

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA\_G\Data\BG012218\  
 Data File : BG032206.D  
 Acq On : 22 Jan 2018 11:01  
 Operator : SJ/JU  
 Sample : SSTD01075  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD01075

Quant Time: Jan 22 12:57:52 2018  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG012218.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 22 12:46:57 2018  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.73  | 152  | 39585    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.61 | 136  | 176865   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 15.36 | 164  | 122489   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 18.10 | 188  | 292810   | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 22.43 | 240  | 364065   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 26.12 | 264  | 384490   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.78  | 96  | 3233   | 5.29  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.84  | 99  | 28894  | 9.66  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 8.01  | 67  | 13273  | 8.88  | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 8.24  | 132 | 26425  | 10.48 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 9.43  | 113 | 27532  | 10.32 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.92  | 128 | 12074  | 9.77  | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.66 | 143 | 15511  | 9.85  | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 11.21 | 165 | 34090  | 11.00 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.73 | 131 | 20993  | 6.72  | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.74 | 166 | 105722 | 9.54  | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 15.06 | 160 | 128054 | 10.68 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.51 | 143 | 11153  | 6.70  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 16.34 | 176 | 96197  | 9.42  | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 16.44 | 200 | 13445  | 6.45  | ng/ul | 0.00 |
| 70) Anthracene-d10             | 18.19 | 188 | 141994 | 10.13 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 20.43 | 212 | 173194 | 10.71 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 25.86 | 264 | 175595 | 9.93  | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |       |        | Qvalue    |
|--------------------------------|-------|-----|-------|--------|-----------|
| 2) 1,4-Dioxane                 | 3.83  | 88  | 2693  | 4.112  | ng/uL# 80 |
| 4) Benzaldehyde                | 7.83  | 77  | 18372 | 12.417 | ng/ul# 78 |
| 6) Phenol                      | 7.87  | 94  | 28734 | 9.908  | ng/ul 92  |
| 8) Bis(2-Chloroethyl)ether     | 8.11  | 93  | 21100 | 10.279 | ng/ul 88  |
| 10) 2-Chlorophenol             | 8.27  | 128 | 24930 | 10.737 | ng/ul# 86 |
| 11) 2-Methylphenol             | 9.16  | 108 | 22499 | 9.951  | ng/ul 99  |
| 12) 2,2'-oxybis(1-Chloropropan | 9.26  | 45  | 20538 | 7.622  | ng/ul# 84 |
| 14) Acetophenone               | 9.56  | 105 | 37335 | 9.628  | ng/ul 92  |
| 15) N-Nitroso-di-n-propylamine | 9.54  | 70  | 17506 | 9.617  | ng/ul# 95 |
| 16) 4-Methylphenol             | 9.49  | 108 | 25744 | 10.263 | ng/ul 91  |
| 17) Hexachloroethane           | 9.85  | 117 | 10160 | 10.937 | ng/ul# 81 |
| 20) Nitrobenzene               | 9.96  | 77  | 25974 | 9.784  | ng/ul# 90 |
| 21) Isophorone                 | 10.49 | 82  | 47672 | 9.397  | ng/ul 95  |
| 23) 2-Nitrophenol              | 10.70 | 139 | 16234 | 10.700 | ng/ul 92  |
| 24) 2,4-Dimethylphenol         | 10.73 | 107 | 31622 | 10.421 | ng/ul 98  |
| 25) Bis(2-Chloroethoxy)methane | 10.97 | 93  | 29454 | 9.995  | ng/ul# 96 |
| 27) 2,4-Dichlorophenol         | 11.24 | 162 | 32634 | 11.052 | ng/ul 97  |
| 28) Naphthalene                | 11.65 | 128 | 82177 | 10.294 | ng/ul 97  |
| 30) 4-Chloroaniline            | 11.75 | 127 | 20558 | 6.911  | ng/ul 97  |
| 31) Hexachlorobutadiene        | 11.93 | 225 | 28964 | 11.337 | ng/ul 99  |
| 32) Caprolactam                | 12.47 | 113 | 7591  | 8.010  | ng/ul# 74 |
| 33) 4-Chloro-3-methylphenol    | 12.82 | 107 | 28241 | 9.815  | ng/ul# 83 |
| 34) 2-Methylnaphthalene        | 13.22 | 142 | 71015 | 10.308 | ng/ul 93  |

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 Operator : SJ/JU  
 Sample : SSTD01075  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD01075

Quant Time: Jan 22 12:57:52 2018  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG012218.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 22 12:46:57 2018  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc   | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|--------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 13.58 | 216  | 57471    | 12.179 | ng/ul# | 92       |
| 37) Hexachlorocyclopentadiene  | 13.55 | 237  | 16493    | 5.961  | ng/ul  | 100      |
| 38) 2,4,6-Trichlorophenol      | 13.81 | 196  | 31623    | 11.228 | ng/ul  | 95       |
| 39) 2,4,5-Trichlorophenol      | 13.88 | 196  | 31360    | 10.480 | ng/ul  | 96       |
| 40) 1,1'-Biphenyl              | 14.20 | 154  | 100275   | 11.263 | ng/ul# | 98       |
| 41) 2-Chloronaphthalene        | 14.25 | 162  | 79038    | 11.007 | ng/ul  | 96       |
| 42) 2-Nitroaniline             | 14.45 | 65   | 14955    | 8.837  | ng/ul  | 93       |
| 44) Dimethylphthalate          | 14.79 | 163  | 102745   | 10.117 | ng/ul  | 98       |
| 45) 2,6-Dinitrotoluene         | 14.92 | 165  | 19622    | 9.625  | ng/ul  | 87       |
| 47) Acenaphthylene             | 15.09 | 152  | 111704   | 9.828  | ng/ul  | 100      |
| 48) 3-Nitroaniline             | 15.25 | 138  | 12789    | 7.548  | ng/ul  | 84       |
| 49) Acenaphthene               | 15.43 | 153  | 80491    | 10.342 | ng/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 15.46 | 184  | 3989     | 3.065  | ng/ul# | 65       |
| 52) 4-Nitrophenol              | 15.53 | 109  | 10950    | 8.085  | ng/ul# | 77       |
| 53) Dibenzofuran               | 15.76 | 168  | 124790   | 10.546 | ng/ul  | 97       |
| 54) 2,4-Dinitrotoluene         | 15.70 | 165  | 28741    | 9.440  | ng/ul# | 79       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.97 | 232  | 30627    | 9.611  | ng/ul# | 91       |
| 56) Diethylphthalate           | 16.13 | 149  | 95418    | 8.949  | ng/ul  | 98       |
| 58) Fluorene                   | 16.40 | 166  | 100449   | 9.784  | ng/ul  | 99       |
| 59) 4-Chlorophenyl-phenylether | 16.38 | 204  | 62654    | 10.217 | ng/ul  | 98       |
| 60) 4-Nitroaniline             | 16.40 | 138  | 15698    | 7.801  | ng/ul  | 89       |
| 63) 4,6-Dinitro-2-methylphenol | 16.46 | 198  | 13386    | 6.700  | ng/ul# | 95       |
| 64) N-Nitrosodiphenylamine     | 16.58 | 169  | 85809    | 11.610 | ng/ul  | 93       |
| 65) 4-Bromophenyl-phenylether  | 17.27 | 248  | 42815    | 11.372 | ng/ul  | 93       |
| 66) Hexachlorobenzene          | 17.40 | 284  | 44711    | 10.442 | ng/ul  | 97       |
| 67) Atrazine                   | 17.51 | 200  | 38038    | 10.787 | ng/ul  | 91       |
| 68) Pentachlorophenol          | 17.74 | 266  | 15759    | 7.050  | ng/ul  | 95       |
| 69) Phenanthrene               | 18.14 | 178  | 155958   | 10.713 | ng/ul  | 98       |
| 71) Anthracene                 | 18.23 | 178  | 161866   | 10.679 | ng/ul  | 96       |
| 72) 1,2,3,4-Tetrachlorobenzene | 14.18 | 216  | 57018    | 13.961 | ng/uL  | 94       |
| 73) Pentachlorobenzene         | 15.67 | 250  | 60772    | 10.013 | ng/uL  | 94       |
| 74) Carbazole                  | 18.49 | 167  | 128120   | 10.322 | ng/ul  | 100      |
| 75) Di-n-butylphthalate        | 19.00 | 149  | 145088   | 10.527 | ng/ul  | 99       |
| 76) Fluoranthene               | 20.10 | 202  | 209646   | 11.268 | ng/ul# | 92       |
| 79) Pyrene                     | 20.46 | 202  | 211830   | 11.210 | ng/ul# | 90       |
| 80) Butylbenzylphthalate       | 21.32 | 149  | 66559    | 11.853 | ng/ul# | 84       |
| 81) 3,3'-Dichlorobenzidine     | 22.30 | 252  | 45146    | 6.778  | ng/ul  | 93       |
| 82) Benzo(a)anthracene         | 22.41 | 228  | 224290   | 11.598 | ng/ul  | 99       |
| 83) Bis(2-ethylhexyl)phthalate | 22.25 | 149  | 95755    | 11.656 | ng/ul  | 99       |
| 84) Chrysene                   | 22.48 | 228  | 205490   | 11.616 | ng/ul  | 98       |
| 86) Di-n-octyl phthalate       | 23.62 | 149  | 162085   | 10.695 | ng/ul  | 100      |
| 87) Benzo(b)fluoranthene       | 24.93 | 252  | 228242   | 10.663 | ng/ul# | 98       |
| 88) Benzo(k)fluoranthene       | 25.00 | 252  | 227625   | 10.908 | ng/ul# | 97       |
| 90) Benzo(a)pyrene             | 25.94 | 252  | 220887   | 10.724 | ng/ul# | 96       |
| 91) Indeno(1,2,3-cd)pyrene     | 30.38 | 276  | 213579   | 8.545  | ng/ul# | 95       |
| 92) Dibenzo(a,h)anthracene     | 30.46 | 278  | 188847   | 8.809  | ng/ul# | 94       |
| 93) Benzo(g,h,i)perylene       | 31.71 | 276  | 94841    | 4.577  | ng/ul# | 95       |

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

