

Data Path : Z:\HPCHEM1\BNA\_G\Data\BG082517\  
 Data File : BG028484.D  
 Acq On : 25 Aug 2017 9:12  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02086

Quant Time: Aug 25 23:04:58 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 25 03:52:53 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 153#  | 0.00     |
| 2    | 1,4-Dioxane                 | 0.451 | 0.480 | -6.4  | 167#  | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.429 | 0.369 | 14.0  | 145   | 0.00     |
| 4    | Benzaldehyde                | 1.269 | 1.108 | 12.7  | 125   | 0.00     |
| 5 S  | Phenol-d5                   | 2.197 | 1.798 | 18.2  | 131   | 0.00     |
| 6    | Phenol                      | 2.183 | 1.795 | 17.8  | 130   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.391 | 1.114 | 19.9  | 122   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.487 | 1.250 | 15.9  | 130   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.386 | 1.320 | 4.8   | 142   | 0.00     |
| 10   | 2-Chlorophenol              | 1.350 | 1.265 | 6.3   | 144   | 0.00     |
| 11   | 2-Methylphenol              | 1.541 | 1.249 | 18.9  | 125   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 3.610 | 2.555 | 29.2# | 105   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.836 | 1.493 | 18.7  | 124   | 0.00     |
| 14   | Acetophenone                | 2.618 | 2.193 | 16.2  | 125   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.488 | 1.179 | 20.8# | 116   | 0.00     |
| 16   | 4-Methylphenol              | 1.740 | 1.388 | 20.2# | 122   | 0.00     |
| 17   | Hexachloroethane            | 0.560 | 0.559 | 0.2   | 150#  | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 131   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.156 | 0.164 | -5.1  | 140   | 0.00     |
| 20   | Nitrobenzene                | 0.453 | 0.421 | 7.1   | 124   | 0.00     |
| 21   | Isophorone                  | 0.860 | 0.764 | 11.2  | 117   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.171 | 0.187 | -9.4  | 144   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.173 | 0.182 | -5.2  | 137   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.430 | 0.414 | 3.7   | 125   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.469 | 0.429 | 8.5   | 119   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.355 | 0.375 | -5.6  | 141   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.323 | 0.327 | -1.2  | 136   | 0.00     |
| 28   | Naphthalene                 | 0.991 | 0.969 | 2.2   | 129   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.395 | 0.427 | -8.1  | 127   | 0.00     |
| 30   | 4-Chloroaniline             | 0.383 | 0.391 | -2.1  | 121   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.238 | 0.264 | -10.9 | 150   | 0.00     |
| 32   | Caprolactam                 | 0.134 | 0.120 | 10.4  | 119   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.412 | 0.393 | 4.6   | 124   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.796 | 0.777 | 2.4   | 129   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 120   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.645 | 0.707 | -9.6  | 138   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.356 | 0.224 | 37.1# | 86    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.402 | 0.457 | -13.7 | 139   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.438 | 0.496 | -13.2 | 139   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.461 | 1.451 | 0.7   | 123   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.154 | 1.156 | -0.2  | 122   | 0.00     |
| 42   | 2-Nitroaniline              | 0.484 | 0.462 | 4.5   | 111   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.779 | 1.670 | 6.1   | 112   | 0.00     |
| 44   | Dimethylphthalate           | 1.638 | 1.493 | 8.9   | 111   | 0.00     |

Data Path : Z:\HPCHEM1\BNA\_G\Data\BG082517\  
 Data File : BG028484.D  
 Acq On : 25 Aug 2017 9:12  
 Operator : SJ/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02086

Quant Time: Aug 25 23:04:58 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Aug 25 03:52:53 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.338 | 0.336 | 0.6   | 116   | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.006 | 1.983 | 1.1   | 120   | 0.00     |
| 47   | Acenaphthylene              | 1.917 | 1.892 | 1.3   | 120   | 0.00     |
| 48   | 3-Nitroaniline              | 0.305 | 0.308 | -1.0  | 118   | 0.00     |
| 49 C | Acenaphthene                | 1.293 | 1.292 | 0.1   | 119   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.191 | 0.116 | 39.3# | 86    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.283 | 0.248 | 12.4  | 108   | 0.00     |
| 52   | 4-Nitrophenol               | 0.293 | 0.256 | 12.6  | 105   | 0.00     |
| 53   | Dibenzofuran                | 1.886 | 1.855 | 1.6   | 119   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.512 | 0.480 | 6.3   | 112   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.408 | 0.441 | -8.1  | 131   | 0.00     |
| 56   | Diethylphthalate            | 1.921 | 1.680 | 12.5  | 107   | 0.00     |
| 57 S | Fluorene-d10                | 1.596 | 1.501 | 6.0   | 114   | 0.00     |
| 58   | Fluorene                    | 1.669 | 1.559 | 6.6   | 114   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.896 | 0.850 | 5.1   | 118   | 0.00     |
| 60   | 4-Nitroaniline              | 0.361 | 0.322 | 10.8  | 110   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 107   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.130 | 0.118 | 9.2   | 109   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.132 | 0.114 | 13.6  | 103   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.526 | 0.556 | -5.7  | 114   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.215 | 0.240 | -11.6 | 121   | 0.00     |
| 66   | Hexachlorobenzene           | 0.229 | 0.258 | -12.7 | 125   | 0.00     |
| 67   | Atrazine                    | 0.229 | 0.239 | -4.4  | 113   | 0.00     |
| 68 C | Pentachlorophenol           | 0.116 | 0.109 | 6.0   | 119   | 0.00     |
| 69   | Phenanthrene                | 1.007 | 1.013 | -0.6  | 109   | 0.00     |
| 70 S | Anthracene-d10              | 0.959 | 0.949 | 1.0   | 110   | 0.00     |
| 71   | Anthracene                  | 1.029 | 1.043 | -1.4  | 111   | 0.00     |
| 72   | Carbazole                   | 0.895 | 0.888 | 0.8   | 109   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.114 | 1.093 | 1.9   | 105   | 0.00     |
| 74 C | Fluoranthene                | 1.223 | 1.229 | -0.5  | 108   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 120   | 0.00     |
| 76 S | Pyrene-d10                  | 1.026 | 0.912 | 11.1  | 105   | 0.00     |
| 77   | Pyrene                      | 1.193 | 1.081 | 9.4   | 108   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.447 | 0.417 | 6.7   | 113   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.386 | 0.399 | -3.4  | 121   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.156 | 1.129 | 2.3   | 119   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.636 | 0.599 | 5.8   | 116   | 0.00     |
| 82   | Chrysene                    | 1.041 | 1.014 | 2.6   | 117   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 126   | 0.00     |
| 84   | Di-n-octyl phthalate        | 1.199 | 1.066 | 11.1  | 116   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.197 | 1.133 | 5.3   | 122   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.169 | 1.134 | 3.0   | 121   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.936 | 0.932 | 0.4   | 127   | 0.00     |

Data Path : Z:\HPCHEM1\BNA\_G\Data\BG082517\  
Data File : BG028484.D  
Acq On : 25 Aug 2017 9:12  
Operator : SJ/JU  
Sample : SSTDCCC020  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
LabSampleId :  
SSTD02086

Quant Time: Aug 25 23:04:58 2017  
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG080217.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Aug 25 03:52:53 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               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.159 | 1.132 | 2.3  | 125   | -0.01    |
| 89   | Indeno(1,2,3-cd)pyrene | 1.357 | 1.295 | 4.6  | 125   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.137 | 1.106 | 2.7  | 126   | -0.01    |
| 91   | Benzo(g,h,i)perylene   | 1.116 | 1.011 | 9.4  | 117   | -0.02    |

(#) = Out of Range

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