

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111820\  
 Data File : BM027832.D  
 Acq On : 18 Nov 2020 15:53  
 Operator : JU/CG  
 Sample : SSTD04005  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD040005

Quant Time: Nov 18 16:28:48 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  | 39804    | 20.00 | ng/ul | 0.00     |
| 20) Naphthalene-d8        | 11.17 | 136  | 174806   | 20.00 | ng/ul | 0.00     |
| 38) Acenaphthene-d10      | 14.95 | 164  | 112172   | 20.00 | ng/ul | 0.00     |
| 64) Phenanthrene-d10      | 17.67 | 188  | 241358   | 20.00 | ng/ul | 0.00     |
| 79) Chrysene-d12          | 21.78 | 240  | 206418   | 20.00 | ng/ul | 0.00     |
| 88) Perylene-d12          | 24.20 | 264  | 187626   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.49  | 96  | 14710  | 14.58 | ng/uL | 0.00 |
| 4) Pyridine-d5                 | 3.96  | 84  | 117974 | 42.09 | ng/ul | 0.00 |
| 7) Phenol-d5                   | 7.47  | 99  | 155212 | 47.44 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)ether-d  | 7.64  | 67  | 101241 | 53.89 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4          | 7.86  | 132 | 115635 | 45.77 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8          | 9.05  | 113 | 126249 | 48.06 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5            | 9.52  | 128 | 58763  | 42.36 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4           | 10.25 | 143 | 66572  | 44.51 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3      | 10.79 | 165 | 125408 | 40.73 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4         | 11.30 | 131 | 169311 | 44.10 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6       | 14.35 | 166 | 375173 | 41.61 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8          | 14.65 | 160 | 452067 | 41.28 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4           | 15.12 | 143 | 69003  | 43.38 | ng/ul | 0.01 |
| 60) Fluorene-d10               | 15.93 | 176 | 323449 | 39.08 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylphenol | 16.04 | 200 | 65168  | 43.07 | ng/ul | 0.00 |
| 73) Anthracene-d10             | 17.77 | 188 | 488051 | 40.56 | ng/ul | 0.00 |
| 81) Pyrene-d10                 | 20.02 | 212 | 510602 | 43.86 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12         | 24.05 | 264 | 419472 | 39.56 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue    |
|--------------------------------|-------|-----|--------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.53  | 88  | 15387  | 13.677 | ng/uL 95  |
| 5) Pyridine                    | 3.98  | 79  | 118386 | 42.497 | ng/ul 98  |
| 6) Benzaldehyde                | 7.45  | 77  | 79872  | 46.171 | ng/ul 96  |
| 8) Phenol                      | 7.49  | 94  | 157186 | 48.206 | ng/ul 95  |
| 10) Bis(2-Chloroethyl)ether    | 7.74  | 93  | 129481 | 51.689 | ng/ul 96  |
| 12) 2-Chlorophenol             | 7.89  | 128 | 118889 | 45.084 | ng/ul 96  |
| 13) 2-Methylphenol             | 8.78  | 108 | 119324 | 47.161 | ng/ul 95  |
| 14) 2,2'-oxybis(1-Chloropropan | 8.87  | 45  | 207858 | 63.769 | ng/ul 99  |
| 16) Acetophenone               | 9.17  | 105 | 206409 | 47.476 | ng/ul 97  |
| 17) N-Nitroso-di-n-propylamine | 9.16  | 70  | 116575 | 54.991 | ng/ul 97  |
| 18) 4-Methylphenol             | 9.11  | 108 | 130345 | 49.575 | ng/ul 97  |
| 19) Hexachloroethane           | 9.45  | 117 | 57488  | 43.864 | ng/ul 95  |
| 22) Nitrobenzene               | 9.56  | 77  | 177006 | 41.932 | ng/ul 97  |
| 23) Isophorone                 | 10.09 | 82  | 325314 | 48.744 | ng/ul 100 |
| 25) 2-Nitrophenol              | 10.28 | 139 | 70085  | 44.143 | ng/ul 97  |
| 26) 2,4-Dimethylphenol         | 10.32 | 107 | 161307 | 42.350 | ng/ul 100 |
| 27) Bis(2-Chloroethoxy)methane | 10.56 | 93  | 182466 | 49.133 | ng/ul 98  |
| 29) 2,4-Dichlorophenol         | 10.81 | 162 | 121139 | 40.660 | ng/ul 99  |
| 30) Naphthalene                | 11.23 | 128 | 395080 | 40.237 | ng/ul 98  |
| 32) 4-Chloroaniline            | 11.32 | 127 | 165954 | 44.329 | ng/ul 99  |
| 33) Hexachlorobutadiene        | 11.50 | 225 | 87709  | 35.539 | ng/ul 98  |
| 34) Caprolactam                | 12.10 | 113 | 43230  | 52.243 | ng/ul 98  |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111820\  
 Data File : BM027832.D  
 Acq On : 18 Nov 2020 15:53  
 Operator : JU/CG  
 Sample : SSTD04005  
 Misc :  
 ALS Vial : 5 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD040005

Quant Time: Nov 18 16:28:48 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  | 147210   | 48.387 | ng/ul | 97       |
| 36) 2-Methylnaphthalene        | 12.81 | 142  | 283068   | 44.589 | ng/ul | 98       |
| 37) 1-Methylnaphthalene        | 13.03 | 142  | 270532   | 42.602 | ng/ul | 99       |
| 39) 1,2,4,5-Tetrachlorobenzene | 13.17 | 216  | 155980   | 37.585 | ng/ul | 96       |
| 40) Hexachlorocyclopentadiene  | 13.15 | 237  | 105070   | 35.288 | ng/ul | 98       |
| 41) 2,4,6-Trichlorophenol      | 13.40 | 196  | 104205   | 40.685 | ng/ul | 97       |
| 42) 2,4,5-Trichlorophenol      | 13.47 | 196  | 112729   | 40.831 | ng/ul | 99       |
| 43) 1,1'-Biphenyl              | 13.80 | 154  | 371715   | 39.274 | ng/ul | 100      |
| 44) 2-Chloronaphthalene        | 13.85 | 162  | 299696   | 39.902 | ng/ul | 99       |
| 45) 2-Nitroaniline             | 14.03 | 65   | 112185   | 45.389 | ng/ul | 99       |
| 47) Dimethylphthalate          | 14.40 | 163  | 387586   | 41.798 | ng/ul | 100      |
| 48) 2,6-Dinitrotoluene         | 14.52 | 165  | 81396    | 44.417 | ng/ul | 96       |
| 50) Acenaphthylene             | 14.68 | 152  | 444135   | 41.238 | ng/ul | 99       |
| 51) 3-Nitroaniline             | 14.85 | 138  | 73202    | 45.708 | ng/ul | 95       |
| 52) Acenaphthene               | 15.02 | 153  | 313058   | 40.618 | ng/ul | 98       |
| 53) 2,4-Dinitrophenol          | 15.05 | 184  | 46383    | 40.096 | ng/ul | 96       |
| 55) 4-Nitrophenol              | 15.13 | 109  | 77568    | 31.366 | ng/ul | 96       |
| 56) Dibenzofuran               | 15.35 | 168  | 434409   | 39.176 | ng/ul | 98       |
| 57) 2,4-Dinitrotoluene         | 15.29 | 165  | 114986   | 43.463 | ng/ul | 95       |
| 58) 2,3,4,6-Tetrachlorophenol  | 15.56 | 232  | 95641    | 40.218 | ng/ul | 97       |
| 59) Diethylphthalate           | 15.75 | 149  | 406578   | 43.134 | ng/ul | 99       |
| 61) Fluorene                   | 15.99 | 166  | 364436   | 39.147 | ng/ul | 99       |
| 62) 4-Chlorophenyl-phenylether | 15.97 | 204  | 186126   | 36.731 | ng/ul | 99       |
| 63) 4-Nitroaniline             | 15.99 | 138  | 73004    | 43.359 | ng/ul | 99       |
| 66) 4,6-Dinitro-2-methylphenol | 16.05 | 198  | 69406    | 42.546 | ng/ul | 96       |
| 67) N-Nitrosodiphenylamine     | 16.18 | 169  | 306463   | 42.224 | ng/ul | 99       |
| 68) 4-Bromophenyl-phenylether  | 16.86 | 248  | 112356   | 39.780 | ng/ul | 98       |
| 69) Hexachlorobenzene          | 16.98 | 284  | 123038   | 38.651 | ng/ul | 99       |
| 70) Atrazine                   | 17.11 | 200  | 122948   | 41.973 | ng/ul | 98       |
| 71) Pentachlorophenol          | 17.32 | 266  | 77846    | 39.240 | ng/ul | 98       |
| 72) Phenanthrene               | 17.72 | 178  | 577194   | 39.926 | ng/ul | 99       |
| 74) Anthracene                 | 17.80 | 178  | 590154   | 41.056 | ng/ul | 99       |
| 75) 1,2,3,4-Tetrachlorobenzene | 13.77 | 216  | 155818   | 39.718 | ng/uL | 98       |
| 76) Pentachlorobenzene         | 15.27 | 250  | 147074   | 38.991 | ng/uL | 98       |
| 77) Carbazole                  | 18.06 | 167  | 506885   | 41.530 | ng/ul | 99       |
| 78) Di-n-butylphthalate        | 18.59 | 149  | 694940   | 49.229 | ng/ul | 100      |
| 80) Fluoranthene               | 19.69 | 202  | 672340   | 45.136 | ng/ul | 99       |
| 82) Pyrene                     | 20.05 | 202  | 682313   | 44.556 | ng/ul | 99       |
| 83) Butylbenzylphthalate       | 20.90 | 149  | 302834   | 54.634 | ng/ul | 99       |
| 84) 3,3'-Dichlorobenzidine     | 21.68 | 252  | 174720   | 39.135 | ng/ul | 99       |
| 85) Benzo(a)anthracene         | 21.76 | 228  | 585734   | 40.664 | ng/ul | 99       |
| 86) Bis(2-ethylhexyl)phthalate | 21.66 | 149  | 423550   | 56.576 | ng/ul | 99       |
| 87) Chrysene                   | 21.82 | 228  | 564996   | 40.663 | ng/ul | 99       |
| 89) Di-n-octyl phthalate       | 22.59 | 149  | 698799   | 54.648 | ng/ul | 100      |
| 90) Benzo(b)fluoranthene       | 23.46 | 252  | 547860   | 37.412 | ng/ul | 98       |
| 91) Benzo(k)fluoranthene       | 23.52 | 252  | 512213   | 37.487 | ng/ul | 98       |
| 93) Benzo(a)pyrene             | 24.10 | 252  | 474738   | 40.413 | ng/ul | 99       |
| 94) Indeno(1,2,3-cd)pyrene     | 26.72 | 276  | 502955   | 33.781 | ng/ul | 96       |
| 95) Dibenzo(a,h)anthracene     | 26.73 | 278  | 425940   | 34.378 | ng/ul | 98       |
| 96) Benzo(g,h,i)perylene       | 27.49 | 276  | 404674   | 32.775 | ng/ul | 99       |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM111820\  
Data File : BM027832.D  
Acq On : 18 Nov 2020 15:53  
Operator : JU/CG  
Sample : SSTD04005  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD040005

Quant Time: Nov 18 16:28:48 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 : BM027832.D  
Acq On : 18 Nov 2020 15:53  
Operator : JU/CG  
Sample : SSTD04005  
Misc :  
ALS Vial : 5 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
SSTD040005

Quant Time: Nov 18 16:28:48 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

