

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080915\  
 Data File : BG018337.D  
 Acq On : 9 Aug 2015 20:24  
 Operator : UM/TP  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02085

Manual Integrations  
 APPROVED

MMDadoda  
 8/10/2015 7:25:35 PM

Quant Time: Aug 10 07:49:29 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080915.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 10 07:38:18 2015  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.07  | 152  | 82121    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.87 | 136  | 371143   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.68 | 164  | 224016   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 17.43 | 188  | 539761   | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.69 | 240  | 526073   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 24.92 | 264  | 537717   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.52  | 96  | 14451  | 7.78  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.21  | 99  | 153376 | 20.58 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.39  | 67  | 95937  | 20.77 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.60  | 132 | 117141 | 20.52 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.76  | 113 | 129534 | 20.68 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.22  | 128 | 59016  | 20.40 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.95  | 143 | 58761  | 19.94 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.49 | 165 | 128830 | 20.60 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.00 | 131 | 167053 | 23.64 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 14.09 | 166 | 386846 | 20.52 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 14.38 | 160 | 491034 | 20.69 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.86 | 143 | 70587  | 20.82 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.67 | 176 | 336681 | 20.51 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.78 | 200 | 51680  | 19.47 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 17.53 | 188 | 535613 | 20.75 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.81 | 212 | 571325 | 20.93 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 24.71 | 264 | 579346 | 20.29 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |       | Ovalue |
|--------------------------------|-------|-----|--------|-------|-------|--------|
| 2) 1,4-Dioxane                 | 3.56  | 88  | 15912  | 8.00  | ng/ul | 90     |
| 4) Benzaldehyde                | 7.20  | 77  | 102923 | 23.54 | ng/ul | 97     |
| 6) Phenol                      | 7.24  | 94  | 157664 | 20.74 | ng/ul | 97     |
| 8) Bis(2-Chloroethyl)ether     | 7.48  | 93  | 124278 | 21.25 | ng/ul | 97     |
| 10) 2-Chlorophenol             | 7.63  | 128 | 120179 | 20.66 | ng/ul | 97     |
| 11) 2-Methylphenol             | 8.50  | 108 | 123693 | 20.98 | ng/ul | 100    |
| 12) 2,2'-oxybis(1-Chloropropan | 8.60  | 45  | 195539 | 20.83 | ng/ul | 98     |
| 14) Acetophenone               | 8.89  | 105 | 192242 | 21.25 | ng/ul | 99     |
| 15) N-Nitroso-di-n-propylamine | 8.87  | 70  | 108810 | 20.96 | ng/ul | 95     |
| 16) 4-Methylphenol             | 8.82  | 108 | 134539 | 20.91 | ng/ul | 89     |
| 17) Hexachloroethane           | 9.16  | 117 | 49920  | 19.91 | ng/ul | 97     |
| 20) Nitrobenzene               | 9.26  | 77  | 150706 | 20.64 | ng/ul | 94     |
| 21) Isophorone                 | 9.79  | 82  | 287500 | 20.47 | ng/ul | 99     |
| 23) 2-Nitrophenol              | 9.98  | 139 | 66385  | 20.29 | ng/ul | 96     |
| 24) 2,4-Dimethylphenol         | 10.03 | 107 | 151188 | 20.62 | ng/ul | 92     |
| 25) Bis(2-Chloroethoxy)methane | 10.28 | 93  | 178215 | 20.36 | ng/ul | 98     |
| 27) 2,4-Dichlorophenol         | 10.51 | 162 | 120693 | 20.60 | ng/ul | 97     |
| 28) Naphthalene                | 10.92 | 128 | 401531 | 20.47 | ng/ul | 98     |
| 30) 4-Chloroaniline            | 11.02 | 127 | 172312 | 23.97 | ng/ul | 95     |
| 31) Hexachlorobutadiene        | 11.22 | 225 | 77170  | 20.86 | ng/ul | 95     |
| 32) Caprolactam                | 11.79 | 113 | 47811  | 20.26 | ng/ul | 85     |
| 33) 4-Chloro-3-methylphenol    | 12.13 | 107 | 140237 | 20.66 | ng/ul | 94     |
| 34) 2-Methylnaphthalene        | 12.52 | 142 | 283744 | 19.97 | ng/ul | 97     |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080915\  
 Data File : BG018337.D  
 Acq On : 9 Aug 2015 20:24  
 Operator : UM/TP  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 SSTD02085

Manual Integrations  
 APPROVED

MMDadoda  
 8/10/2015 7:25:35 PM

Quant Time: Aug 10 07:49:29 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080915.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Aug 10 07:38:18 2015  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 35) 1-Methylnaphthalene        | 12.74 | 142  | 279177   | 20.13 | ng/ul  | 100      |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.88 | 216  | 144006   | 20.05 | ng/ul  | 98       |
| 38) Hexachlorocyclopentadiene  | 12.87 | 237  | 85346    | 19.96 | ng/ul  | 89       |
| 39) 2,4,6-Trichlorophenol      | 13.12 | 196  | 100685   | 20.58 | ng/ul  | 95       |
| 40) 2,4,5-Trichlorophenol      | 13.18 | 196  | 111359   | 21.17 | ng/ul  | 97       |
| 41) 1,1'-Biphenyl              | 13.52 | 154  | 390260   | 20.87 | ng/ul  | 96       |
| 42) 2-Chloronaphthalene        | 13.57 | 162  | 296500   | 20.42 | ng/ul  | 96       |
| 43) 2-Nitroaniline             | 13.75 | 65   | 97319    | 20.76 | ng/ul  | 97       |
| 45) Dimethylphthalate          | 14.14 | 163  | 388630   | 20.90 | ng/ul  | 99       |
| 46) 2,6-Dinitrotoluene         | 14.25 | 165  | 79202    | 21.39 | ng/ul  | 98       |
| 48) Acenaphthylene             | 14.41 | 152  | 486430   | 20.43 | ng/ul  | 98       |
| 49) 3-Nitroaniline             | 14.58 | 138  | 86006    | 22.77 | ng/ul  | 94       |
| 50) Acenaphthene               | 14.75 | 153  | 344357   | 20.62 | ng/ul  | 97       |
| 51) 2,4-Dinitrophenol          | 14.78 | 184  | 32458    | 17.98 | ng/ul  | 92       |
| 53) 4-Nitrophenol              | 14.87 | 109  | 61744    | 20.95 | ng/ul  | 92       |
| 54) Dibenzofuran               | 15.08 | 168  | 450009   | 20.62 | ng/ul  | 99       |
| 55) 2,4-Dinitrotoluene         | 15.03 | 165  | 113538   | 21.49 | ng/ul  | 97       |
| 56) 2,3,4,6-Tetrachlorophenol  | 15.30 | 232  | 93538    | 20.41 | ng/ul# | 96       |
| 57) Diethylphthalate           | 15.50 | 149  | 409676   | 20.74 | ng/ul  | 98       |
| 59) Fluorene                   | 15.73 | 166  | 386431   | 20.59 | ng/ul  | 97       |
| 60) 4-Chlorophenyl-phenylether | 15.73 | 204  | 180443   | 20.31 | ng/ul  | 98       |
| 61) 4-Nitroaniline             | 15.74 | 138  | 89083    | 21.35 | ng/ul  | 98       |
| 64) 4,6-Dinitro-2-methylphenol | 15.79 | 198  | 54600    | 19.28 | ng/ul# | 95       |
| 65) N-Nitrosodiphenylamine     | 15.93 | 169  | 330445   | 20.35 | ng/ul  | 96       |
| 66) 4-Bromophenyl-phenylether  | 16.61 | 248  | 118873   | 20.37 | ng/ul  | 98       |
| 67) Hexachlorobenzene          | 16.73 | 284  | 126379   | 20.47 | ng/ul  | 98       |
| 68) Atrazine                   | 16.88 | 200  | 132984   | 21.88 | ng/ul  | 98       |
| 69) Pentachlorophenol          | 17.07 | 266  | 85245    | 20.62 | ng/ul  | 99       |
| 70) Phenanthrene               | 17.47 | 178  | 600023   | 20.49 | ng/ul  | 99       |
| 72) Anthracene                 | 17.56 | 178  | 619190   | 20.75 | ng/ul  | 99       |
| 73) 1,2,3,4-Tetrachlorobenzene | 13.49 | 216  | 153513   | 20.56 | ng/uL  | 96       |
| 74) Pentachlorobenzene         | 15.01 | 250  | 144102   | 20.21 | ng/uL  | 90       |
| 75) Carbazole                  | 17.83 | 167  | 586380   | 22.41 | ng/ul  | 98       |
| 76) Di-n-butylphthalate        | 18.39 | 149  | 718423   | 21.02 | ng/ul  | 99       |
| 77) Fluoranthene               | 19.47 | 202  | 718376   | 22.97 | ng/ul  | 99       |
| 80) Pvrene                     | 19.83 | 202  | 737195   | 20.72 | ng/ul  | 99       |
| 81) Butylbenzylphthalate       | 20.72 | 149  | 314781   | 20.92 | ng/ul  | 93       |
| 82) 3,3'-Dichlorobenzidine     | 21.59 | 252  | 249707   | 22.86 | ng/ul  | 97       |
| 83) Benzo(a)anthracene         | 21.67 | 228  | 679118   | 20.82 | ng/ul  | 97       |
| 84) Bis(2-ethylhexyl)phthalate | 21.60 | 149  | 453048   | 21.31 | ng/ul  | 99       |
| 85) Chrysene                   | 21.74 | 228  | 636604   | 20.87 | ng/ul  | 98       |
| 87) Di-n-octyl phthalate       | 22.82 | 149  | 760173   | 21.91 | ng/ul  | 100      |
| 88) Benzo(b)fluoranthene       | 23.90 | 252  | 678799   | 20.09 | ng/ul  | 99       |
| 89) Benzo(k)fluoranthene       | 23.97 | 252  | 680890   | 20.93 | ng/ul  | 97       |
| 91) Benzo(a)pyrene             | 24.77 | 252  | 661264   | 20.59 | ng/ul  | 99       |
| 92) Indeno(1,2,3-cd)pyrene     | 28.61 | 276  | 753032m  | 20.26 | ng/ul  |          |
| 93) Dibenzo(a,h)anthracene     | 28.68 | 278  | 631529   | 20.45 | ng/ul  | 99       |
| 94) Benzo(a,h,i)perylene       | 29.76 | 276  | 623193   | 19.97 | ng/ul  | 97       |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080915\  
Data File : BG018337.D  
Acq On : 9 Aug 2015 20:24  
Operator : UM/TP  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
SSTD02085

Manual Integrations  
APPROVED

MMDadoda  
8/10/2015 7:25:35 PM

Quant Time: Aug 10 07:49:29 2015  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080915.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Aug 10 07:38:18 2015  
Response via : Initial Calibration

| Internal Standards | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|--------------------|------|------|----------|------|-------|----------|
|--------------------|------|------|----------|------|-------|----------|

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

