

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101116\  
 Data File : BG024264.D  
 Acq On : 11 Oct 2016 12:28  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02015

Quant Time: Oct 12 04:32:26 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Oct 12 04:32:24 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    | 73    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.456 | 0.532 | -16.7  | 83    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.421 | 0.428 | -1.7   | 78    | 0.00     |
| 4    | Benzaldehyde                | 1.313 | 1.447 | -10.2  | 70    | 0.00     |
| 5 S  | Phenol-d5                   | 1.866 | 1.958 | -4.9   | 73    | 0.00     |
| 6    | Phenol                      | 1.894 | 1.989 | -5.0   | 72    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.261 | 1.268 | -0.6   | 71    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.394 | 1.410 | -1.1   | 68    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.285 | 1.375 | -7.0   | 73    | 0.00     |
| 10   | 2-Chlorophenol              | 1.302 | 1.311 | -0.7   | 70    | 0.00     |
| 11   | 2-Methylphenol              | 1.403 | 1.507 | -7.4   | 75    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.784 | 2.887 | -3.7   | 70    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.509 | 1.632 | -8.2   | 75    | 0.00     |
| 14   | Acetophenone                | 2.240 | 2.388 | -6.6   | 72    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.351 | 1.510 | -11.8  | 76    | 0.00     |
| 16   | 4-Methylphenol              | 1.488 | 1.634 | -9.8   | 77    | 0.00     |
| 17   | Hexachloroethane            | 0.602 | 0.626 | -4.0   | 75    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 78    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.147 | -3.5   | 76    | 0.00     |
| 20   | Nitrobenzene                | 0.449 | 0.474 | -5.6   | 76    | 0.00     |
| 21   | Isophorone                  | 0.816 | 0.916 | -12.3  | 78    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.159 | 0.183 | -15.1  | 86    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.173 | 0.189 | -9.2   | 79    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.426 | 0.462 | -8.5   | 78    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.448 | 0.484 | -8.0   | 76    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.346 | 0.373 | -7.8   | 78    | 0.01     |
| 27 C | 2,4-Dichlorophenol          | 0.332 | 0.354 | -6.6   | 77    | 0.00     |
| 28   | Naphthalene                 | 0.963 | 1.044 | -8.4   | 79    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.363 | 0.441 | -21.5# | 85    | 0.00     |
| 30   | 4-Chloroaniline             | 0.369 | 0.424 | -14.9  | 81    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.266 | 0.294 | -10.5  | 80    | 0.00     |
| 32   | Caprolactam                 | 0.126 | 0.149 | -18.3  | 86    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.406 | 0.464 | -14.3  | 84    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.766 | 0.840 | -9.7   | 82    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 84    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.590 | 0.634 | -7.5   | 84    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.432 | 0.428 | 0.9    | 78    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.382 | 0.401 | -5.0   | 83    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.420 | 0.461 | -9.8   | 87    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.295 | 1.397 | -7.9   | 83    | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.015 | 1.086 | -7.0   | 84    | 0.00     |
| 42   | 2-Nitroaniline              | 0.386 | 0.446 | -15.5  | 89    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.396 | 1.580 | -13.2  | 89    | 0.00     |
| 44   | Dimethylphthalate           | 1.369 | 1.527 | -11.5  | 87    | 0.00     |

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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          | 0.266 | 0.311 | -16.9  | 89    | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.600 | 1.795 | -12.2  | 87    | 0.00     |
| 47   | Acenaphthylene              | 1.539 | 1.731 | -12.5  | 88    | 0.00     |
| 48   | 3-Nitroaniline              | 0.251 | 0.302 | -20.3# | 94    | 0.00     |
| 49 C | Acenaphthene                | 1.100 | 1.213 | -10.3  | 87    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.178 | 0.209 | -17.4  | 98    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.248 | 0.281 | -13.3  | 90    | 0.00     |
| 52   | 4-Nitrophenol               | 0.300 | 0.357 | -19.0  | 92    | 0.00     |
| 53   | Dibenzofuran                | 1.619 | 1.870 | -15.5  | 89    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.394 | 0.462 | -17.3  | 91    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.413 | 0.491 | -18.9  | 93    | 0.00     |
| 56   | Diethylphthalate            | 1.443 | 1.640 | -13.7  | 87    | 0.00     |
| 57 S | Fluorene-d10                | 1.320 | 1.499 | -13.6  | 87    | 0.00     |
| 58   | Fluorene                    | 1.339 | 1.506 | -12.5  | 87    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.743 | 0.824 | -10.9  | 87    | 0.00     |
| 60   | 4-Nitroaniline              | 0.289 | 0.323 | -11.8  | 88    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 87    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.120 | 0.126 | -5.0   | 94    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.126 | 0.140 | -11.1  | 93    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.566 | 0.594 | -4.9   | 86    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.232 | 0.256 | -10.3  | 94    | 0.00     |
| 66   | Hexachlorobenzene           | 0.258 | 0.285 | -10.5  | 92    | 0.00     |
| 67   | Atrazine                    | 0.231 | 0.256 | -10.8  | 91    | 0.00     |
| 68 C | Pentachlorophenol           | 0.157 | 0.174 | -10.8  | 93    | 0.00     |
| 69   | Phenanthrene                | 1.010 | 1.123 | -11.2  | 91    | 0.00     |
| 70 S | Anthracene-d10              | 0.903 | 0.979 | -8.4   | 88    | 0.00     |
| 71   | Anthracene                  | 1.028 | 1.140 | -10.9  | 90    | 0.00     |
| 72   | Carbazole                   | 0.807 | 0.957 | -18.6  | 88    | 0.00     |
| 73   | Di-n-butylphthalate         | 0.986 | 1.113 | -12.9  | 90    | 0.00     |
| 74 C | Fluoranthene                | 1.020 | 1.298 | -27.3# | 91    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 90    | 0.00     |
| 76 S | Pyrene-d10                  | 0.933 | 1.008 | -8.0   | 89    | 0.00     |
| 77   | Pyrene                      | 1.068 | 1.166 | -9.2   | 90    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.429 | 0.450 | -4.9   | 89    | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.372 | 0.419 | -12.6  | 92    | 0.00     |
| 80   | Benzo(a)anthracene          | 1.110 | 1.213 | -9.3   | 92    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.607 | 0.619 | -2.0   | 85    | 0.01     |
| 82   | Chrysene                    | 1.057 | 1.133 | -7.2   | 89    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 94    | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.938 | 1.050 | -11.9  | 91    | 0.01     |
| 85   | Benzo(b)fluoranthene        | 1.110 | 1.144 | -3.1   | 91    | 0.01     |
| 86   | Benzo(k)fluoranthene        | 1.070 | 1.156 | -8.0   | 95    | 0.01     |
| 87 S | Benzo(a)pyrene-d12          | 0.900 | 0.954 | -6.0   | 93    | 0.00     |

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|------|------------------------|-------|-------|-------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.058 | 1.115 | -5.4  | 92    | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.253 | 1.380 | -10.1 | 97    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.059 | 1.140 | -7.6  | 96    | 0.01     |
| 91   | Benzo(a,h,i)perylene   | 1.053 | 1.123 | -6.6  | 95    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 1