

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012918\  
 Data File : BM013998.D  
 Acq On : 30 Jan 2018 15:34  
 Operator : SJ/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 SSTD02009

Manual Integrations  
 APPROVED

Sohil  
 1/30/2018 6:08:22 PM

Quant Time: Jan 30 16:29:40 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 30 07:49:34 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.57  | 152  | 73045    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.34 | 136  | 282477   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.22 | 164  | 195933   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.97 | 188  | 514982   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.18 | 240  | 635387   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.37 | 264  | 631172   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.13  | 96  | 12095  | 6.54  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.76  | 99  | 100541 | 17.93 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.92  | 67  | 63313  | 18.63 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.10  | 132 | 87441  | 19.14 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.29  | 113 | 83113  | 19.28 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.72  | 128 | 42774  | 19.51 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 46048  | 19.88 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.98  | 165 | 94547  | 21.13 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.50 | 131 | 101381 | 26.12 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.64 | 166 | 335127 | 20.75 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.91 | 160 | 428113 | 20.79 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.46 | 143 | 43062  | 16.41 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.22 | 176 | 320154 | 22.54 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.37 | 200 | 56582  | 18.25 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.07 | 188 | 517180 | 20.62 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.38 | 212 | 621253 | 21.00 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.23 | 264 | 608625 | 20.98 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |        | Ovalue |    |
|--------------------------------|-------|-----|--------|--------|--------|----|
| 2) 1,4-Dioxane                 | 3.16  | 88  | 14264  | 7.046  | ng/uL# | 66 |
| 4) Benzaldehyde                | 6.73  | 77  | 80182  | 20.662 | ng/ul  | 91 |
| 6) Phenol                      | 6.79  | 94  | 100376 | 17.847 | ng/ul  | 93 |
| 8) Bis(2-Chloroethyl)ether     | 7.00  | 93  | 77177  | 17.536 | ng/ul  | 97 |
| 10) 2-Chlorophenol             | 7.14  | 128 | 85765  | 18.833 | ng/ul  | 99 |
| 11) 2-Methylphenol             | 8.02  | 108 | 78173  | 18.855 | ng/ul  | 87 |
| 12) 2,2'-oxybis(1-Chloropropan | 8.10  | 45  | 74861  | 16.717 | ng/ul# | 71 |
| 14) Acetophenone               | 8.39  | 105 | 141699 | 19.864 | ng/ul  | 96 |
| 15) N-Nitroso-di-n-propylamine | 8.37  | 70  | 71445  | 20.529 | ng/ul# | 74 |
| 16) 4-Methylphenol             | 8.35  | 108 | 86544  | 19.522 | ng/ul  | 99 |
| 17) Hexachloroethane           | 8.62  | 117 | 46176  | 20.831 | ng/ul  | 85 |
| 20) Nitrobenzene               | 8.77  | 77  | 119625 | 21.127 | ng/ul  | 99 |
| 21) Isophorone                 | 9.29  | 82  | 189952 | 20.763 | ng/ul# | 91 |
| 23) 2-Nitrophenol              | 9.47  | 139 | 51495  | 21.334 | ng/ul  | 96 |
| 24) 2,4-Dimethylphenol         | 9.53  | 107 | 111300 | 21.871 | ng/ul  | 96 |
| 25) Bis(2-Chloroethoxy)methane | 9.77  | 93  | 112610 | 20.191 | ng/ul  | 95 |
| 27) 2,4-Dichlorophenol         | 10.00 | 162 | 87262  | 20.237 | ng/ul# | 85 |
| 28) Naphthalene                | 10.39 | 128 | 296833 | 21.122 | ng/ul  | 99 |
| 30) 4-Chloroaniline            | 10.52 | 127 | 97675  | 25.375 | ng/ul  | 90 |
| 31) Hexachlorobutadiene        | 10.67 | 225 | 86400  | 24.410 | ng/ul  | 96 |
| 32) Caprolactam                | 11.30 | 113 | 25745m | 21.462 | ng/ul  |    |
| 33) 4-Chloro-3-methylphenol    | 11.66 | 107 | 96062  | 21.976 | ng/ul  | 92 |
| 34) 2-Methylnaphthalene        | 12.02 | 142 | 222776 | 21.262 | ng/ul  | 99 |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012918\  
 Data File : BM013998.D  
 Acq On : 30 Jan 2018 15:34  
 Operator : SJ/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02009

Manual Integrations  
 APPROVED

Sohil  
 1/30/2018 6:08:22 PM

Quant Time: Jan 30 16:29:40 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jan 30 07:49:34 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.39 | 216  | 150459   | 20.724 | ng/ul  | 98       |
| 37) Hexachlorocyclopentadiene  | 12.36 | 237  | 43598    | 14.465 | ng/ul  | 94       |
| 38) 2,4,6-Trichlorophenol      | 12.65 | 196  | 89140    | 21.212 | ng/ul  | 98       |
| 39) 2,4,5-Trichlorophenol      | 12.73 | 196  | 90690    | 20.740 | ng/ul  | 93       |
| 40) 1,1'-Biphenyl              | 13.05 | 154  | 314650   | 19.293 | ng/ul  | 97       |
| 41) 2-Chloronaphthalene        | 13.09 | 162  | 254571   | 19.968 | ng/ul  | 97       |
| 42) 2-Nitroaniline             | 13.32 | 65   | 76617    | 21.786 | ng/ul  | 95       |
| 44) Dimethylphthalate          | 13.69 | 163  | 323293   | 20.362 | ng/ul  | 99       |
| 45) 2,6-Dinitrotoluene         | 13.82 | 165  | 63769    | 20.726 | ng/ul# | 86       |
| 47) Acenaphthylene             | 13.94 | 152  | 383401   | 20.349 | ng/ul  | 99       |
| 48) 3-Nitroaniline             | 14.16 | 138  | 47481    | 19.622 | ng/ul  | 96       |
| 49) Acenaphthene               | 14.29 | 153  | 269075   | 20.620 | ng/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 14.38 | 184  | 22586    | 12.806 | ng/ul# | 84       |
| 52) 4-Nitrophenol              | 14.47 | 109  | 51388    | 19.276 | ng/ul# | 73       |
| 53) Dibenzofuran               | 14.63 | 168  | 413922   | 21.086 | ng/ul  | 99       |
| 54) 2,4-Dinitrotoluene         | 14.62 | 165  | 99244    | 22.127 | ng/ul# | 68       |
| 55) 2,3,4,6-Tetrachlorophenol  | 14.86 | 232  | 99061    | 24.085 | ng/ul  | 96       |
| 56) Diethylphthalate           | 15.06 | 149  | 352126   | 22.038 | ng/ul  | 94       |
| 58) Fluorene                   | 15.27 | 166  | 343139   | 21.339 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.27 | 204  | 197874   | 22.773 | ng/ul  | 98       |
| 60) 4-Nitroaniline             | 15.33 | 138  | 55772    | 19.699 | ng/ul# | 75       |
| 63) 4,6-Dinitro-2-methylphenol | 15.38 | 198  | 56774    | 17.352 | ng/ul  | 98       |
| 64) N-Nitrosodiphenylamine     | 15.50 | 169  | 304550   | 20.107 | ng/ul  | 97       |
| 65) 4-Bromophenyl-phenylether  | 16.17 | 248  | 124292   | 20.435 | ng/ul  | 92       |
| 66) Hexachlorobenzene          | 16.28 | 284  | 136930   | 20.914 | ng/ul# | 91       |
| 67) Atrazine                   | 16.46 | 200  | 132360   | 21.896 | ng/ul  | 97       |
| 68) Pentachlorophenol          | 16.63 | 266  | 60135    | 17.284 | ng/ul  | 91       |
| 69) Phenanthrene               | 17.02 | 178  | 588516   | 20.559 | ng/ul  | 99       |
| 71) Anthracene                 | 17.11 | 178  | 600914   | 20.352 | ng/ul  | 98       |
| 72) Carbazole                  | 17.39 | 167  | 489393   | 20.080 | ng/ul  | 98       |
| 73) Di-n-butylphthalate        | 17.95 | 149  | 599146   | 20.902 | ng/ul  | 98       |
| 74) Fluoranthene               | 19.04 | 202  | 741025   | 21.686 | ng/ul# | 97       |
| 77) Pyrene                     | 19.41 | 202  | 781382   | 20.229 | ng/ul# | 96       |
| 78) Butylbenzylphthalate       | 20.33 | 149  | 284587   | 20.915 | ng/ul  | 95       |
| 79) 3,3'-Dichlorobenzidine     | 21.11 | 252  | 239996   | 22.832 | ng/ul  | 96       |
| 80) Benzo(a)anthracene         | 21.17 | 228  | 795239   | 20.713 | ng/ul  | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.10 | 149  | 406535   | 19.914 | ng/ul  | 99       |
| 82) Chrysene                   | 21.22 | 228  | 738137   | 20.957 | ng/ul  | 99       |
| 84) Di-n-octyl phthalate       | 21.97 | 149  | 703192   | 18.487 | ng/ul  | 98       |
| 85) Benzo(b)fluoranthene       | 22.71 | 252  | 800279   | 20.646 | ng/ul  | 96       |
| 86) Benzo(k)fluoranthene       | 22.76 | 252  | 800620   | 21.369 | ng/ul# | 97       |
| 88) Benzo(a)pyrene             | 23.27 | 252  | 765574   | 20.907 | ng/ul# | 98       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.54 | 276  | 869197   | 20.109 | ng/ul  | 98       |
| 90) Dibenzo(a,h)anthracene     | 25.55 | 278  | 738806   | 20.241 | ng/ul# | 99       |
| 91) Benzo(g,h,i)perylene       | 26.20 | 276  | 707718   | 19.894 | ng/ul  | 97       |

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

