

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013118\  
 Data File : BM014028.D  
 Acq On : 31 Jan 2018 11:30  
 Operator : SJ/JU  
 Sample : SSTD00513  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD00513

Manual Integrations  
 APPROVED

Sohil  
 2/1/2018 6:04:06 PM

Quant Time: Jan 31 15:48:59 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 31 13:46:12 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.56  | 152  | 80526    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.33 | 136  | 340010   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.22 | 164  | 240758   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.97 | 188  | 589177   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.18 | 240  | 621661   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.36 | 264  | 613829   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.12  | 96  | 3563   | 1.75 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 0.00  | 99  | 0d     | 0.00 | ng/ul |      |
| 7) Bis-(2-Chloroethyl)ether-d  | 0.00  | 67  | 0d     | 0.00 | ng/ul |      |
| 9) 2-Chlorophenol-d4           | 7.10  | 132 | 21566  | 4.28 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 0.00  | 113 | 0d     | 0.00 | ng/ul |      |
| 19) Nitrobenzene-d5            | 8.72  | 128 | 12089  | 4.58 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.44  | 143 | 12233  | 4.39 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.99  | 165 | 24503m | 4.55 | ng/ul | 0.02 |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d     | 0.00 | ng/ul |      |
| 43) Dimethylphthalate-d6       | 13.64 | 166 | 99743  | 5.03 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.91 | 160 | 115382 | 4.56 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 0.00  | 143 | 0d     | 0.00 | ng/ul |      |
| 57) Fluorene-d10               | 15.22 | 176 | 92804  | 5.32 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 0.00  | 200 | 0d     | 0.00 | ng/ul |      |
| 70) Anthracene-d10             | 17.07 | 188 | 137264 | 4.78 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.38 | 212 | 155537 | 5.37 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.22 | 264 | 142105 | 5.04 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Ovalue |    |
|--------------------------------|-------|-----|--------|-------|--------|----|
| 2) 1,4-Dioxane                 | 3.16  | 88  | 4202   | 1.883 | ng/uL# | 63 |
| 10) 2-Chlorophenol             | 7.13  | 128 | 21918  | 4.366 | ng/ul# | 85 |
| 15) N-Nitroso-di-n-propylamine | 8.37  | 70  | 21199  | 5.525 | ng/ul# | 62 |
| 17) Hexachloroethane           | 8.62  | 117 | 12060  | 4.935 | ng/ul  | 89 |
| 20) Nitrobenzene               | 8.77  | 77  | 34887  | 5.119 | ng/ul  | 96 |
| 21) Isophorone                 | 9.29  | 82  | 54393  | 4.939 | ng/ul# | 96 |
| 23) 2-Nitrophenol              | 9.47  | 139 | 12377  | 4.260 | ng/ul# | 84 |
| 24) 2,4-Dimethylphenol         | 9.54  | 107 | 29600  | 4.832 | ng/ul  | 93 |
| 25) Bis(2-Chloroethoxy)methane | 9.77  | 93  | 29269  | 4.360 | ng/ul# | 88 |
| 27) 2,4-Dichlorophenol         | 10.01 | 162 | 22694  | 4.372 | ng/ul# | 77 |
| 28) Naphthalene                | 10.39 | 128 | 84220  | 4.979 | ng/ul  | 95 |
| 31) Hexachlorobutadiene        | 10.66 | 225 | 24609  | 5.776 | ng/ul  | 95 |
| 33) 4-Chloro-3-methylphenol    | 11.66 | 107 | 25333  | 4.815 | ng/ul# | 88 |
| 34) 2-Methylnaphthalene        | 12.02 | 142 | 62849  | 4.983 | ng/ul  | 97 |
| 36) 1,2,4,5-Tetrachlorobenzene | 12.39 | 216 | 43246  | 4.848 | ng/ul# | 92 |
| 38) 2,4,6-Trichlorophenol      | 12.65 | 196 | 21881  | 4.237 | ng/ul  | 97 |
| 39) 2,4,5-Trichlorophenol      | 12.74 | 196 | 25522m | 4.750 | ng/ul  |    |
| 40) 1,1'-Biphenyl              | 13.04 | 154 | 90134  | 4.498 | ng/ul  | 98 |
| 41) 2-Chloronaphthalene        | 13.09 | 162 | 70349  | 4.491 | ng/ul  | 97 |
| 42) 2-Nitroaniline             | 13.32 | 65  | 18841  | 4.360 | ng/ul  | 91 |
| 44) Dimethylphthalate          | 13.69 | 163 | 92023  | 4.717 | ng/ul  | 98 |
| 45) 2,6-Dinitrotoluene         | 13.82 | 165 | 16410  | 4.340 | ng/ul# | 94 |
| 47) Acenaphthylene             | 13.93 | 152 | 106988 | 4.621 | ng/ul  | 99 |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM013118\  
 Data File : BM014028.D  
 Acq On : 31 Jan 2018 11:30  
 Operator : SJ/JU  
 Sample : SSTD00513  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD00513

Manual Integrations  
 APPROVED

Sohil  
 2/1/2018 6:04:06 PM

Quant Time: Jan 31 15:48:59 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM013118.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 31 13:46:12 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 49) Acenaphthene               | 14.28 | 153  | 76597    | 4.777 | ng/ul  | 98       |
| 53) Dibenzofuran               | 14.63 | 168  | 118229   | 4.901 | ng/ul  | 92       |
| 54) 2,4-Dinitrotoluene         | 14.62 | 165  | 24129    | 4.378 | ng/ul# | 33       |
| 55) 2,3,4,6-Tetrachlorophenol  | 14.86 | 232  | 22799    | 4.511 | ng/ul# | 98       |
| 56) Diethylphthalate           | 15.06 | 149  | 98814    | 5.033 | ng/ul# | 89       |
| 58) Fluorene                   | 15.27 | 166  | 97327    | 4.926 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.27 | 204  | 55856    | 5.232 | ng/ul# | 90       |
| 64) N-Nitrosodiphenylamine     | 15.49 | 169  | 83308    | 4.808 | ng/ul  | 96       |
| 65) 4-Bromophenyl-phenylether  | 16.17 | 248  | 37007    | 5.318 | ng/ul  | 92       |
| 66) Hexachlorobenzene          | 16.28 | 284  | 38978    | 5.204 | ng/ul# | 83       |
| 69) Phenanthrene               | 17.02 | 178  | 162270   | 4.955 | ng/ul  | 97       |
| 71) Anthracene                 | 17.11 | 178  | 159725   | 4.728 | ng/ul  | 96       |
| 73) Di-n-butylphthalate        | 17.95 | 149  | 149237   | 4.551 | ng/ul  | 97       |
| 77) Pyrene                     | 19.41 | 202  | 192297   | 5.088 | ng/ul  | 97       |
| 78) Butylbenzylphthalate       | 20.32 | 149  | 64204    | 4.823 | ng/ul  | 96       |
| 80) Benzo(a)anthracene         | 21.16 | 228  | 192414   | 5.122 | ng/ul  | 96       |
| 81) Bis(2-ethylhexyl)phthalate | 21.10 | 149  | 87748    | 4.393 | ng/ul  | 98       |
| 82) Chrysene                   | 21.22 | 228  | 184914   | 5.366 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 22.72 | 252  | 174287   | 4.623 | ng/ul  | 97       |
| 86) Benzo(k)fluoranthene       | 22.76 | 252  | 188254   | 5.167 | ng/ul  | 95       |
| 88) Benzo(a)pyrene             | 23.27 | 252  | 167535   | 4.705 | ng/ul  | 97       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.54 | 276  | 192125   | 4.570 | ng/ul  | 96       |
| 90) Dibenzo(a,h)anthracene     | 25.56 | 278  | 161744   | 4.556 | ng/ul# | 98       |
| 91) Benzo(g,h,i)perylene       | 26.22 | 276  | 156318   | 4.518 | ng/ul# | 91       |

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

