

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
 Data File : BM027833.D  
 Acq On : 18 Nov 2020 16:30  
 Operator : JU/CG  
 Sample : SSTD08006  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080006

Quant Time: Nov 18 17:06:13 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM111820.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 18 16:04:33 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.34  | 152  | 51287    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 11.18 | 136  | 216107   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.96 | 164  | 130439   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.67 | 188  | 255405   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.78 | 240  | 168833   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 24.20 | 264  | 149031   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |        |       |      |
|--------------------------------|-------|-----|---------|--------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.49  | 96  | 37587   | 28.91  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.96  | 84  | 301927  | 83.61  | ng/ul | 0.00 |
| 7) Phenol-d5                   | 7.48  | 99  | 389072  | 92.29  | ng/ul | 0.01 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.65  | 67  | 247357  | 102.19 | ng/ul | 0.01 |
| 11) 2-Chlorophenol-d4          | 7.86  | 132 | 290010  | 89.09  | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 9.05  | 113 | 309755  | 91.52  | ng/ul | 0.01 |
| 21) Nitrobenzene-d5            | 9.52  | 128 | 144769  | 84.42  | ng/ul | 0.01 |
| 24) 2-Nitrophenol-d4           | 10.25 | 143 | 164667  | 89.05  | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.79 | 165 | 306983  | 80.65  | ng/ul | 0.01 |
| 31) 4-Chloroaniline-d4         | 11.30 | 131 | 392947  | 82.78  | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 14.36 | 166 | 832564  | 79.40  | ng/ul | 0.01 |
| 49) Acenaphthylene-d8          | 14.66 | 160 | 1036310 | 81.38  | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 15.13 | 143 | 146540  | 79.22  | ng/ul | 0.02 |
| 60) Fluorene-d10               | 15.94 | 176 | 717303  | 74.53  | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 16.05 | 200 | 140914  | 88.02  | ng/ul | 0.01 |
| 73) Anthracene-d10             | 17.77 | 188 | 1012453 | 79.50  | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 20.02 | 212 | 935526  | 98.25  | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 24.06 | 264 | 647847  | 76.91  | ng/ul | 0.01 |

## Target Compounds

|                                |       |     |        |         | Ovalue   |
|--------------------------------|-------|-----|--------|---------|----------|
| 2) 1,4-Dioxane                 | 3.53  | 88  | 40312  | 27.809  | ng/uL 97 |
| 5) Pyridine                    | 3.98  | 79  | 301329 | 83.949  | ng/ul 95 |
| 6) Benzaldehyde                | 7.46  | 77  | 189477 | 85.007  | ng/ul 98 |
| 8) Phenol                      | 7.50  | 94  | 393468 | 93.652  | ng/ul 98 |
| 10) Bis(2-Chloroethyl)ether    | 7.75  | 93  | 317497 | 98.368  | ng/ul 96 |
| 12) 2-Chlorophenol             | 7.90  | 128 | 297466 | 87.546  | ng/ul 99 |
| 13) 2-Methylphenol             | 8.78  | 108 | 298856 | 91.672  | ng/ul 98 |
| 14) 2,2'-oxybis(1-Chloropropan | 8.88  | 45  | 505764 | 120.423 | ng/ul 98 |
| 16) Acetophenone               | 9.18  | 105 | 498656 | 89.015  | ng/ul 98 |
| 17) N-Nitroso-di-n-propylamine | 9.17  | 70  | 280448 | 102.674 | ng/ul 99 |
| 18) 4-Methylphenol             | 9.12  | 108 | 317200 | 93.631  | ng/ul 97 |
| 19) Hexachloroethane           | 9.45  | 117 | 144362 | 85.487  | ng/ul 97 |
| 22) Nitrobenzene               | 9.57  | 77  | 431817 | 82.747  | ng/ul 98 |
| 23) Isophorone                 | 10.10 | 82  | 765639 | 92.795  | ng/ul 99 |
| 25) 2-Nitrophenol              | 10.28 | 139 | 172292 | 87.779  | ng/ul 99 |
| 26) 2,4-Dimethylphenol         | 10.33 | 107 | 382239 | 81.176  | ng/ul 99 |
| 27) Bis(2-Chloroethoxy)methane | 10.57 | 93  | 439050 | 95.630  | ng/ul 99 |
| 29) 2,4-Dichlorophenol         | 10.82 | 162 | 295699 | 80.282  | ng/ul 99 |
| 30) Naphthalene                | 11.23 | 128 | 957552 | 78.884  | ng/ul 99 |
| 32) 4-Chloroaniline            | 11.33 | 127 | 383167 | 82.790  | ng/ul 99 |
| 33) Hexachlorobutadiene        | 11.51 | 225 | 214043 | 70.153  | ng/ul 98 |
| 34) Caprolactam                | 12.13 | 113 | 93942  | 91.831  | ng/ul 93 |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
 Data File : BM027833.D  
 Acq On : 18 Nov 2020 16:30  
 Operator : JU/CG  
 Sample : SSTD08006  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD080006

Quant Time: Nov 18 17:06:13 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM111820.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Nov 18 16:04:33 2020  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc    | Units | Dev(Min) |
|--------------------------------|-------|------|----------|---------|-------|----------|
| 35) 4-Chloro-3-methylphenol    | 12.43 | 107  | 344019   | 91.466  | ng/ul | 97       |
| 36) 2-Methylnaphthalene        | 12.82 | 142  | 672149   | 85.643  | ng/ul | 97       |
| 37) 1-Methylnaphthalene        | 13.03 | 142  | 644366   | 82.079  | ng/ul | 100      |
| 39) 1,2,4,5-Tetrachlorobenzene | 13.17 | 216  | 367323   | 76.114  | ng/ul | 98       |
| 40) Hexachlorocyclopentadiene  | 13.15 | 237  | 258996   | 74.803  | ng/ul | 97       |
| 41) 2,4,6-Trichlorophenol      | 13.40 | 196  | 241894   | 81.218  | ng/ul | 98       |
| 42) 2,4,5-Trichlorophenol      | 13.47 | 196  | 257606   | 80.240  | ng/ul | 98       |
| 43) 1,1'-Biphenyl              | 13.80 | 154  | 874415   | 79.450  | ng/ul | 99       |
| 44) 2-Chloronaphthalene        | 13.85 | 162  | 695957   | 79.684  | ng/ul | 99       |
| 45) 2-Nitroaniline             | 14.04 | 65   | 256693   | 89.312  | ng/ul | 98       |
| 47) Dimethylphthalate          | 14.40 | 163  | 856542   | 79.436  | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene         | 14.53 | 165  | 184382   | 86.524  | ng/ul | 97       |
| 50) Acenaphthylene             | 14.69 | 152  | 1015631  | 81.096  | ng/ul | 99       |
| 51) 3-Nitroaniline             | 14.85 | 138  | 160787   | 86.336  | ng/ul | 100      |
| 52) Acenaphthene               | 15.02 | 153  | 716649   | 79.960  | ng/ul | 99       |
| 53) 2,4-Dinitrophenol          | 15.05 | 184  | 107657   | 80.032  | ng/ul | 96       |
| 55) 4-Nitrophenol              | 15.15 | 109  | 163235   | 56.763  | ng/ul | 95       |
| 56) Dibenzofuran               | 15.35 | 168  | 970753   | 75.285  | ng/ul | 99       |
| 57) 2,4-Dinitrotoluene         | 15.30 | 165  | 249844   | 81.212  | ng/ul | 98       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.57 | 232  | 215323   | 77.865  | ng/ul | 99       |
| 59) Diethylphthalate           | 15.75 | 149  | 880060   | 80.290  | ng/ul | 99       |
| 61) Fluorene                   | 15.99 | 166  | 809432   | 74.771  | ng/ul | 100      |
| 62) 4-Chlorophenyl-phenylether | 15.97 | 204  | 419314   | 71.161  | ng/ul | 97       |
| 63) 4-Nitroaniline             | 16.01 | 138  | 154240   | 78.778  | ng/ul | 96       |
| 66) 4,6-Dinitro-2-methylphenol | 16.06 | 198  | 148266   | 85.888  | ng/ul | 100      |
| 67) N-Nitrosodiphenylamine     | 16.19 | 169  | 671402   | 87.417  | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether  | 16.86 | 248  | 246164   | 82.363  | ng/ul | 97       |
| 69) Hexachlorobenzene          | 16.99 | 284  | 266024   | 78.972  | ng/ul | 98       |
| 70) Atrazine                   | 17.12 | 200  | 250632   | 80.856  | ng/ul | 97       |
| 71) Pentachlorophenol          | 17.32 | 266  | 165616   | 78.891  | ng/ul | 96       |
| 72) Phenanthrene               | 17.72 | 178  | 1207450  | 78.928  | ng/ul | 100      |
| 74) Anthracene                 | 17.81 | 178  | 1227273  | 80.683  | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.77 | 216  | 363842   | 87.643  | ng/uL | 98       |
| 76) Pentachlorobenzene         | 15.27 | 250  | 335887   | 84.151  | ng/uL | 97       |
| 77) Carbazole                  | 18.06 | 167  | 1001320  | 77.529  | ng/ul | 99       |
| 78) Di-n-butylphthalate        | 18.60 | 149  | 1369867  | 91.703  | ng/ul | 99       |
| 80) Fluoranthene               | 19.69 | 202  | 1255457  | 103.046 | ng/ul | 99       |
| 82) Pyrene                     | 20.05 | 202  | 1236227  | 98.699  | ng/ul | 98       |
| 83) Butylbenzylphthalate       | 20.90 | 149  | 511214   | 112.759 | ng/ul | 98       |
| 84) 3,3'-Dichlorobenzidine     | 21.69 | 252  | 249703   | 68.381  | ng/ul | 100      |
| 85) Benzo(a)anthracene         | 21.76 | 228  | 949683   | 80.609  | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.66 | 149  | 686340   | 112.088 | ng/ul | 98       |
| 87) Chrysene                   | 21.82 | 228  | 911098   | 80.170  | ng/ul | 99       |
| 89) Di-n-octyl phthalate       | 22.59 | 149  | 1073404  | 105.682 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene       | 23.47 | 252  | 844096   | 72.569  | ng/ul | 99       |
| 91) Benzo(k)fluoranthene       | 23.52 | 252  | 783609   | 72.201  | ng/ul | 100      |
| 93) Benzo(a)pyrene             | 24.11 | 252  | 738887   | 79.188  | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene     | 26.74 | 276  | 851994   | 72.044  | ng/ul | 96       |
| 95) Dibenzo(a,h)anthracene     | 26.74 | 278  | 714052   | 72.557  | ng/ul | 99       |
| 96) Benzo(g,h,i)perylene       | 27.51 | 276  | 703566   | 71.739  | ng/ul | 98       |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
Data File : BM027833.D  
Acq On : 18 Nov 2020 16:30  
Operator : JU/CG  
Sample : SSTD08006  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD080006

Quant Time: Nov 18 17:06:13 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SFAM-EPA-BM111820.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Wed Nov 18 16:04:33 2020  
Response via : Initial Calibration

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

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

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

