

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
 Data File : BM027830.D  
 Acq On : 18 Nov 2020 14:41  
 Operator : JU/CG  
 Sample : SSTD01003  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010003

Quant Time: Nov 18 16:08:02 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  | 40275    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 11.17 | 136  | 172530   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.95 | 164  | 112159   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.67 | 188  | 243114   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.77 | 240  | 203965   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 24.20 | 264  | 195583   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.49  | 96  | 3738   | 3.66  | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.96  | 84  | 27108  | 9.56  | ng/ul | 0.00 |
| 7) Phenol-d5                   | 7.46  | 99  | 35874  | 10.84 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.64  | 67  | 24493  | 12.89 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.85  | 132 | 27428  | 10.73 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 9.03  | 113 | 29331  | 11.04 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 9.51  | 128 | 13513  | 9.87  | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 10.24 | 143 | 15140  | 10.26 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.78 | 165 | 28611  | 9.42  | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 11.29 | 131 | 40245  | 10.62 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 14.34 | 166 | 92939  | 10.31 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.65 | 160 | 110022 | 10.05 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 15.11 | 143 | 16142  | 10.15 | ng/ul | 0.00 |
| 60) Fluorene-d10               | 15.93 | 176 | 77807  | 9.40  | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 16.03 | 200 | 13928  | 9.14  | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.76 | 188 | 119754 | 9.88  | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 20.02 | 212 | 125099 | 10.87 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 24.04 | 264 | 104797 | 9.48  | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |       |        | Ovalue    |
|--------------------------------|-------|-----|-------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.53  | 88  | 3903  | 3.429  | ng/uL 90  |
| 5) Pyridine                    | 3.99  | 79  | 28053 | 9.952  | ng/ul 98  |
| 6) Benzaldehyde                | 7.45  | 77  | 19253 | 10.999 | ng/ul 96  |
| 8) Phenol                      | 7.49  | 94  | 37606 | 11.398 | ng/ul 96  |
| 10) Bis(2-Chloroethyl)ether    | 7.73  | 93  | 31389 | 12.384 | ng/ul 99  |
| 12) 2-Chlorophenol             | 7.89  | 128 | 28159 | 10.553 | ng/ul 96  |
| 13) 2-Methylphenol             | 8.77  | 108 | 28285 | 11.048 | ng/ul 95  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.87  | 45  | 49865 | 15.119 | ng/ul 100 |
| 16) Acetophenone               | 9.17  | 105 | 49740 | 11.307 | ng/ul 98  |
| 17) N-Nitroso-di-n-propylamine | 9.15  | 70  | 28322 | 13.204 | ng/ul 97  |
| 18) 4-Methylphenol             | 9.10  | 108 | 30955 | 11.636 | ng/ul 98  |
| 19) Hexachloroethane           | 9.45  | 117 | 13778 | 10.390 | ng/ul 92  |
| 22) Nitrobenzene               | 9.55  | 77  | 42079 | 10.100 | ng/ul 100 |
| 23) Isophorone                 | 10.08 | 82  | 78332 | 11.892 | ng/ul 99  |
| 25) 2-Nitrophenol              | 10.27 | 139 | 16375 | 10.450 | ng/ul 98  |
| 26) 2,4-Dimethylphenol         | 10.32 | 107 | 38020 | 10.114 | ng/ul 99  |
| 27) Bis(2-Chloroethoxy)methane | 10.56 | 93  | 44218 | 12.064 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 10.81 | 162 | 29018 | 9.868  | ng/ul 94  |
| 30) Naphthalene                | 11.22 | 128 | 95316 | 9.835  | ng/ul 98  |
| 32) 4-Chloroaniline            | 11.32 | 127 | 39306 | 10.638 | ng/ul 96  |
| 33) Hexachlorobutadiene        | 11.50 | 225 | 21509 | 8.830  | ng/ul 99  |
| 34) Caprolactam                | 12.07 | 113 | 10037 | 12.290 | ng/ul 93  |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
 Data File : BM027830.D  
 Acq On : 18 Nov 2020 14:41  
 Operator : JU/CG  
 Sample : SSTD01003  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD010003

Quant Time: Nov 18 16:08:02 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    | 12.42 | 107  | 35506    | 11.824 | ng/ul  | 93       |
| 36) 2-Methylnaphthalene        | 12.81 | 142  | 69716    | 11.127 | ng/ul  | 95       |
| 37) 1-Methylnaphthalene        | 13.03 | 142  | 66159    | 10.556 | ng/ul# | 100      |
| 39) 1,2,4,5-Tetrachlorobenzene | 13.17 | 216  | 37380    | 9.008  | ng/ul  | 93       |
| 40) Hexachlorocyclopentadiene  | 13.15 | 237  | 23518    | 7.900  | ng/ul  | 92       |
| 41) 2,4,6-Trichlorophenol      | 13.40 | 196  | 24517    | 9.573  | ng/ul  | 95       |
| 42) 2,4,5-Trichlorophenol      | 13.47 | 196  | 26585    | 9.630  | ng/ul  | 99       |
| 43) 1,1'-Biphenyl              | 13.79 | 154  | 90797    | 9.594  | ng/ul  | 100      |
| 44) 2-Chloronaphthalene        | 13.84 | 162  | 72503    | 9.654  | ng/ul  | 99       |
| 45) 2-Nitroaniline             | 14.03 | 65   | 25974    | 10.510 | ng/ul  | 97       |
| 47) Dimethylphthalate          | 14.39 | 163  | 96908    | 10.452 | ng/ul  | 99       |
| 48) 2,6-Dinitrotoluene         | 14.52 | 165  | 18761    | 10.239 | ng/ul  | 96       |
| 50) Acenaphthylene             | 14.68 | 152  | 108922   | 10.115 | ng/ul  | 99       |
| 51) 3-Nitroaniline             | 14.84 | 138  | 16732    | 10.449 | ng/ul  | 99       |
| 52) Acenaphthene               | 15.02 | 153  | 76742    | 9.958  | ng/ul  | 99       |
| 53) 2,4-Dinitrophenol          | 15.04 | 184  | 9300     | 8.040  | ng/ul# | 88       |
| 55) 4-Nitrophenol              | 15.12 | 109  | 18558    | 7.505  | ng/ul  | 98       |
| 56) Dibenzofuran               | 15.34 | 168  | 106549   | 9.610  | ng/ul  | 99       |
| 57) 2,4-Dinitrotoluene         | 15.29 | 165  | 26829    | 10.142 | ng/ul  | 98       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.56 | 232  | 22788    | 9.584  | ng/ul  | 99       |
| 59) Diethylphthalate           | 15.73 | 149  | 98321    | 10.432 | ng/ul  | 98       |
| 61) Fluorene                   | 15.99 | 166  | 89366    | 9.601  | ng/ul  | 99       |
| 62) 4-Chlorophenyl-phenylether | 15.97 | 204  | 45744    | 9.028  | ng/ul  | 97       |
| 63) 4-Nitroaniline             | 15.99 | 138  | 15694    | 9.322  | ng/ul  | 97       |
| 66) 4,6-Dinitro-2-methylphenol | 16.04 | 198  | 14670    | 8.928  | ng/ul  | 100      |
| 67) N-Nitrosodiphenylamine     | 16.18 | 169  | 74500    | 10.190 | ng/ul  | 99       |
| 68) 4-Bromophenyl-phenylether  | 16.86 | 248  | 27362    | 9.618  | ng/ul  | 98       |
| 69) Hexachlorobenzene          | 16.98 | 284  | 30246    | 9.433  | ng/ul  | 98       |
| 70) Atrazine                   | 17.10 | 200  | 29122    | 9.870  | ng/ul  | 98       |
| 71) Pentachlorophenol          | 17.32 | 266  | 17517    | 8.766  | ng/ul  | 98       |
| 72) Phenanthrene               | 17.71 | 178  | 143626   | 9.863  | ng/ul  | 97       |
| 74) Anthracene                 | 17.80 | 178  | 145559   | 10.053 | ng/ul  | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.77 | 216  | 37482    | 9.485  | ng/uL  | 99       |
| 76) Pentachlorobenzene         | 15.26 | 250  | 37214    | 9.795  | ng/uL  | 100      |
| 77) Carbazole                  | 18.06 | 167  | 122058   | 9.928  | ng/ul  | 100      |
| 78) Di-n-butylphthalate        | 18.59 | 149  | 167063   | 11.749 | ng/ul  | 99       |
| 80) Fluoranthene               | 19.69 | 202  | 162612   | 11.048 | ng/ul  | 99       |
| 82) Pyrene                     | 20.04 | 202  | 163272   | 10.790 | ng/ul  | 98       |
| 83) Butylbenzylphthalate       | 20.90 | 149  | 71546    | 13.063 | ng/ul  | 99       |
| 84) 3,3'-Dichlorobenzidine     | 21.68 | 252  | 46913    | 10.634 | ng/ul  | 97       |
| 85) Benzo(a)anthracene         | 21.76 | 228  | 142021   | 9.978  | ng/ul  | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.66 | 149  | 99245    | 13.416 | ng/ul  | 99       |
| 87) Chrysene                   | 21.81 | 228  | 136915   | 9.972  | ng/ul  | 99       |
| 89) Di-n-octyl phthalate       | 22.59 | 149  | 163286   | 12.250 | ng/ul  | 100      |
| 90) Benzo(b)fluoranthene       | 23.46 | 252  | 135157   | 8.854  | ng/ul  | 99       |
| 91) Benzo(k)fluoranthene       | 23.51 | 252  | 126300   | 8.867  | ng/ul  | 97       |
| 93) Benzo(a)pyrene             | 24.09 | 252  | 118985   | 9.717  | ng/ul  | 98       |
| 94) Indeno(1,2,3-cd)pyrene     | 26.70 | 276  | 137412   | 8.854  | ng/ul  | 99       |
| 95) Dibenzo(a,h)anthracene     | 26.71 | 278  | 116008   | 8.982  | ng/ul  | 99       |
| 96) Benzo(g,h,i)perylene       | 27.48 | 276  | 112748   | 8.760  | ng/ul  | 99       |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
Data File : BM027830.D  
Acq On : 18 Nov 2020 14:41  
Operator : JU/CG  
Sample : SSTD01003  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD010003

Quant Time: Nov 18 16:08:02 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

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM111820\  
Data File : BM027830.D  
Acq On : 18 Nov 2020 14:41  
Operator : JU/CG  
Sample : SSTD01003  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD010003

Quant Time: Nov 18 16:08:02 2020  
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