

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101016\  
 Data File : BG024236.D  
 Acq On : 10 Oct 2016 11:07  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02011

Manual Integrations  
 APPROVED

sohil  
 10/11/2016 7:05:16 PM

Quant Time: Oct 10 13:35:26 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 07 06:34:32 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.30  | 152  | 156487   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.13 | 136  | 658724   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.91 | 164  | 555178   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.66 | 188  | 1209895m | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.96 | 240  | 1261474  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.41 | 264  | 1320500  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.65  | 96  | 25042   | 7.60  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.42  | 99  | 286340  | 19.61 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.61  | 67  | 187102  | 18.96 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.82  | 132 | 197885  | 19.68 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.99  | 113 | 237063  | 20.08 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.47  | 128 | 100766  | 21.56 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.19 | 143 | 115354  | 22.06 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.73 | 165 | 234234  | 20.56 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.25 | 131 | 283839  | 23.72 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.31 | 166 | 800122  | 20.64 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.61 | 160 | 925467  | 20.84 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 15.07 | 143 | 137257  | 19.97 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.90 | 176 | 737964  | 20.14 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 16.00 | 200 | 150886m | 20.75 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.75 | 188 | 1121523 | 20.52 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 20.02 | 212 | 1236245 | 21.00 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.17 | 264 | 1198204 | 20.17 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.69  | 88  | 26721  | 7.48  | ng/uL# | 63     |
| 4) Benzaldehyde                | 7.43  | 77  | 219003 | 21.31 | ng/ul# | 74     |
| 6) Phenol                      | 7.45  | 94  | 298061 | 20.11 | ng/ul# | 61     |
| 8) Bis(2-Chloroethyl)ether     | 7.71  | 93  | 203118 | 18.62 | ng/ul# | 77     |
| 10) 2-Chlorophenol             | 7.85  | 128 | 198928 | 19.52 | ng/ul# | 81     |
| 11) 2-Methylphenol             | 8.72  | 108 | 214526 | 19.55 | ng/ul  | 91     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.82  | 45  | 428529 | 19.67 | ng/ul# | 88     |
| 14) Acetophenone               | 9.12  | 105 | 353054 | 20.14 | ng/ul# | 68     |
| 15) N-Nitroso-di-n-propylamine | 9.10  | 70  | 208356 | 19.71 | ng/ul# | 74     |
| 16) 4-Methylphenol             | 9.05  | 108 | 231824 | 19.91 | ng/ul  | 97     |
| 17) Hexachloroethane           | 9.39  | 117 | 89635  | 19.02 | ng/ul# | 81     |
| 20) Nitrobenzene               | 9.52  | 77  | 315723 | 21.34 | ng/ul# | 77     |
| 21) Isophorone                 | 10.04 | 82  | 568428 | 21.14 | ng/ul# | 93     |
| 23) 2-Nitrophenol              | 10.22 | 139 | 121948 | 21.40 | ng/ul# | 52     |
| 24) 2,4-Dimethylphenol         | 10.26 | 107 | 291668 | 20.77 | ng/ul  | 94     |
| 25) Bis(2-Chloroethoxy)methane | 10.51 | 93  | 303879 | 20.62 | ng/ul  | 98     |
| 27) 2,4-Dichlorophenol         | 10.76 | 162 | 234768 | 21.50 | ng/ul  | 96     |
| 28) Naphthalene                | 11.18 | 128 | 681993 | 21.50 | ng/ul  | 97     |
| 30) 4-Chloroaniline            | 11.28 | 127 | 274597 | 22.61 | ng/ul  | 99     |
| 31) Hexachlorobutadiene        | 11.45 | 225 | 190195 | 21.74 | ng/ul  | 97     |
| 32) Caprolactam                | 12.04 | 113 | 81124  | 19.54 | ng/ul# | 12     |
| 33) 4-Chloro-3-methylphenol    | 12.37 | 107 | 288756 | 21.61 | ng/ul# | 79     |
| 34) 2-Methylnaphthalene        | 12.76 | 142 | 533531 | 21.16 | ng/ul  | 96     |

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 Data File : BG024236.D  
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 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 SSTD02011

Manual Integrations  
 APPROVED

sohil  
 10/11/2016 7:05:16 PM

Quant Time: Oct 10 13:35:26 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Oct 07 06:34:32 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 13.12 | 216  | 337810   | 20.63 | ng/ul  | 98       |
| 37) Hexachlorocyclopentadiene  | 13.09 | 237  | 228108   | 19.00 | ng/ul  | 95       |
| 38) 2,4,6-Trichlorophenol      | 13.35 | 196  | 217390   | 20.48 | ng/ul  | 95       |
| 39) 2,4,5-Trichlorophenol      | 13.41 | 196  | 235534   | 20.19 | ng/ul  | 93       |
| 40) 1,1'-Biphenyl              | 13.75 | 154  | 731339   | 20.34 | ng/ul  | 96       |
| 41) 2-Chloronaphthalene        | 13.80 | 162  | 576746   | 20.46 | ng/ul  | 95       |
| 42) 2-Nitroaniline             | 14.00 | 65   | 231512   | 21.58 | ng/ul# | 58       |
| 44) Dimethylphthalate          | 14.35 | 163  | 759260   | 19.98 | ng/ul  | 97       |
| 45) 2,6-Dinitrotoluene         | 14.48 | 165  | 159755   | 21.66 | ng/ul# | 70       |
| 47) Acenaphthylene             | 14.64 | 152  | 898074   | 21.02 | ng/ul  | 98       |
| 48) 3-Nitroaniline             | 14.81 | 138  | 151981   | 21.81 | ng/ul# | 54       |
| 49) Acenaphthene               | 14.98 | 153  | 624653   | 20.45 | ng/ul  | 98       |
| 50) 2,4-Dinitrophenol          | 15.01 | 184  | 97125    | 19.63 | ng/ul# | 67       |
| 52) 4-Nitrophenol              | 15.09 | 109  | 169083   | 20.30 | ng/ul# | 60       |
| 53) Dibenzofuran               | 15.31 | 168  | 943694m  | 20.99 | ng/ul  |          |
| 54) 2,4-Dinitrotoluene         | 15.26 | 165  | 226612   | 20.74 | ng/ul# | 56       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.52 | 232  | 233206   | 20.37 | ng/ul  | 93       |
| 56) Diethylphthalate           | 15.71 | 149  | 814342   | 20.32 | ng/ul  | 96       |
| 58) Fluorene                   | 15.95 | 166  | 760385   | 20.46 | ng/ul  | 100      |
| 59) 4-Chlorophenyl-phenylether | 15.94 | 204  | 411896   | 19.97 | ng/ul  | 95       |
| 60) 4-Nitroaniline             | 15.97 | 138  | 164138   | 20.46 | ng/ul# | 43       |
| 63) 4,6-Dinitro-2-methylphenol | 16.02 | 198  | 154524m  | 20.28 | ng/ul  |          |
| 64) N-Nitrosodiphenylamine     | 16.15 | 169  | 695942   | 20.34 | ng/ul  | 96       |
| 65) 4-Bromophenyl-phenylether  | 16.83 | 248  | 280015   | 19.98 | ng/ul  | 94       |
| 66) Hexachlorobenzene          | 16.94 | 284  | 320784   | 20.52 | ng/ul  | 98       |
| 67) Atrazine                   | 17.09 | 200  | 299465   | 21.40 | ng/ul# | 94       |
| 68) Pentachlorophenol          | 17.29 | 266  | 195349   | 20.56 | ng/ul  | 88       |
| 69) Phenanthrene               | 17.70 | 178  | 1282418m | 20.99 | ng/ul  |          |
| 71) Anthracene                 | 17.79 | 178  | 1296200  | 20.85 | ng/ul  | 99       |
| 72) Carbazole                  | 18.05 | 167  | 1086103  | 22.26 | ng/ul  | 98       |
| 73) Di-n-butylphthalate        | 18.58 | 149  | 1282903  | 21.50 | ng/ul# | 97       |
| 74) Fluoranthene               | 19.69 | 202  | 1439926m | 23.34 | ng/ul  |          |
| 77) Pyrene                     | 20.05 | 202  | 1463636  | 21.72 | ng/ul# | 86       |
| 78) Butylbenzylphthalate       | 20.92 | 149  | 562348   | 20.77 | ng/ul# | 83       |
| 79) 3,3'-Dichlorobenzidine     | 21.84 | 252  | 516957   | 22.00 | ng/ul# | 94       |
| 80) Benzo(a)anthracene         | 21.93 | 228  | 1489612  | 21.28 | ng/ul  | 99       |
| 81) Bis(2-ethylhexyl)phthalate | 21.81 | 149  | 803148   | 20.98 | ng/ul# | 99       |
| 82) Chrysene                   | 22.00 | 228  | 1410399  | 21.17 | ng/ul  | 98       |
| 84) Di-n-octyl phthalate       | 23.10 | 149  | 1364488  | 22.02 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 24.30 | 252  | 1465640  | 19.99 | ng/ul# | 92       |
| 86) Benzo(k)fluoranthene       | 24.38 | 252  | 1436406  | 20.34 | ng/ul# | 91       |
| 88) Benzo(a)pyrene             | 25.24 | 252  | 1422811  | 20.38 | ng/ul# | 92       |
| 89) Indeno(1,2,3-cd)pyrene     | 29.38 | 276  | 1729810  | 20.91 | ng/ul# | 79       |
| 90) Dibenzo(a,h)anthracene     | 29.45 | 278  | 1453376  | 20.78 | ng/ul# | 84       |
| 91) Benzo(g,h,i)perylene       | 30.63 | 276  | 1420617  | 20.43 | ng/ul# | 78       |

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

