

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM082416\  
 Data File : BM007275.D  
 Acq On : 24 Aug 2016 10:18  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02065

Manual Integrations  
 APPROVED

sohil  
 8/24/2016 7:28:18 PM

Quant Time: Aug 24 13:50:35 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM082316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 23 13:53:51 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.80  | 152  | 165347   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.59 | 136  | 833141   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.43 | 164  | 614523   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.17 | 188  | 1656769  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.36 | 240  | 2326474  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.63 | 264  | 2577008  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.31  | 96  | 25052   | 7.74  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 305114  | 19.53 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.13  | 67  | 166668  | 20.31 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.33  | 132 | 231932  | 20.09 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.50  | 113 | 271228  | 19.91 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.96  | 128 | 121008  | 19.78 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.67  | 143 | 144213  | 20.19 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.20 | 165 | 293214  | 20.62 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 380436  | 21.45 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.84 | 166 | 1063499 | 20.06 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.12 | 160 | 1240233 | 20.18 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 202531  | 20.24 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.43 | 176 | 935051  | 20.39 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.55 | 200 | 208582  | 20.02 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.27 | 188 | 1581142 | 20.43 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.56 | 212 | 1974874 | 20.30 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.48 | 264 | 2421211 | 20.40 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |         |       |       | Qvalue |
|--------------------------------|-------|-----|---------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.35  | 88  | 27778   | 7.70  | ng/uL | 99     |
| 4) Benzaldehyde                | 6.95  | 77  | 196711  | 22.11 | ng/ul | 95     |
| 6) Phenol                      | 6.99  | 94  | 322907  | 19.93 | ng/ul | 93     |
| 8) Bis(2-Chloroethyl)ether     | 7.23  | 93  | 221430  | 19.62 | ng/ul | 96     |
| 10) 2-Chlorophenol             | 7.36  | 128 | 238060  | 20.14 | ng/ul | 97     |
| 11) 2-Methylphenol             | 8.24  | 108 | 250168  | 19.76 | ng/ul | 99     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.33  | 45  | 252332  | 19.87 | ng/ul | 97     |
| 14) Acetophenone               | 8.62  | 105 | 435370  | 20.84 | ng/ul | 100    |
| 15) N-Nitroso-di-n-propylamine | 8.60  | 70  | 224723  | 20.24 | ng/ul | 96     |
| 16) 4-Methylphenol             | 8.56  | 108 | 280313  | 19.73 | ng/ul | 95     |
| 17) Hexachloroethane           | 8.87  | 117 | 95084   | 20.68 | ng/ul | 99     |
| 20) Nitrobenzene               | 9.00  | 77  | 317903  | 20.23 | ng/ul | 97     |
| 21) Isophorone                 | 9.52  | 82  | 644035  | 20.12 | ng/ul | 99     |
| 23) 2-Nitrophenol              | 9.70  | 139 | 156264  | 20.13 | ng/ul | 96     |
| 24) 2,4-Dimethylphenol         | 9.76  | 107 | 355614  | 20.76 | ng/ul | 99     |
| 25) Bis(2-Chloroethoxy)methane | 10.00 | 93  | 354814  | 19.67 | ng/ul | 98     |
| 27) 2,4-Dichlorophenol         | 10.23 | 162 | 290796  | 20.59 | ng/ul | 100    |
| 28) Naphthalene                | 10.63 | 128 | 824370  | 19.90 | ng/ul | 99     |
| 30) 4-Chloroaniline            | 10.75 | 127 | 391322  | 21.43 | ng/ul | 96     |
| 31) Hexachlorobutadiene        | 10.92 | 225 | 201308  | 21.09 | ng/ul | 97     |
| 32) Caprolactam                | 11.50 | 113 | 111516m | 19.60 | ng/ul |        |
| 33) 4-Chloro-3-methylphenol    | 11.87 | 107 | 359010  | 20.53 | ng/ul | 97     |
| 34) 2-Methylnaphthalene        | 12.25 | 142 | 694503  | 20.48 | ng/ul | 100    |

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM082416\  
 Data File : BM007275.D  
 Acq On : 24 Aug 2016 10:18  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02065

Manual Integrations  
 APPROVED

sohil  
 8/24/2016 7:28:18 PM

Quant Time: Aug 24 13:50:35 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM082316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Aug 23 13:53:51 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.62 | 216  | 418759   | 20.46 | ng/ul  | 98       |
| 37) Hexachlorocyclopentadiene  | 12.60 | 237  | 238673   | 19.34 | ng/ul  | 96       |
| 38) 2,4,6-Trichlorophenol      | 12.86 | 196  | 289363   | 19.99 | ng/ul  | 98       |
| 39) 2,4,5-Trichlorophenol      | 12.93 | 196  | 306545   | 20.32 | ng/ul  | 98       |
| 40) 1,1'-Biphenyl              | 13.26 | 154  | 961865   | 20.18 | ng/ul  | 98       |
| 41) 2-Chloronaphthalene        | 13.30 | 162  | 756156   | 20.24 | ng/ul  | 99       |
| 42) 2-Nitroaniline             | 13.51 | 65   | 253328   | 20.76 | ng/ul  | 96       |
| 44) Dimethylphthalate          | 13.89 | 163  | 1068889  | 20.24 | ng/ul  | 99       |
| 45) 2,6-Dinitrotoluene         | 14.01 | 165  | 223937   | 20.75 | ng/ul  | 91       |
| 47) Acenaphthylene             | 14.15 | 152  | 1275645  | 20.17 | ng/ul  | 100      |
| 48) 3-Nitroaniline             | 14.34 | 138  | 230506   | 20.92 | ng/ul  | 97       |
| 49) Acenaphthene               | 14.50 | 153  | 832053   | 20.22 | ng/ul  | 99       |
| 50) 2,4-Dinitrophenol          | 14.55 | 184  | 140718   | 19.23 | ng/ul  | 88       |
| 52) 4-Nitrophenol              | 14.65 | 109  | 226024   | 21.23 | ng/ul  | 86       |
| 53) Dibenzofuran               | 14.83 | 168  | 1243485  | 20.19 | ng/ul  | 95       |
| 54) 2,4-Dinitrotoluene         | 14.80 | 165  | 346659   | 20.95 | ng/ul  | 97       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.06 | 232  | 313839   | 20.28 | ng/ul# | 96       |
| 56) Diethylphthalate           | 15.26 | 149  | 1114541  | 20.23 | ng/ul  | 99       |
| 58) Fluorene                   | 15.48 | 166  | 1037294  | 20.45 | ng/ul  | 97       |
| 59) 4-Chlorophenyl-phenylether | 15.47 | 204  | 544819   | 20.55 | ng/ul  | 95       |
| 60) 4-Nitroaniline             | 15.50 | 138  | 260070   | 20.80 | ng/ul  | 86       |
| 63) 4,6-Dinitro-2-methylphenol | 15.56 | 198  | 224620   | 20.20 | ng/ul  | 100      |
| 64) N-Nitrosodiphenylamine     | 15.69 | 169  | 938507   | 20.29 | ng/ul  | 100      |
| 65) 4-Bromophenyl-phenylether  | 16.37 | 248  | 378800   | 20.25 | ng/ul  | 98       |
| 66) Hexachlorobenzene          | 16.48 | 284  | 425085   | 20.29 | ng/ul  | 94       |
| 67) Atrazine                   | 16.64 | 200  | 405095   | 20.74 | ng/ul  | 98       |
| 68) Pentachlorophenol          | 16.83 | 266  | 265860   | 19.67 | ng/ul  | 99       |
| 69) Phenanthrene               | 17.22 | 178  | 1808454  | 20.33 | ng/ul  | 99       |
| 71) Anthracene                 | 17.31 | 178  | 1871695  | 20.38 | ng/ul  | 99       |
| 72) Carbazole                  | 17.58 | 167  | 1664470  | 20.86 | ng/ul  | 99       |
| 73) Di-n-butylphthalate        | 18.14 | 149  | 2028981  | 20.21 | ng/ul  | 99       |
| 74) Fluoranthene               | 19.23 | 202  | 2418876  | 21.67 | ng/ul  | 97       |
| 77) Pyrene                     | 19.59 | 202  | 2527099  | 20.25 | ng/ul  | 98       |
| 78) Butylbenzylphthalate       | 20.49 | 149  | 1022755  | 19.69 | ng/ul  | 98       |
| 79) 3,3'-Dichlorobenzidine     | 21.27 | 252  | 981207   | 21.24 | ng/ul  | 98       |
| 80) Benzo(a)anthracene         | 21.34 | 228  | 2805561  | 20.46 | ng/ul  | 100      |
| 81) Bis(2-ethylhexyl)phthalate | 21.27 | 149  | 1497568  | 19.88 | ng/ul  | 100      |
| 82) Chrysene                   | 21.39 | 228  | 2512437  | 20.23 | ng/ul  | 99       |
| 84) Di-n-octyl phthalate       | 22.16 | 149  | 2718301  | 21.18 | ng/ul  | 99       |
| 85) Benzo(b)fluoranthene       | 22.94 | 252  | 3055882  | 20.06 | ng/ul  | 98       |
| 86) Benzo(k)fluoranthene       | 22.99 | 252  | 2995114  | 21.27 | ng/ul  | 100      |
| 88) Benzo(a)pyrene             | 23.53 | 252  | 3010417  | 20.65 | ng/ul  | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 25.92 | 276  | 3724812  | 20.82 | ng/ul  | 99       |
| 90) Dibenzo(a,h)anthracene     | 25.93 | 278  | 3111116  | 20.90 | ng/ul  | 98       |
| 91) Benzo(g,h,i)perylene       | 26.63 | 276  | 3156776  | 20.78 | ng/ul  | 96       |

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

