

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG080516\  
 Data File : BG023519.D  
 Acq On : 6 Aug 2016 11:47  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02013

Manual Integrations  
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8/8/2016 6:56:56 PM

Quant Time: Aug 08 03:18:55 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 01 04:44:16 2004  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.12  | 152  | 126849   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.95 | 136  | 570501   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.78 | 164  | 420185   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.54 | 188  | 1051180  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.86 | 240  | 1045714  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 25.24 | 264  | 1036467  | 20.00 | ng/ul | 0.00     |

#### System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.49  | 96  | 23756   | 7.96  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.27  | 99  | 252178  | 19.73 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.43  | 67  | 164561  | 20.10 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.65  | 132 | 172992  | 19.84 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.83  | 113 | 201904  | 20.24 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.29  | 128 | 94717   | 21.06 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 10.02 | 143 | 107007  | 20.72 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.57 | 165 | 206370  | 21.18 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.08 | 131 | 267558  | 23.41 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.18 | 166 | 714900  | 20.78 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.47 | 160 | 829542  | 21.42 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.99 | 143 | 120265  | 20.48 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.78 | 176 | 655461  | 20.62 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.90 | 200 | 139333  | 20.66 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.64 | 188 | 1016597 | 21.47 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.93 | 212 | 1112713 | 21.02 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 25.00 | 264 | 988568  | 21.07 | ng/ul | 0.00 |

#### Target Compounds

|                                |       |     |        |       |        | Qvalue |
|--------------------------------|-------|-----|--------|-------|--------|--------|
| 2) 1,4-Dioxane                 | 3.53  | 88  | 24179  | 7.72  | ng/uL# | 77     |
| 4) Benzaldehyde                | 7.25  | 77  | 168385 | 21.62 | ng/ul  | 88     |
| 6) Phenol                      | 7.30  | 94  | 270827 | 20.35 | ng/ul# | 73     |
| 8) Bis(2-Chloroethyl)ether     | 7.53  | 93  | 200089 | 20.71 | ng/ul# | 86     |
| 10) 2-Chlorophenol             | 7.68  | 128 | 178482 | 20.10 | ng/ul# | 85     |
| 11) 2-Methylphenol             | 8.56  | 108 | 193084 | 19.27 | ng/ul  | 91     |
| 12) 2,2'-oxybis(1-Chloropropan | 8.65  | 45  | 277990 | 21.03 | ng/ul  | 99     |
| 14) Acetophenone               | 8.95  | 105 | 331819 | 20.74 | ng/ul# | 86     |
| 15) N-Nitroso-di-n-propylamine | 8.93  | 70  | 194824 | 20.24 | ng/ul# | 90     |
| 16) 4-Methylphenol             | 8.89  | 108 | 222498 | 19.81 | ng/ul  | 96     |
| 17) Hexachloroethane           | 9.21  | 117 | 76256  | 19.90 | ng/ul  | 98     |
| 20) Nitrobenzene               | 9.34  | 77  | 273067 | 20.72 | ng/ul# | 85     |
| 21) Isophorone                 | 9.85  | 82  | 518960 | 21.41 | ng/ul  | 95     |
| 23) 2-Nitrophenol              | 10.05 | 139 | 117132 | 21.01 | ng/ul# | 70     |
| 24) 2,4-Dimethylphenol         | 10.11 | 107 | 257125 | 21.02 | ng/ul# | 88     |
| 25) Bis(2-Chloroethoxy)methane | 10.34 | 93  | 298542 | 21.35 | ng/ul  | 97     |
| 27) 2,4-Dichlorophenol         | 10.59 | 162 | 204446 | 20.81 | ng/ul  | 95     |
| 28) Naphthalene                | 11.00 | 128 | 608135 | 21.00 | ng/ul  | 96     |
| 30) 4-Chloroaniline            | 11.11 | 127 | 256552 | 22.62 | ng/ul  | 97     |
| 31) Hexachlorobutadiene        | 11.28 | 225 | 133165 | 20.35 | ng/ul  | 96     |
| 32) Caprolactam                | 11.87 | 113 | 77847  | 19.08 | ng/ul# | 37     |
| 33) 4-Chloro-3-methylphenol    | 12.23 | 107 | 258110 | 20.41 | ng/ul  | 83     |
| 34) 2-Methylnaphthalene        | 12.60 | 142 | 493462 | 21.01 | ng/ul  | 95     |

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 Data File : BG023519.D  
 Acq On : 6 Aug 2016 11:47  
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 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 SSTD02013

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Quant Time: Aug 08 03:18:55 2016  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Sep 01 04:44:16 2004  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 36) 1,2,4,5-Tetrachlorobenzene | 12.97 | 216  | 274894   | 21.31 | ng/ul  | 97       |
| 37) Hexachlorocyclopentadiene  | 12.94 | 237  | 144304   | 20.22 | ng/ul  | 98       |
| 38) 2,4,6-Trichlorophenol      | 13.21 | 196  | 190490   | 21.32 | ng/ul  | 95       |
| 39) 2,4,5-Trichlorophenol      | 13.29 | 196  | 202535   | 21.15 | ng/ul  | 94       |
| 40) 1,1'-Biphenyl              | 13.61 | 154  | 659316   | 21.35 | ng/ul  | 97       |
| 41) 2-Chloronaphthalene        | 13.65 | 162  | 513687   | 21.39 | ng/ul  | 96       |
| 42) 2-Nitroaniline             | 13.86 | 65   | 209881   | 21.41 | ng/ul# | 61       |
| 44) Dimethylphthalate          | 14.22 | 163  | 704361   | 20.68 | ng/ul  | 99       |
| 45) 2,6-Dinitrotoluene         | 14.35 | 165  | 154016   | 20.85 | ng/ul# | 88       |
| 47) Acenaphthylene             | 14.50 | 152  | 862964   | 21.31 | ng/ul  | 99       |
| 48) 3-Nitroaniline             | 14.69 | 138  | 150407   | 21.66 | ng/ul# | 55       |
| 49) Acenaphthene               | 14.85 | 153  | 577623   | 20.93 | ng/ul  | 95       |
| 50) 2,4-Dinitrophenol          | 14.90 | 184  | 85169    | 18.69 | ng/ul# | 76       |
| 52) 4-Nitrophenol              | 15.01 | 109  | 117988   | 20.06 | ng/ul# | 77       |
| 53) Dibenzofuran               | 15.18 | 168  | 855562   | 20.93 | ng/ul  | 90       |
| 54) 2,4-Dinitrotoluene         | 15.15 | 165  | 234482   | 20.95 | ng/ul# | 74       |
| 55) 2,3,4,6-Tetrachlorophenol  | 15.41 | 232  | 196065   | 19.93 | ng/ul# | 96       |
| 56) Diethylphthalate           | 15.59 | 149  | 741376   | 20.82 | ng/ul  | 97       |
| 58) Fluorene                   | 15.83 | 166  | 704015   | 20.67 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.82 | 204  | 357402   | 20.86 | ng/ul  | 93       |
| 60) 4-Nitroaniline             | 15.86 | 138  | 166731   | 20.70 | ng/ul# | 56       |
| 63) 4,6-Dinitro-2-methylphenol | 15.91 | 198  | 149470   | 21.11 | ng/ul# | 85       |
| 64) N-Nitrosodiphenylamine     | 16.03 | 169  | 641992   | 21.85 | ng/ul  | 99       |
| 65) 4-Bromophenyl-phenylether  | 16.72 | 248  | 252601   | 20.95 | ng/ul  | 97       |
| 66) Hexachlorobenzene          | 16.84 | 284  | 276084   | 21.50 | ng/ul  | 95       |
| 67) Atrazine                   | 16.99 | 200  | 259816   | 21.41 | ng/ul  | 96       |
| 68) Pentachlorophenol          | 17.19 | 266  | 144837m  | 20.44 | ng/ul  |          |
| 69) Phenanthrene               | 17.58 | 178  | 1166105  | 21.42 | ng/ul  | 100      |
| 71) Anthracene                 | 17.68 | 178  | 1204983  | 21.69 | ng/ul  | 98       |
| 72) Carbazole                  | 17.95 | 167  | 1067557  | 23.85 | ng/ul  | 99       |
| 73) Di-n-butylphthalate        | 18.49 | 149  | 1253037  | 22.35 | ng/ul# | 94       |
| 74) Fluoranthene               | 19.60 | 202  | 1377690  | 24.72 | ng/ul# | 92       |
| 77) Pyrene                     | 19.96 | 202  | 1385094  | 21.43 | ng/ul  | 95       |
| 78) Butylbenzylphthalate       | 20.84 | 149  | 551651m  | 22.25 | ng/ul  |          |
| 79) 3,3'-Dichlorobenzidine     | 21.75 | 252  | 422351   | 22.92 | ng/ul# | 97       |
| 80) Benzo(a)anthracene         | 21.84 | 228  | 1263575  | 21.46 | ng/ul  | 97       |
| 81) Bis(2-ethylhexyl)phthalate | 21.72 | 149  | 764040   | 23.17 | ng/ul# | 98       |
| 82) Chrysene                   | 21.91 | 228  | 1157870  | 21.62 | ng/ul  | 96       |
| 84) Di-n-octyl phthalate       | 22.98 | 149  | 1260403  | 24.08 | ng/ul  | 100      |
| 85) Benzo(b)fluoranthene       | 24.16 | 252  | 1263593  | 21.43 | ng/ul# | 95       |
| 86) Benzo(k)fluoranthene       | 24.23 | 252  | 1208586  | 20.82 | ng/ul# | 92       |
| 88) Benzo(a)pyrene             | 25.07 | 252  | 1203169  | 21.14 | ng/ul# | 95       |
| 89) Indeno(1,2,3-cd)pyrene     | 29.09 | 276  | 1412650m | 21.12 | ng/ul  |          |
| 90) Dibenzo(a,h)anthracene     | 29.16 | 278  | 1200815m | 21.05 | ng/ul  |          |
| 91) Benzo(g,h,i)perylene       | 30.31 | 276  | 1151760  | 20.77 | ng/ul# | 86       |

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

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 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleID :  
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