   | 19.98 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.70  | 99  | 439928  | 45.55 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.86  | 67  | 303660  | 55.08 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.05  | 132 | 331774  | 42.32 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.22  | 113 | 334670  | 42.17 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.65  | 128 | 159164  | 43.66 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.36  | 143 | 163044  | 42.07 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.90  | 165 | 303676  | 38.32 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.41 | 131 | 370937  | 40.31 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.57 | 166 | 855728  | 39.39 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.84 | 160 | 1042970 | 40.07 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.37 | 143 | 168056  | 40.63 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.15 | 176 | 746611  | 39.15 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.28 | 200 | 144444  | 39.39 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.00 | 188 | 1123543 | 40.02 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.30 | 212 | 1246801 | 40.24 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.09 | 264 | 1379317 | 40.49 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |        | Ovalue    |
|--------------------------------|-------|-----|---------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.11  | 88  | 61275   | 19.509 | ng/uL 90  |
| 4) Benzaldehyde                | 6.66  | 77  | 312451  | 61.079 | ng/ul 96  |
| 6) Phenol                      | 6.72  | 94  | 491260  | 48.707 | ng/ul 99  |
| 8) Bis(2-Chloroethyl)ether     | 6.95  | 93  | 380778  | 47.673 | ng/ul 99  |
| 10) 2-Chlorophenol             | 7.08  | 128 | 357069  | 43.272 | ng/ul 99  |
| 11) 2-Methylphenol             | 7.96  | 108 | 339431  | 43.982 | ng/ul 98  |
| 12) 2,2'-oxybis(1-Chloropropan | 8.04  | 45  | 684715  | 64.249 | ng/ul 99  |
| 14) Acetophenone               | 8.32  | 105 | 561708  | 45.887 | ng/ul 99  |
| 15) N-Nitroso-di-n-propylamine | 8.32  | 70  | 315390  | 51.430 | ng/ul 96  |
| 16) 4-Methylphenol             | 8.28  | 108 | 371303  | 44.072 | ng/ul 99  |
| 17) Hexachloroethane           | 8.57  | 117 | 151130  | 44.055 | ng/ul 98  |
| 20) Nitrobenzene               | 8.69  | 77  | 465701  | 52.439 | ng/ul 98  |
| 21) Isophorone                 | 9.22  | 82  | 794139  | 48.423 | ng/ul 100 |
| 23) 2-Nitrophenol              | 9.40  | 139 | 185645  | 42.654 | ng/ul 93  |
| 24) 2,4-Dimethylphenol         | 9.47  | 107 | 416396  | 47.179 | ng/ul 99  |
| 25) Bis(2-Chloroethoxy)methane | 9.70  | 93  | 481683  | 45.825 | ng/ul 99  |
| 27) 2,4-Dichlorophenol         | 9.93  | 162 | 316347  | 40.234 | ng/ul 96  |
| 28) Naphthalene                | 10.32 | 128 | 1106277 | 41.417 | ng/ul 99  |
| 30) 4-Chloroaniline            | 10.43 | 127 | 382842  | 41.002 | ng/ul 99  |
| 31) Hexachlorobutadiene        | 10.62 | 225 | 203257  | 39.568 | ng/ul 98  |
| 32) Caprolactam                | 11.20 | 113 | 57803   | 22.231 | ng/ul 91  |
| 33) 4-Chloro-3-methylphenol    | 11.57 | 107 | 372626  | 45.588 | ng/ul 99  |
| 34) 2-Methylnaphthalene        | 11.94 | 142 | 754918  | 39.569 | ng/ul 99  |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051320\  
 Data File : BM025978.D  
 Acq On : 13 May 2020 13:00  
 Operator : MA/SJ  
 Sample : SSTD04055  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD04055

Quant Time: May 13 13:45:32 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed May 13 13:00:56 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units | Dev(Min) |
|--------------------------------|-------|------|----------|--------|-------|----------|
| 35) 1-Methylnaphthalene        | 12.16 | 142  | 695618   | 36.563 | ng/ul | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.32 | 216  | 378863   | 41.905 | ng/ul | 99       |
| 38) Hexachlorocyclopentadiene  | 12.30 | 237  | 239576   | 38.829 | ng/ul | 99       |
| 39) 2,4,6-Trichlorophenol      | 12.57 | 196  | 234141   | 40.736 | ng/ul | 97       |
| 40) 2,4,5-Trichlorophenol      | 12.64 | 196  | 258733   | 41.354 | ng/ul | 97       |
| 41) 1,1'-Biphenyl              | 12.97 | 154  | 945764   | 41.407 | ng/ul | 100      |
| 42) 2-Chloronaphthalene        | 13.01 | 162  | 729675   | 40.561 | ng/ul | 99       |
| 43) 2-Nitroaniline             | 13.22 | 65   | 270954   | 60.265 | ng/ul | 97       |
| 45) Dimethylphthalate          | 13.62 | 163  | 900102   | 39.802 | ng/ul | 99       |
| 46) 2,6-Dinitrotoluene         | 13.73 | 165  | 183154   | 41.177 | ng/ul | 95       |
| 48) Acenaphthylene             | 13.87 | 152  | 1124827  | 39.901 | ng/ul | 100      |
| 49) 3-Nitroaniline             | 14.06 | 138  | 179204   | 44.658 | ng/ul | 98       |
| 50) Acenaphthene               | 14.22 | 153  | 773855   | 40.382 | ng/ul | 98       |
| 51) 2,4-Dinitrophenol          | 14.27 | 184  | 99111    | 37.142 | ng/ul | 95       |
| 53) 4-Nitrophenol              | 14.38 | 109  | 147822   | 45.719 | ng/ul | 97       |
| 54) Dibenzofuran               | 14.55 | 168  | 1085640  | 39.935 | ng/ul | 99       |
| 55) 2,4-Dinitrotoluene         | 14.53 | 165  | 267498   | 41.134 | ng/ul | 94       |
| 56) 2,3,4,6-Tetrachlorophenol  | 14.79 | 232  | 219132   | 38.643 | ng/ul | 99       |
| 57) Diethylphthalate           | 15.00 | 149  | 907344   | 39.144 | ng/ul | 99       |
| 59) Fluorene                   | 15.20 | 166  | 883051   | 38.616 | ng/ul | 100      |
| 60) 4-Chlorophenyl-phenylether | 15.21 | 204  | 436626   | 37.307 | ng/ul | 98       |
| 61) 4-Nitroaniline             | 15.23 | 138  | 214399   | 44.901 | ng/ul | 97       |
| 64) 4,6-Dinitro-2-methylphenol | 15.29 | 198  | 151250   | 37.719 | ng/ul | 97       |
| 65) N-Nitrosodiphenylamine     | 15.42 | 169  | 757677   | 40.797 | ng/ul | 99       |
| 66) 4-Bromophenyl-phenylether  | 16.10 | 248  | 261438   | 35.274 | ng/ul | 99       |
| 67) Hexachlorobenzene          | 16.22 | 284  | 290193   | 31.889 | ng/ul | 100      |
| 68) Atrazine                   | 16.39 | 200  | 268087   | 38.274 | ng/ul | 99       |
| 69) Pentachlorophenol          | 16.56 | 266  | 179848   | 33.838 | ng/ul | 98       |
| 70) Phenanthrene               | 16.94 | 178  | 1398065  | 39.521 | ng/ul | 99       |
| 72) Anthracene                 | 17.03 | 178  | 1418576  | 39.348 | ng/ul | 100      |
| 73) 1,2,3,4-Tetrachlorobenzene | 12.94 | 216  | 375422   | 40.766 | ng/uL | 100      |
| 74) Pentachlorobenzene         | 14.48 | 250  | 351282   | 34.268 | ng/uL | 99       |
| 75) Carbazole                  | 17.30 | 167  | 1272894  | 40.400 | ng/ul | 99       |
| 76) Di-n-butylphthalate        | 17.89 | 149  | 1522808  | 41.396 | ng/ul | 99       |
| 77) Fluoranthene               | 18.96 | 202  | 1605516  | 38.503 | ng/ul | 98       |
| 80) Pvrene                     | 19.33 | 202  | 1683291  | 38.858 | ng/ul | 99       |
| 81) Butylbenzylphthalate       | 20.26 | 149  | 672654   | 40.771 | ng/ul | 95       |
| 82) 3,3'-Dichlorobenzidine     | 21.02 | 252  | 527011   | 35.959 | ng/ul | 99       |
| 83) Benzo(a)anthracene         | 21.08 | 228  | 1673046  | 39.343 | ng/ul | 100      |
| 84) Bis(2-ethylhexyl)phthalate | 21.04 | 149  | 1043612  | 40.779 | ng/ul | 100      |
| 85) Chrysene                   | 21.13 | 228  | 1619741  | 38.822 | ng/ul | 99       |
| 87) Di-n-octyl phthalate       | 21.90 | 149  | 1748448  | 43.849 | ng/ul | 100      |
| 88) Benzo(b)fluoranthene       | 22.60 | 252  | 1711800  | 41.228 | ng/ul | 99       |
| 89) Benzo(k)fluoranthene       | 22.64 | 252  | 1691045  | 41.344 | ng/ul | 100      |
| 91) Benzo(a)pyrene             | 23.13 | 252  | 1512420  | 40.027 | ng/ul | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 25.31 | 276  | 1952871  | 39.136 | ng/ul | 99       |
| 93) Dibenzo(a,h)anthracene     | 25.33 | 278  | 1642505  | 38.871 | ng/ul | 100      |
| 94) Benzo(a,h,i)perylene       | 25.95 | 276  | 1624448  | 39.500 | ng/ul | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM051320\  
Data File : BM025978.D  
Acq On : 13 May 2020 13:00  
Operator : MA/SJ  
Sample : SSTD04055  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD04055

Quant Time: May 13 13:45:32 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed May 13 13:00:56 2020  
Response via : Initial Calibration

| Internal Standards                                                         | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|----------------------------------------------------------------------------|------|------|----------|------|-------|----------|
| -----                                                                      |      |      |          |      |       |          |
| (#) = qualifier out of range (m) = manual integration (+) = signals summed |      |      |          |      |       |          |

