

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012417\  
 Data File : BM008856.D  
 Acq On : 25 Jan 2017 01:53  
 Operator : UM/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleID :  
 SSTD02047

Manual Integrations  
 APPROVED

mohammad  
 1/25/2017 3:46:28 PM

Quant Time: Jan 25 04:13:06 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM012317.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 25 03:34:06 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.93  | 152  | 66732    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.74 | 136  | 287035   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.57 | 164  | 234923   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.32 | 188  | 543334   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.48 | 240  | 642421   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.84 | 264  | 541363   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.39  | 96  | 11845  | 8.77  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.08  | 99  | 118588 | 20.98 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.26  | 67  | 100779 | 22.43 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.46  | 132 | 77374  | 21.01 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.63  | 113 | 89916  | 20.32 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.10  | 128 | 39888  | 21.85 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.82  | 143 | 45082  | 20.30 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.35 | 165 | 99485  | 21.40 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.87 | 131 | 105712 | 20.89 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.97 | 166 | 336134 | 21.84 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.26 | 160 | 400290 | 21.98 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.75 | 143 | 51379  | 19.96 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.56 | 176 | 313893 | 21.47 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.67 | 200 | 60150  | 18.59 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.42 | 188 | 515883 | 21.71 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.70 | 212 | 597169 | 21.19 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.69 | 264 | 502735 | 21.93 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.42  | 88  | 14221  | 8.69  | ng/uL# | 29     |
| 4) Benzaldehyde                | 7.08  | 77  | 118891 | 25.57 | ng/ul# | 68     |
| 6) Phenol                      | 7.11  | 94  | 127813 | 20.74 | ng/ul# | 47     |
| 8) Bis(2-Chloroethyl)ether     | 7.36  | 93  | 96114  | 20.82 | ng/ul# | 69     |
| 10) 2-Chlorophenol             | 7.49  | 128 | 84028  | 22.09 | ng/ul# | 69     |
| 11) 2-Methylphenol             | 8.36  | 108 | 92110  | 20.84 | ng/ul  | 95     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.46  | 45  | 182991 | 21.09 | ng/ul# | 91     |
| 14) Acetophenone               | 8.76  | 105 | 169762 | 20.93 | ng/ul# | 61     |
| 15) N-Nitroso-di-n-propylamine | 8.74  | 70  | 109281 | 20.67 | ng/ul# | 62     |
| 16) 4-Methylphenol             | 8.69  | 108 | 100468 | 21.10 | ng/ul  | 98     |
| 17) Hexachloroethane           | 9.02  | 117 | 49931  | 21.28 | ng/ul  | 94     |
| 20) Nitrobenzene               | 9.14  | 77  | 176890 | 21.98 | ng/ul# | 84     |
| 21) Isophorone                 | 9.66  | 82  | 273757 | 21.07 | ng/ul# | 89     |
| 23) 2-Nitrophenol              | 9.85  | 139 | 48247  | 20.95 | ng/ul# | 16     |
| 24) 2,4-Dimethylphenol         | 9.90  | 107 | 142421 | 21.69 | ng/ul# | 84     |
| 25) Bis(2-Chloroethoxy)methane | 10.14 | 93  | 140940 | 21.77 | ng/ul  | 95     |
| 27) 2,4-Dichlorophenol         | 10.37 | 162 | 99909  | 21.88 | ng/ul  | 93     |
| 28) Naphthalene                | 10.79 | 128 | 281411 | 21.24 | ng/ul  | 98     |
| 30) 4-Chloroaniline            | 10.90 | 127 | 117031 | 22.57 | ng/ul  | 93     |
| 31) Hexachlorobutadiene        | 11.06 | 225 | 98869  | 22.39 | ng/ul  | 94     |
| 32) Caprolactam                | 11.68 | 113 | 28125m | 20.03 | ng/ul  |        |
| 33) 4-Chloro-3-methylphenol    | 12.00 | 107 | 124088 | 20.95 | ng/ul# | 90     |
| 34) 2-Methylnaphthalene        | 12.39 | 142 | 226631 | 21.06 | ng/ul  | 98     |

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012417\  
 Data File : BM008856.D  
 Acq On : 25 Jan 2017 01:53  
 Operator : UM/SJ  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SSTD02047

Manual Integrations  
 APPROVED

mohammad  
 1/25/2017 3:46:28 PM

Quant Time: Jan 25 04:13:06 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM012317.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 25 03:34:06 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.76 | 216  | 154117   | 21.36 | ng/ul# | 93       |
| 37) Hexachlorocyclopentadiene  | 12.73 | 237  | 115432   | 20.69 | ng/ul# | 90       |
| 38) 2,4,6-Trichlorophenol      | 13.00 | 196  | 94538    | 21.70 | ng/ul  | 98       |
| 39) 2,4,5-Trichlorophenol      | 13.06 | 196  | 99840    | 21.34 | ng/ul  | 94       |
| 40) 1,1'-Biphenyl              | 13.40 | 154  | 322024   | 22.18 | ng/ul  | 97       |
| 41) 2-Chloronaphthalene        | 13.45 | 162  | 249529   | 21.84 | ng/ul  | 92       |
| 42) 2-Nitroaniline             | 13.65 | 65   | 127243   | 22.44 | ng/ul# | 63       |
| 44) Dimethylphthalate          | 14.02 | 163  | 322671   | 20.97 | ng/ul  | 99       |
| 45) 2,6-Dinitrotoluene         | 14.15 | 165  | 61170    | 20.75 | ng/ul# | 42       |
| 47) Acenaphthylene             | 14.29 | 152  | 384000   | 21.88 | ng/ul  | 98       |
| 48) 3-Nitroaniline             | 14.47 | 138  | 61474    | 22.05 | ng/ul# | 65       |
| 49) Acenaphthene               | 14.63 | 153  | 267490   | 21.67 | ng/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 14.68 | 184  | 34485    | 16.29 | ng/ul# | 28       |
| 52) 4-Nitrophenol              | 14.77 | 109  | 107649   | 20.28 | ng/ul# | 66       |
| 53) Dibenzofuran               | 14.97 | 168  | 407558   | 21.47 | ng/ul  | 88       |
| 54) 2,4-Dinitrotoluene         | 14.93 | 165  | 96748    | 21.38 | ng/ul# | 71       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.19 | 232  | 100561   | 21.40 | ng/ul# | 92       |
| 56) Diethylphthalate           | 15.38 | 149  | 358271   | 21.41 | ng/ul  | 94       |
| 58) Fluorene                   | 15.62 | 166  | 339708   | 21.42 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.61 | 204  | 200349   | 21.81 | ng/ul  | 91       |
| 60) 4-Nitroaniline             | 15.63 | 138  | 68285    | 21.97 | ng/ul# | 22       |
| 63) 4,6-Dinitro-2-methylphenol | 15.69 | 198  | 66076    | 19.74 | ng/ul# | 54       |
| 64) N-Nitrosodiphenylamine     | 15.82 | 169  | 299639   | 21.80 | ng/ul  | 97       |
| 65) 4-Bromophenyl-phenylether  | 16.50 | 248  | 131837   | 22.35 | ng/ul  | 92       |
| 66) Hexachlorobenzene          | 16.62 | 284  | 138340   | 21.62 | ng/ul# | 89       |
| 67) Atrazine                   | 16.77 | 200  | 129812   | 20.63 | ng/ul  | 94       |
| 68) Pentachlorophenol          | 16.96 | 266  | 85752    | 20.42 | ng/ul  | 96       |
| 69) Phenanthrene               | 17.36 | 178  | 574692   | 21.47 | ng/ul  | 99       |
| 71) Anthracene                 | 17.45 | 178  | 596455   | 21.79 | ng/ul  | 98       |
| 72) Carbazole                  | 17.72 | 167  | 512837   | 21.36 | ng/ul  | 98       |
| 73) Di-n-butylphthalate        | 18.27 | 149  | 605606   | 20.97 | ng/ul# | 96       |
| 74) Fluoranthene               | 19.37 | 202  | 717961   | 20.64 | ng/ul# | 96       |
| 77) Pyrene                     | 19.73 | 202  | 752017   | 21.82 | ng/ul# | 94       |
| 78) Butylbenzylphthalate       | 20.62 | 149  | 271126   | 21.13 | ng/ul# | 77       |
| 79) 3,3'-Dichlorobenzidine     | 21.40 | 252  | 261514   | 21.77 | ng/ul  | 95       |
| 80) Benzo(a)anthracene         | 21.47 | 228  | 744258   | 21.54 | ng/ul  | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.38 | 149  | 373910   | 20.97 | ng/ul  | 98       |
| 82) Chrysene                   | 21.52 | 228  | 681924   | 21.23 | ng/ul  | 96       |
| 84) Di-n-octyl phthalate       | 22.29 | 149  | 613744   | 19.53 | ng/ul# | 82       |
| 85) Benzo(b)fluoranthene       | 23.13 | 252  | 697651   | 22.18 | ng/ul# | 98       |
| 86) Benzo(k)fluoranthene       | 23.17 | 252  | 637499   | 21.34 | ng/ul# | 98       |
| 88) Benzo(a)pyrene             | 23.74 | 252  | 633742   | 21.96 | ng/ul# | 97       |
| 89) Indeno(1,2,3-cd)pyrene     | 26.26 | 276  | 671473   | 23.18 | ng/ul# | 91       |
| 90) Dibenzo(a,h)anthracene     | 26.27 | 278  | 565019   | 22.53 | ng/ul# | 95       |
| 91) Benzo(g,h,i)perylene       | 27.00 | 276  | 555178   | 23.57 | ng/ul# | 93       |

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

