

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121216\  
 Data File : BG025111.D  
 Acq On : 12 Dec 2016 10:36  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 SSTD02031

Manual Integrations  
 APPROVED

sohil  
 12/13/2016 6:06:20 PM

Quant Time: Dec 13 00:24:06 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 12 03:23:22 2016  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.24  | 152  | 93859    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 11.07 | 136  | 401874   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.86 | 164  | 355408   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.60 | 188  | 813035   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.90 | 240  | 1014075  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.32 | 264  | 1017567  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |       |
|--------------------------------|-------|-----|--------|-------|-------|-------|
| 3) 1,4-Dioxane-d8              | 3.59  | 96  | 15287  | 8.41  | ng/uL | -0.01 |
| 5) Phenol-d5                   | 7.39  | 99  | 173580 | 21.44 | ng/ul | 0.00  |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.56  | 67  | 104981 | 20.96 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.77  | 132 | 123150 | 21.75 | ng/ul | 0.00  |
| 13) 4-Methylphenol-d8          | 8.94  | 113 | 136548 | 20.18 | ng/ul | 0.00  |
| 19) Nitrobenzene-d5            | 9.42  | 128 | 58146  | 20.43 | ng/ul | 0.00  |
| 22) 2-Nitrophenol-d4           | 10.14 | 143 | 73159  | 20.62 | ng/ul | 0.00  |
| 26) 2,4-Dichlorophenol-d3      | 10.68 | 165 | 150191 | 22.01 | ng/ul | 0.00  |
| 29) 4-Chloroaniline-d4         | 11.20 | 131 | 172178 | 25.68 | ng/ul | 0.00  |
| 43) Dimethylphthalate-d6       | 14.25 | 166 | 487167 | 19.93 | ng/ul | 0.00  |
| 46) Acenaphthylene-d8          | 14.56 | 160 | 574513 | 21.45 | ng/ul | 0.00  |
| 51) 4-Nitrophenol-d4           | 15.05 | 143 | 81415  | 19.38 | ng/ul | 0.00  |
| 57) Fluorene-d10               | 15.85 | 176 | 477534 | 20.75 | ng/ul | 0.00  |
| 62) 4,6-Dinitro-2-methylphenol | 15.96 | 200 | 97533  | 19.40 | ng/ul | 0.00  |
| 70) Anthracene-d10             | 17.70 | 188 | 684356 | 19.94 | ng/ul | 0.00  |
| 76) Pyrene-d10                 | 19.98 | 212 | 855769 | 20.49 | ng/ul | 0.00  |
| 87) Benzo(a)pyrene-d12         | 25.08 | 264 | 871845 | 21.15 | ng/ul | 0.00  |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.63  | 88  | 17133  | 7.24  | ng/uL  | 98     |
| 4) Benzaldehyde                | 7.37  | 77  | 130921 | 21.48 | ng/ul  | 94     |
| 6) Phenol                      | 7.42  | 94  | 184005 | 22.14 | ng/ul  | 92     |
| 8) Bis(2-Chloroethyl)ether     | 7.65  | 93  | 120513 | 19.82 | ng/ul  | 96     |
| 10) 2-Chlorophenol             | 7.80  | 128 | 118370 | 20.79 | ng/ul# | 84     |
| 11) 2-Methylphenol             | 8.68  | 108 | 129703 | 21.20 | ng/ul  | 98     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.76  | 45  | 232777 | 22.51 | ng/ul  | 98     |
| 14) Acetophenone               | 9.07  | 105 | 207300 | 20.13 | ng/ul# | 89     |
| 15) N-Nitroso-di-n-propylamine | 9.04  | 70  | 114841 | 19.31 | ng/ul  | 92     |
| 16) 4-Methylphenol             | 9.01  | 108 | 142922 | 20.99 | ng/ul  | 98     |
| 17) Hexachloroethane           | 9.32  | 117 | 48299  | 19.15 | ng/ul  | 91     |
| 20) Nitrobenzene               | 9.46  | 77  | 179894 | 20.57 | ng/ul  | 100    |
| 21) Isophorone                 | 9.98  | 82  | 309548 | 18.70 | ng/ul# | 94     |
| 23) 2-Nitrophenol              | 10.17 | 139 | 76674  | 21.31 | ng/ul  | 89     |
| 24) 2,4-Dimethylphenol         | 10.22 | 107 | 171417 | 20.83 | ng/ul  | 94     |
| 25) Bis(2-Chloroethoxy)methane | 10.45 | 93  | 171769 | 20.37 | ng/ul# | 93     |
| 27) 2,4-Dichlorophenol         | 10.71 | 162 | 149016 | 22.49 | ng/ul  | 97     |
| 28) Naphthalene                | 11.12 | 128 | 376689 | 20.16 | ng/ul  | 97     |
| 30) 4-Chloroaniline            | 11.23 | 127 | 166903 | 24.35 | ng/ul# | 86     |
| 31) Hexachlorobutadiene        | 11.38 | 225 | 115234 | 20.23 | ng/ul  | 93     |
| 32) Caprolactam                | 11.99 | 113 | 52285  | 19.01 | ng/ul  | 98     |
| 33) 4-Chloro-3-methylphenol    | 12.33 | 107 | 179019 | 21.92 | ng/ul  | 98     |
| 34) 2-Methylnaphthalene        | 12.70 | 142 | 305958 | 20.75 | ng/ul  | 91     |

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02031

Manual Integrations  
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Quant Time: Dec 13 00:24:06 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 12 03:23:22 2016  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 13.06 | 216  | 215759   | 20.14 | ng/ul  | 93       |
| 37) Hexachlorocyclopentadiene  | 13.03 | 237  | 91558    | 14.56 | ng/ul  | 96       |
| 38) 2,4,6-Trichlorophenol      | 13.30 | 196  | 143985   | 21.40 | ng/ul  | 91       |
| 39) 2,4,5-Trichlorophenol      | 13.38 | 196  | 162160   | 21.99 | ng/ul  | 97       |
| 40) 1,1'-Biphenyl              | 13.70 | 154  | 442736   | 20.55 | ng/ul  | 97       |
| 41) 2-Chloronaphthalene        | 13.74 | 162  | 346948   | 20.23 | ng/ul  | 99       |
| 42) 2-Nitroaniline             | 13.95 | 65   | 151184   | 22.52 | ng/ul  | 94       |
| 44) Dimethylphthalate          | 14.30 | 163  | 476823   | 19.85 | ng/ul  | 96       |
| 45) 2,6-Dinitrotoluene         | 14.44 | 165  | 101625   | 19.69 | ng/ul  | 93       |
| 47) Acenaphthylene             | 14.59 | 152  | 531922   | 20.43 | ng/ul  | 99       |
| 48) 3-Nitroaniline             | 14.77 | 138  | 103911   | 23.24 | ng/ul# | 91       |
| 49) Acenaphthene               | 14.92 | 153  | 378167   | 21.20 | ng/ul  | 99       |
| 50) 2,4-Dinitrophenol          | 14.98 | 184  | 54269    | 16.18 | ng/ul  | 89       |
| 52) 4-Nitrophenol              | 15.08 | 109  | 107569   | 21.29 | ng/ul  | 100      |
| 53) Dibenzofuran               | 15.25 | 168  | 584702   | 20.73 | ng/ul  | 97       |
| 54) 2,4-Dinitrotoluene         | 15.22 | 165  | 160658   | 20.65 | ng/ul# | 96       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.48 | 232  | 167282   | 21.47 | ng/ul  | 92       |
| 56) Diethylphthalate           | 15.65 | 149  | 491508   | 19.32 | ng/ul  | 96       |
| 58) Fluorene                   | 15.91 | 166  | 479199   | 20.87 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.88 | 204  | 263171   | 20.35 | ng/ul  | 89       |
| 60) 4-Nitroaniline             | 15.93 | 138  | 106958   | 19.84 | ng/ul  | 97       |
| 63) 4,6-Dinitro-2-methylphenol | 15.98 | 198  | 102080   | 19.60 | ng/ul# | 94       |
| 64) N-Nitrosodiphenylamine     | 16.10 | 169  | 453237   | 21.67 | ng/ul  | 92       |
| 65) 4-Bromophenyl-phenylether  | 16.78 | 248  | 195206   | 21.26 | ng/ul  | 91       |
| 66) Hexachlorobenzene          | 16.90 | 284  | 220529   | 21.33 | ng/ul  | 97       |
| 67) Atrazine                   | 17.04 | 200  | 190711   | 19.18 | ng/ul  | 97       |
| 68) Pentachlorophenol          | 17.25 | 266  | 118168   | 19.03 | ng/ul# | 87       |
| 69) Phenanthrene               | 17.64 | 178  | 798863   | 20.59 | ng/ul  | 97       |
| 71) Anthracene                 | 17.74 | 178  | 812549   | 20.37 | ng/ul  | 99       |
| 72) Carbazole                  | 18.01 | 167  | 697319   | 20.95 | ng/ul  | 98       |
| 73) Di-n-butylphthalate        | 18.53 | 149  | 826307   | 19.66 | ng/ul  | 99       |
| 74) Fluoranthene               | 19.64 | 202  | 1008485  | 21.87 | ng/ul  | 99       |
| 77) Pyrene                     | 20.01 | 202  | 1006952  | 20.66 | ng/ul# | 97       |
| 78) Butylbenzylphthalate       | 20.86 | 149  | 390951   | 20.48 | ng/ul  | 95       |
| 79) 3,3'-Dichlorobenzidine     | 21.78 | 252  | 398400   | 22.54 | ng/ul  | 96       |
| 80) Benzo(a)anthracene         | 21.87 | 228  | 1078656  | 20.47 | ng/ul  | 98       |
| 81) Bis(2-ethylhexyl)phthalate | 21.73 | 149  | 512051   | 19.49 | ng/ul  | 98       |
| 82) Chrysene                   | 21.95 | 228  | 1016430  | 20.63 | ng/ul  | 97       |
| 84) Di-n-octyl phthalate       | 22.99 | 149  | 895328   | 21.67 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 24.22 | 252  | 1052342m | 20.86 | ng/ul  |          |
| 86) Benzo(k)fluoranthene       | 24.29 | 252  | 1017105m | 21.21 | ng/ul  |          |
| 88) Benzo(a)pyrene             | 25.15 | 252  | 1019644  | 21.02 | ng/ul# | 99       |
| 89) Indeno(1,2,3-cd)pyrene     | 29.25 | 276  | 1065598m | 17.67 | ng/ul  |          |
| 90) Dibenzo(a,h)anthracene     | 29.30 | 278  | 886522m  | 17.45 | ng/ul  |          |
| 91) Benzo(g,h,i)perylene       | 30.48 | 276  | 845922m  | 16.81 | ng/ul  |          |

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

