

Data Path : Z:\HPCHEM1\BNA G\DATA\BG082817\  
 Data File : BG082815.D  
 Acq On : 28 Aug 2017 12:37  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02089

Quant Time: Aug 29 01:27:46 2017  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sat Aug 26 01:54:26 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                 | 0.451 | 0.477 | -5.8   | 140   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.429 | 0.418 | 2.6    | 139   | 0.00     |
| 4    | Benzaldehyde                | 1.269 | 1.242 | 2.1    | 118   | 0.00     |
| 5 S  | Phenol-d5                   | 2.197 | 1.897 | 13.7   | 117   | 0.00     |
| 6    | Phenol                      | 2.183 | 1.884 | 13.7   | 115   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.391 | 1.219 | 12.4   | 112   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.487 | 1.314 | 11.6   | 115   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.386 | 1.358 | 2.0    | 123   | 0.00     |
| 10   | 2-Chlorophenol              | 1.350 | 1.381 | -2.3   | 132   | 0.00     |
| 11   | 2-Methylphenol              | 1.541 | 1.347 | 12.6   | 114   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 3.610 | 2.704 | 25.1#  | 94    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.836 | 1.587 | 13.6   | 111   | 0.00     |
| 14   | Acetophenone                | 2.618 | 2.270 | 13.3   | 109   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.488 | 1.276 | 14.2   | 105   | 0.00     |
| 16   | 4-Methylphenol              | 1.740 | 1.494 | 14.1   | 110   | 0.00     |
| 17   | Hexachloroethane            | 0.560 | 0.542 | 3.2    | 123   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 111   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.156 | 0.167 | -7.1   | 120   | 0.00     |
| 20   | Nitrobenzene                | 0.453 | 0.442 | 2.4    | 110   | 0.00     |
| 21   | Isophorone                  | 0.860 | 0.799 | 7.1    | 103   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.171 | 0.198 | -15.8  | 129   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.173 | 0.201 | -16.2  | 128   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.430 | 0.436 | -1.4   | 111   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.469 | 0.444 | 5.3    | 104   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.355 | 0.387 | -9.0   | 123   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.323 | 0.346 | -7.1   | 122   | 0.00     |
| 28   | Naphthalene                 | 0.991 | 1.029 | -3.8   | 116   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.395 | 0.450 | -13.9  | 113   | 0.00     |
| 30   | 4-Chloroaniline             | 0.383 | 0.423 | -10.4  | 111   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.238 | 0.279 | -17.2  | 134   | 0.00     |
| 32   | Caprolactam                 | 0.134 | 0.123 | 8.2    | 103   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.412 | 0.421 | -2.2   | 112   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.796 | 0.787 | 1.1    | 111   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 104   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.645 | 0.745 | -15.5  | 126   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.356 | 0.434 | -21.9# | 145   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.402 | 0.468 | -16.4  | 123   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.438 | 0.470 | -7.3   | 114   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.461 | 1.523 | -4.2   | 112   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.154 | 1.193 | -3.4   | 109   | 0.00     |
| 42   | 2-Nitroaniline              | 0.484 | 0.471 | 2.7    | 98    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.779 | 1.694 | 4.8    | 99    | 0.00     |
| 44   | Dimethylphthalate           | 1.638 | 1.571 | 4.1    | 101   | 0.00     |

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 Data File : BG028515.D  
 Acq On : 28 Aug 2017 12:37  
 Operator : SJ/JU  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02089

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 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M  
 Quant Title : SVOA CALIBRATION  
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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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.338 | 0.345 | -2.1  | 104   | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.006 | 2.081 | -3.7  | 110   | 0.00     |
| 47   | Acenaphthylene              | 1.917 | 1.933 | -0.8  | 107   | 0.00     |
| 48   | 3-Nitroaniline              | 0.305 | 0.325 | -6.6  | 108   | 0.00     |
| 49 C | Acenaphthene                | 1.293 | 1.286 | 0.5   | 103   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.191 | 0.175 | 8.4   | 113   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.283 | 0.246 | 13.1  | 93    | 0.00     |
| 52   | 4-Nitrophenol               | 0.293 | 0.235 | 19.8  | 84    | 0.00     |
| 53   | Dibenzofuran                | 1.886 | 1.878 | 0.4   | 105   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.512 | 0.524 | -2.3  | 106   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.408 | 0.456 | -11.8 | 118   | 0.00     |
| 56   | Diethylphthalate            | 1.921 | 1.736 | 9.6   | 96    | 0.00     |
| 57 S | Fluorene-d10                | 1.596 | 1.585 | 0.7   | 105   | 0.00     |
| 58   | Fluorene                    | 1.669 | 1.620 | 2.9   | 103   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.896 | 0.901 | -0.6  | 109   | 0.00     |
| 60   | 4-Nitroaniline              | 0.361 | 0.347 | 3.9   | 103   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.130 | 0.117 | 10.0  | 98    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.132 | 0.121 | 8.3   | 99    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.526 | 0.545 | -3.6  | 101   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.215 | 0.233 | -8.4  | 107   | 0.00     |
| 66   | Hexachlorobenzene           | 0.229 | 0.243 | -6.1  | 106   | 0.00     |
| 67   | Atrazine                    | 0.229 | 0.230 | -0.4  | 99    | 0.00     |
| 68 C | Pentachlorophenol           | 0.116 | 0.100 | 13.8  | 99    | 0.00     |
| 69   | Phenanthrene                | 1.007 | 0.997 | 1.0   | 97    | 0.00     |
| 70 S | Anthracene-d10              | 0.959 | 0.971 | -1.3  | 102   | 0.00     |
| 71   | Anthracene                  | 1.029 | 1.051 | -2.1  | 101   | 0.00     |
| 72   | Carbazole                   | 0.895 | 0.873 | 2.5   | 97    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.114 | 1.102 | 1.1   | 96    | 0.00     |
| 74 C | Fluoranthene                | 1.223 | 1.256 | -2.7  | 100   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 76 S | Pyrene-d10                  | 1.026 | 0.941 | 8.3   | 99    | 0.00     |
| 77   | Pyrene                      | 1.193 | 1.092 | 8.5   | 100   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.447 | 0.431 | 3.6   | 107   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.386 | 0.413 | -7.0  | 115   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.156 | 1.147 | 0.8   | 111   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.636 | 0.603 | 5.2   | 107   | 0.00     |
| 82   | Chrysene                    | 1.041 | 1.031 | 1.0   | 109   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 112   | 0.00     |
| 84   | Di-n-octyl phthalate        | 1.199 | 1.107 | 7.7   | 107   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.197 | 1.198 | -0.1  | 115   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.169 | 1.188 | -1.6  | 112   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.936 | 0.951 | -1.6  | 115   | 0.00     |

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Misc :  
ALS Vial : 2 Sample Multiplier: 1

Instrument :  
BNA\_G  
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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.159 | 1.181 | -1.9 | 116   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.357 | 1.268 | 6.6  | 108   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.137 | 1.032 | 9.2  | 105   | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 1.116 | 0.981 | 12.1 | 100   | 0.01     |

(#) = Out of Range

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