

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050117\  
 Data File : BG026818.D  
 Acq On : 1 May 2017 22:30  
 Operator : SJ/MA  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
 SSTD02069

Manual Integrations  
 APPROVED

mohammad  
 5/3/2017 12:12:44 PM

Quant Time: May 02 03:19:45 2017  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 02 01:49:56 2017  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.01  | 152  | 145570   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.81 | 136  | 753714   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.62 | 164  | 433381   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.36 | 188  | 812629   | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.60 | 240  | 758561   | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 24.76 | 264  | 784915   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.42  | 96  | 23928   | 7.40  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.18  | 99  | 292800  | 22.54 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.33  | 67  | 165197  | 21.62 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.55  | 132 | 228461  | 20.72 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.72  | 113 | 239042  | 22.17 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.16  | 128 | 121354  | 19.26 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.89  | 143 | 137242  | 20.03 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.44 | 165 | 230155  | 21.14 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.95 | 131 | 325714m | 22.35 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.02 | 166 | 673572  | 19.31 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.32 | 160 | 890163  | 20.32 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.84 | 143 | 115233  | 18.33 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.61 | 176 | 596839  | 19.60 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.73 | 200 | 94182   | 19.09 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.46 | 188 | 799399  | 20.27 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.73 | 212 | 775662  | 19.37 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 24.55 | 264 | 754957  | 20.34 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       |        | Ovalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.46  | 88  | 25959  | 7.04  | ng/uL# | 84     |
| 4) Benzaldehyde                | 7.14  | 77  | 115338 | 18.37 | ng/ul# | 66     |
| 6) Phenol                      | 7.21  | 94  | 300977 | 22.86 | ng/ul# | 76     |
| 8) Bis(2-Chloroethyl)ether     | 7.43  | 93  | 209095 | 20.57 | ng/ul  | 94     |
| 10) 2-Chlorophenol             | 7.58  | 128 | 223007 | 20.55 | ng/ul# | 82     |
| 11) 2-Methylphenol             | 8.46  | 108 | 234812 | 22.53 | ng/ul  | 97     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.54  | 45  | 259650 | 24.47 | ng/ul# | 91     |
| 14) Acetophenone               | 8.82  | 105 | 351131 | 22.17 | ng/ul# | 74     |
| 15) N-Nitroso-di-n-propylamine | 8.81  | 70  | 166393 | 21.30 | ng/ul# | 73     |
| 16) 4-Methylphenol             | 8.79  | 108 | 256383 | 22.49 | ng/ul  | 93     |
| 17) Hexachloroethane           | 9.09  | 117 | 81377  | 19.94 | ng/ul# | 75     |
| 20) Nitrobenzene               | 9.21  | 77  | 247100 | 20.33 | ng/ul# | 78     |
| 21) Isophorone                 | 9.73  | 82  | 465086 | 19.93 | ng/ul# | 91     |
| 23) 2-Nitrophenol              | 9.92  | 139 | 142425 | 19.77 | ng/ul# | 61     |
| 24) 2,4-Dimethylphenol         | 9.99  | 107 | 268512 | 20.44 | ng/ul# | 79     |
| 25) Bis(2-Chloroethoxy)methane | 10.20 | 93  | 311393 | 19.54 | ng/ul  | 95     |
| 27) 2,4-Dichlorophenol         | 10.46 | 162 | 223378 | 20.57 | ng/ul  | 95     |
| 28) Naphthalene                | 10.86 | 128 | 777959 | 19.83 | ng/ul  | 98     |
| 30) 4-Chloroaniline            | 10.97 | 127 | 310138 | 22.56 | ng/ul  | 94     |
| 31) Hexachlorobutadiene        | 11.14 | 225 | 103267 | 18.80 | ng/ul  | 87     |
| 32) Caprolactam                | 11.71 | 113 | 87312  | 21.52 | ng/ul# | 72     |
| 33) 4-Chloro-3-methylphenol    | 12.10 | 107 | 269218 | 22.87 | ng/ul  | 86     |
| 34) 2-Methylnaphthalene        | 12.46 | 142 | 596408 | 20.48 | ng/ul  | 96     |

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 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02069

Manual Integrations  
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Quant Time: May 02 03:19:45 2017  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG042717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue May 02 01:49:56 2017  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.83 | 216  | 224056   | 19.61 | ng/ul# | 92       |
| 37) Hexachlorocyclopentadiene  | 12.80 | 237  | 92875    | 18.45 | ng/ul  | 99       |
| 38) 2,4,6-Trichlorophenol      | 13.07 | 196  | 162942   | 20.73 | ng/ul  | 96       |
| 39) 2,4,5-Trichlorophenol      | 13.15 | 196  | 171324   | 21.09 | ng/ul  | 97       |
| 40) 1,1'-Biphenyl              | 13.45 | 154  | 726182   | 19.93 | ng/ul  | 98       |
| 41) 2-Chloronaphthalene        | 13.50 | 162  | 542724   | 19.82 | ng/ul  | 97       |
| 42) 2-Nitroaniline             | 13.71 | 65   | 173952   | 22.23 | ng/ul# | 72       |
| 44) Dimethylphthalate          | 14.07 | 163  | 665350   | 19.48 | ng/ul  | 98       |
| 45) 2,6-Dinitrotoluene         | 14.19 | 165  | 143052   | 19.85 | ng/ul# | 82       |
| 47) Acenaphthylene             | 14.35 | 152  | 904256   | 20.26 | ng/ul  | 98       |
| 48) 3-Nitroaniline             | 14.53 | 138  | 155013   | 20.44 | ng/ul# | 76       |
| 49) Acenaphthene               | 14.68 | 153  | 619343   | 19.95 | ng/ul  | 96       |
| 50) 2,4-Dinitrophenol          | 14.74 | 184  | 74350m   | 20.18 | ng/ul  |          |
| 52) 4-Nitrophenol              | 14.85 | 109  | 66858    | 18.13 | ng/ul# | 37       |
| 53) Dibenzofuran               | 15.02 | 168  | 815867   | 19.74 | ng/ul  | 94       |
| 54) 2,4-Dinitrotoluene         | 14.98 | 165  | 204680   | 19.85 | ng/ul# | 74       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.25 | 232  | 125056   | 19.59 | ng/ul# | 75       |
| 56) Diethylphthalate           | 15.42 | 149  | 695719   | 19.47 | ng/ul  | 98       |
| 58) Fluorene                   | 15.66 | 166  | 663698   | 19.68 | ng/ul  | 100      |
| 59) 4-Chlorophenyl-phenylether | 15.65 | 204  | 284700   | 19.36 | ng/ul  | 93       |
| 60) 4-Nitroaniline             | 15.69 | 138  | 145830   | 17.45 | ng/ul# | 50       |
| 63) 4,6-Dinitro-2-methylphenol | 15.75 | 198  | 99868    | 19.43 | ng/ul# | 92       |
| 64) N-Nitrosodiphenylamine     | 15.86 | 169  | 551885   | 20.66 | ng/ul  | 98       |
| 65) 4-Bromophenyl-phenylether  | 16.54 | 248  | 163623   | 20.18 | ng/ul# | 83       |
| 66) Hexachlorobenzene          | 16.67 | 284  | 170892   | 19.57 | ng/ul# | 88       |
| 67) Atrazine                   | 16.81 | 200  | 172312   | 20.64 | ng/ul# | 90       |
| 68) Pentachlorophenol          | 17.02 | 266  | 79458    | 19.92 | ng/ul  | 94       |
| 69) Phenanthrene               | 17.40 | 178  | 934373   | 20.41 | ng/ul  | 100      |
| 71) Anthracene                 | 17.49 | 178  | 961561   | 20.48 | ng/ul  | 100      |
| 72) Carbazole                  | 17.76 | 167  | 874646   | 22.02 | ng/ul  | 98       |
| 73) Di-n-butylphthalate        | 18.31 | 149  | 1127284  | 21.82 | ng/ul# | 96       |
| 74) Fluoranthene               | 19.41 | 202  | 968261   | 22.78 | ng/ul# | 81       |
| 77) Pyrene                     | 19.76 | 202  | 994577   | 19.54 | ng/ul# | 76       |
| 78) Butylbenzylphthalate       | 20.64 | 149  | 540054m  | 22.85 | ng/ul  |          |
| 79) 3,3'-Dichlorobenzidine     | 21.50 | 252  | 310570   | 21.30 | ng/ul# | 95       |
| 80) Benzo(a)anthracene         | 21.59 | 228  | 892132   | 20.11 | ng/ul  | 98       |
| 81) Bis(2-ethylhexyl)phthalate | 21.48 | 149  | 764262   | 22.66 | ng/ul# | 99       |
| 82) Chrysene                   | 21.65 | 228  | 828867   | 20.11 | ng/ul  | 98       |
| 84) Di-n-octyl phthalate       | 22.66 | 149  | 1308861  | 24.27 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 23.76 | 252  | 916374   | 20.43 | ng/ul# | 94       |
| 86) Benzo(k)fluoranthene       | 23.82 | 252  | 867443   | 19.68 | ng/ul# | 93       |
| 88) Benzo(a)pyrene             | 24.62 | 252  | 883754   | 20.02 | ng/ul# | 93       |
| 89) Indeno(1,2,3-cd)pyrene     | 28.37 | 276  | 1056960  | 20.20 | ng/ul# | 80       |
| 90) Dibenzo(a,h)anthracene     | 28.42 | 278  | 888570   | 20.03 | ng/ul# | 88       |
| 91) Benzo(g,h,i)perylene       | 29.50 | 276  | 874386   | 20.15 | ng/ul# | 80       |

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

