

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012517\  
 Data File : BG025633.D  
 Acq On : 25 Jan 2017 12:53  
 Operator : SJ/MA  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02017

Quant Time: Jan 25 18:10:44 2017  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG011817.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jan 25 02:56:13 2017  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 20% Max. Rel. Area : 150%

|      | Compound                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 129   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.472 | 0.524 | -11.0 | 150#  | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.400 | 0.363 | 9.3   | 124   | 0.00     |
| 4    | Benzaldehyde                | 1.180 | 1.392 | -18.0 | 120   | 0.00     |
| 5 S  | Phenol-d5                   | 1.714 | 1.720 | -0.4  | 126   | 0.00     |
| 6    | Phenol                      | 1.801 | 1.812 | -0.6  | 128   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.107 | 1.132 | -2.3  | 120   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.286 | 1.350 | -5.0  | 128   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.209 | 1.241 | -2.6  | 125   | 0.00     |
| 10   | 2-Chlorophenol              | 1.197 | 1.180 | 1.4   | 120   | 0.00     |
| 11   | 2-Methylphenol              | 1.341 | 1.360 | -1.4  | 131   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.443 | 2.628 | -7.6  | 132   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.457 | 1.550 | -6.4  | 136   | 0.00     |
| 14   | Acetophenone                | 2.276 | 2.417 | -6.2  | 133   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.362 | 1.424 | -4.6  | 125   | 0.00     |
| 16   | 4-Methylphenol              | 1.471 | 1.491 | -1.4  | 125   | 0.00     |
| 17   | Hexachloroethane            | 0.558 | 0.584 | -4.7  | 133   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 124   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.149 | -4.9  | 121   | 0.00     |
| 20   | Nitrobenzene                | 0.454 | 0.510 | -12.3 | 131   | 0.00     |
| 21   | Isophorone                  | 0.853 | 0.908 | -6.4  | 126   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.169 | 0.181 | -7.1  | 124   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.173 | 0.187 | -8.1  | 132   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.416 | 0.434 | -4.3  | 123   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.430 | 0.477 | -10.9 | 133   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.334 | 0.363 | -8.7  | 130   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.324 | 0.355 | -9.6  | 128   | 0.00     |
| 28   | Naphthalene                 | 0.923 | 1.031 | -11.7 | 132   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.373 | 0.442 | -18.5 | 128   | 0.00     |
| 30   | 4-Chloroaniline             | 0.376 | 0.435 | -15.7 | 127   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.283 | 0.301 | -6.4  | 127   | 0.00     |
| 32   | Caprolactam                 | 0.136 | 0.150 | -10.3 | 133   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.394 | 0.439 | -11.4 | 126   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.754 | 0.830 | -10.1 | 129   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 134   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.602 | 0.635 | -5.5  | 127   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.238 | 0.264 | -10.9 | 180#  | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.365 | 0.362 | 0.8   | 119   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.382 | 0.393 | -2.9  | 125   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.239 | 1.276 | -3.0  | 126   | 0.00     |
| 41   | 2-Chloronaphthalene         | 0.962 | 1.034 | -7.5  | 134   | 0.00     |
| 42   | 2-Nitroaniline              | 0.417 | 0.439 | -5.3  | 128   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.412 | 1.485 | -5.2  | 124   | 0.00     |
| 44   | Dimethylphthalate           | 1.389 | 1.440 | -3.7  | 125   | 0.00     |

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
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 LabSampleId :  
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Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 20% Max. Rel. Area : 150%

|      | Compound                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.286 | 0.318 | -11.2 | 135   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.526 | 1.659 | -8.7  | 131   | 0.00     |
| 47   | Acenaphthylene              | 1.485 | 1.614 | -8.7  | 130   | 0.00     |
| 48   | 3-Nitroaniline              | 0.259 | 0.306 | -18.1 | 140   | 0.00     |
| 49 C | Acenaphthene                | 1.058 | 1.108 | -4.7  | 126   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.168 | 0.161 | 4.2   | 139   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.194 | 0.183 | 5.7   | 116   | 0.00     |
| 52   | 4-Nitrophenol               | 0.268 | 0.249 | 7.1   | 118   | 0.00     |
| 53   | Dibenzofuran                | 1.626 | 1.747 | -7.4  | 131   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.436 | 0.469 | -7.6  | 135   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.405 | 0.381 | 5.9   | 114   | 0.00     |
| 56   | Diethylphthalate            | 1.501 | 1.530 | -1.9  | 123   | 0.00     |
| 57 S | Fluorene-d10                | 1.331 | 1.399 | -5.1  | 128   | 0.00     |
| 58   | Fluorene                    | 1.328 | 1.433 | -7.9  | 132   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.749 | 0.784 | -4.7  | 126   | 0.00     |
| 60   | 4-Nitroaniline              | 0.250 | 0.279 | -11.6 | 130   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 123   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.123 | 0.123 | 0.0   | 131   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.127 | 0.119 | 6.3   | 115   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.516 | 0.560 | -8.5  | 125   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.232 | 0.243 | -4.7  | 121   | 0.00     |
| 66   | Hexachlorobenzene           | 0.264 | 0.277 | -4.9  | 125   | 0.00     |
| 67   | Atrazine                    | 0.242 | 0.262 | -8.3  | 124   | 0.00     |
| 68 C | Pentachlorophenol           | 0.110 | 0.085 | 22.7  | 107   | 0.00     |
| 69   | Phenanthrene                | 0.972 | 1.063 | -9.4  | 125   | 0.00     |
| 70 S | Anthracene-d10              | 0.852 | 0.930 | -9.2  | 126   | 0.00     |
| 71   | Anthracene                  | 0.977 | 1.082 | -10.7 | 125   | 0.00     |
| 72   | Carbazole                   | 0.811 | 0.901 | -11.1 | 121   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.048 | 1.118 | -6.7  | 122   | 0.00     |
| 74 C | Fluoranthene                | 1.148 | 1.299 | -13.2 | 120   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 117   | 0.00     |
| 76 S | Pyrene-d10                  | 0.832 | 0.917 | -10.2 | 117   | 0.00     |
| 77   | Pyrene                      | 0.983 | 1.094 | -11.3 | 118   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.382 | 0.428 | -12.0 | 123   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.367 | 0.436 | -18.8 | 121   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.057 | 1.156 | -9.4  | 118   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.534 | 0.609 | -14.0 | 123   | 0.00     |
| 82   | Chrysene                    | 0.994 | 1.107 | -11.4 | 119   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 119   | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.858 | 0.997 | -16.2 | 124   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.035 | 1.088 | -5.1  | 119   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 0.971 | 1.075 | -10.7 | 120   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.841 | 0.895 | -6.4  | 119   | 0.00     |

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Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
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Max. RRF Dev : 20% Max. Rel. Area : 150%

|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 0.979 | 1.074 | -9.7 | 123   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.240 | 1.307 | -5.4 | 120   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.033 | 1.090 | -5.5 | 117   | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 1.018 | 1.093 | -7.4 | 121   | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 0