

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122816\  
 Data File : BG025394.D  
 Acq On : 28 Dec 2016 12:50  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02026

Quant Time: Dec 29 00:07:28 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 20 03:02:46 2016  
 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    | 72    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.504 | 0.399 | 20.8#  | 59    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.387 | 0.359 | 7.2    | 67    | 0.00     |
| 4    | Benzaldehyde                | 1.299 | 1.346 | -3.6   | 71    | 0.00     |
| 5 S  | Phenol-d5                   | 1.725 | 1.700 | 1.4    | 71    | 0.00     |
| 6    | Phenol                      | 1.771 | 1.772 | -0.1   | 70    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.067 | 1.106 | -3.7   | 74    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.296 | 1.309 | -1.0   | 74    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.207 | 1.175 | 2.7    | 69    | 0.00     |
| 10   | 2-Chlorophenol              | 1.213 | 1.248 | -2.9   | 73    | 0.00     |
| 11   | 2-Methylphenol              | 1.304 | 1.352 | -3.7   | 75    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.204 | 2.424 | -10.0  | 77    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.442 | 1.505 | -4.4   | 76    | 0.00     |
| 14   | Acetophenone                | 2.194 | 2.155 | 1.8    | 71    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.267 | 1.356 | -7.0   | 75    | 0.00     |
| 16   | 4-Methylphenol              | 1.451 | 1.399 | 3.6    | 74    | 0.00     |
| 17   | Hexachloroethane            | 0.537 | 0.569 | -6.0   | 74    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 66    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.154 | -8.5   | 72    | 0.00     |
| 20   | Nitrobenzene                | 0.435 | 0.496 | -14.0  | 77    | 0.00     |
| 21   | Isophorone                  | 0.824 | 0.918 | -11.4  | 72    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.177 | 0.197 | -11.3  | 75    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.179 | 0.198 | -10.6  | 73    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.410 | 0.465 | -13.4  | 73    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.420 | 0.453 | -7.9   | 70    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.340 | 0.358 | -5.3   | 69    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.330 | 0.371 | -12.4  | 73    | 0.00     |
| 28   | Naphthalene                 | 0.930 | 0.997 | -7.2   | 71    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.334 | 0.434 | -29.9# | 85    | 0.00     |
| 30   | 4-Chloroaniline             | 0.341 | 0.427 | -25.2# | 81    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.283 | 0.334 | -18.0  | 77    | 0.00     |
| 32   | Caprolactam                 | 0.137 | 0.146 | -6.6   | 72    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.406 | 0.439 | -8.1   | 71    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.734 | 0.801 | -9.1   | 70    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 65    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.603 | 0.697 | -15.6  | 78    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.354 | 0.306 | 13.6   | 62    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.379 | 0.419 | -10.6  | 73    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.415 | 0.419 | -1.0   | 65    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.213 | 1.388 | -14.4  | 74    | 0.00     |
| 41   | 2-Chloronaphthalene         | 0.965 | 1.046 | -8.4   | 73    | 0.00     |
| 42   | 2-Nitroaniline              | 0.378 | 0.455 | -20.4# | 76    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.376 | 1.588 | -15.4  | 73    | 0.00     |
| 44   | Dimethylphthalate           | 1.352 | 1.522 | -12.6  | 73    | 0.00     |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122816\  
 Data File : BG025394.D  
 Acq On : 28 Dec 2016 12:50  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02026

Quant Time: Dec 29 00:07:28 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Dec 20 03:02:46 2016  
 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) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.291 | 0.311 | -6.9   | 68    | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.507 | 1.707 | -13.3  | 73    | 0.00     |
| 47   | Acenaphthylene              | 1.465 | 1.623 | -10.8  | 71    | 0.00     |
| 48   | 3-Nitroaniline              | 0.252 | 0.306 | -21.4# | 76    | 0.00     |
| 49 C | Acenaphthene                | 1.004 | 1.133 | -12.8  | 74    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.189 | 0.189 | 0.0    | 73    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.236 | 0.226 | 4.2    | 61    | 0.00     |
| 52   | 4-Nitrophenol               | 0.284 | 0.296 | -4.2   | 69    | 0.00     |
| 53   | Dibenzofuran                | 1.587 | 1.776 | -11.9  | 73    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.438 | 0.479 | -9.4   | 71    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.439 | 0.447 | -1.8   | 66    | 0.00     |
| 56   | Diethylphthalate            | 1.432 | 1.647 | -15.0  | 73    | 0.00     |
| 57 S | Fluorene-d10                | 1.295 | 1.471 | -13.6  | 73    | 0.00     |
| 58   | Fluorene                    | 1.292 | 1.442 | -11.6  | 71    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.728 | 0.818 | -12.4  | 73    | 0.00     |
| 60   | 4-Nitroaniline              | 0.303 | 0.309 | -2.0   | 67    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 69    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.124 | 0.117 | 5.6    | 68    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.128 | 0.125 | 2.3    | 70    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.515 | 0.561 | -8.9   | 73    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.226 | 0.245 | -8.4   | 74    | 0.00     |
| 66   | Hexachlorobenzene           | 0.254 | 0.288 | -13.4  | 76    | 0.00     |
| 67   | Atrazine                    | 0.245 | 0.280 | -14.3  | 76    | 0.00     |
| 68 C | Pentachlorophenol           | 0.153 | 0.106 | 30.7#  | 49    | 0.00     |
| 69   | Phenanthrene                | 0.955 | 1.078 | -12.9  | 74    | 0.00     |
| 70 S | Anthracene-d10              | 0.844 | 0.944 | -11.8  | 75    | 0.00     |
| 71   | Anthracene                  | 0.981 | 1.094 | -11.5  | 72    | 0.00     |
| 72   | Carbazole                   | 0.819 | 0.945 | -15.4  | 73    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.034 | 1.194 | -15.5  | 76    | 0.00     |
| 74 C | Fluoranthene                | 1.134 | 1.449 | -27.8# | 77    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 72    | 0.00     |
| 76 S | Pyrene-d10                  | 0.824 | 0.909 | -10.3  | 77    | 0.00     |
| 77   | Pyrene                      | 0.961 | 1.074 | -11.8  | 76    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.376 | 0.393 | -4.5   | 72    | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.349 | 0.398 | -14.0  | 78    | 0.00     |
| 80   | Benzo(a)anthracene          | 1.039 | 1.171 | -12.7  | 78    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.518 | 0.571 | -10.2  | 77    | 0.00     |
| 82   | Chrysene                    | 0.972 | 1.086 | -11.7  | 78    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 71    | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.812 | 0.972 | -19.7  | 78    | 0.00     |
| 85   | Benzo(b)fluoranthene        | 0.992 | 1.137 | -14.6  | 78    | 0.00     |
| 86   | Benzo(k)fluoranthene        | 0.943 | 1.069 | -13.4  | 78    | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.810 | 0.927 | -14.4  | 79    | 0.00     |

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122816\  
Data File : BG025394.D  
Acq On : 28 Dec 2016 12:50  
Operator : UM/SJ  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02026

Quant Time: Dec 29 00:07:28 2016  
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Tue Dec 20 03:02:46 2016  
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               | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|------------------------|-------|-------|-------|-------|----------|
| 88 C | Benzo(a)pyrene         | 0.954 | 1.087 | -13.9 | 78    | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.185 | 1.359 | -14.7 | 80    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 0.998 | 1.146 | -14.8 | 81    | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 0.989 | 1.122 | -13.4 | 79    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 2