

Data Path : Z:\HPCHEM1\BNA G\DATA\BG121616\  
 Data File : BG025267.D  
 Acq On : 17 Dec 2016 9:46  
 Operator : UM/SJ  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02015

Quant Time: Dec 18 02:05:47 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Dec 18 00:39:32 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    | 71    | 0.01     |
| 2    | 1,4-Dioxane                 | 0.504 | 0.516 | -2.4   | 75    | 0.02     |
| 3 S  | 1,4-Dioxane-d8              | 0.387 | 0.423 | -9.3   | 79    | 0.01     |
| 4    | Benzaldehyde                | 1.299 | 1.512 | -16.4  | 79    | 0.01     |
| 5 S  | Phenol-d5                   | 1.725 | 1.892 | -9.7   | 78    | 0.01     |
| 6    | Phenol                      | 1.771 | 2.014 | -13.7  | 79    | 0.01     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.067 | 1.171 | -9.7   | 77    | 0.01     |
| 8    | Bis(2-Chloroethyl)ether     | 1.296 | 1.419 | -9.5   | 80    | 0.01     |
| 9 S  | 2-Chlorophenol-d4           | 1.207 | 1.384 | -14.7  | 81    | 0.01     |
| 10   | 2-Chlorophenol              | 1.213 | 1.415 | -16.7  | 82    | 0.02     |
| 11   | 2-Methylphenol              | 1.304 | 1.455 | -11.6  | 80    | 0.01     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.204 | 2.584 | -17.2  | 82    | 0.02     |
| 13 S | 4-Methylphenol-d8           | 1.442 | 1.605 | -11.3  | 81    | 0.01     |
| 14   | Acetophenone                | 2.194 | 2.341 | -6.7   | 77    | 0.01     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.267 | 1.382 | -9.1   | 76    | 0.00     |
| 16   | 4-Methylphenol              | 1.451 | 1.558 | -7.4   | 81    | 0.01     |
| 17   | Hexachloroethane            | 0.537 | 0.665 | -23.8# | 86    | 0.01     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 69    | 0.02     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.161 | -13.4  | 79    | 0.00     |
| 20   | Nitrobenzene                | 0.435 | 0.500 | -14.9  | 82    | 0.01     |
| 21   | Isophorone                  | 0.824 | 0.894 | -8.5   | 74    | 0.01     |
| 22 S | 2-Nitrophenol-d4            | 0.177 | 0.193 | -9.0   | 77    | 0.01     |
| 23 C | 2-Nitrophenol               | 0.179 | 0.203 | -13.4  | 79    | 0.02     |
| 24   | 2,4-Dimethylphenol          | 0.410 | 0.460 | -12.2  | 75    | 0.01     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.420 | 0.464 | -10.5  | 75    | 0.01     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.340 | 0.376 | -10.6  | 76    | 0.02     |
| 27 C | 2,4-Dichlorophenol          | 0.330 | 0.381 | -15.5  | 79    | 0.00     |
| 28   | Naphthalene                 | 0.930 | 1.024 | -10.1  | 76    | 0.01     |
| 29 S | 4-Chloroaniline-d4          | 0.334 | 0.432 | -29.3# | 89    | 0.01     |
| 30   | 4-Chloroaniline             | 0.341 | 0.435 | -27.6# | 86    | 0.02     |
| 31 C | Hexachlorobutadiene         | 0.283 | 0.324 | -14.5  | 78    | 0.01     |
| 32   | Caprolactam                 | 0.137 | 0.126 | 8.0    | 65    | 0.01     |
| 33 C | 4-Chloro-3-methylphenol     | 0.406 | 0.445 | -9.6   | 75    | 0.01     |
| 34   | 2-Methylnaphthalene         | 0.734 | 0.788 | -7.4   | 72    | 0.02     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 63    | 0.01     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.603 | 0.721 | -19.6  | 77    | 0.01     |
| 37   | Hexachlorocyclopentadiene   | 0.354 | 0.205 | 42.1#  | 40    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.379 | 0.469 | -23.7  | 79    | 0.01     |
| 39   | 2,4,5-Trichlorophenol       | 0.415 | 0.495 | -19.3  | 74    | 0.01     |
| 40   | 1,1'-Biphenyl               | 1.213 | 1.408 | -16.1  | 72    | 0.01     |
| 41   | 2-Chloronaphthalene         | 0.965 | 1.105 | -14.5  | 74    | 0.00     |
| 42   | 2-Nitroaniline              | 0.378 | 0.476 | -25.9# | 76    | 0.01     |
| 43 S | Dimethylphthalate-d6        | 1.376 | 1.497 | -8.8   | 66    | 0.00     |
| 44   | Dimethylphthalate           | 1.352 | 1.490 | -10.2  | 69    | 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.291 | 0.317 | -8.9   | 67    | 0.01     |
| 46 S | Acenaphthylene-d8           | 1.507 | 1.721 | -14.2  | 71    | 0.01     |
| 47   | Acenaphthylene              | 1.465 | 1.704 | -16.3  | 71    | 0.00     |
| 48   | 3-Nitroaniline              | 0.252 | 0.302 | -19.8  | 73    | 0.00     |
| 49 C | Acenaphthene                | 1.004 | 1.130 | -12.5  | 71    | 0.01     |
| 50   | 2,4-Dinitrophenol           | 0.189 | 0.120 | 36.5#  | 45    | 0.01     |
| 51 S | 4-Nitrophenol-d4            | 0.236 | 0.278 | -17.8  | 73    | 0.01     |
| 52   | 4-Nitrophenol               | 0.284 | 0.337 | -18.7  | 76    | 0.01     |
| 53   | Dibenzofuran                | 1.587 | 1.782 | -12.3  | 70    | 0.01     |
| 54   | 2,4-Dinitrotoluene          | 0.438 | 0.474 | -8.2   | 67    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.439 | 0.514 | -17.1  | 73    | 0.01     |
| 56   | Diethylphthalate            | 1.432 | 1.470 | -2.7   | 63    | 0.01     |
| 57 S | Fluorene-d10                | 1.295 | 1.445 | -11.6  | 69    | 0.01     |
| 58   | Fluorene                    | 1.292 | 1.428 | -10.5  | 68    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.728 | 0.814 | -11.8  | 70    | 0.00     |
| 60   | 4-Nitroaniline              | 0.303 | 0.341 | -12.5  | 71    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 60    | 0.01     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.124 | 0.141 | -13.7  | 71    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.128 | 0.154 | -20.3# | 74    | 0.01     |
| 64   | N-Nitrosodiphenylamine      | 0.515 | 0.600 | -16.5  | 68    | 0.01     |
| 65   | 4-Bromophenyl-phenylether   | 0.226 | 0.272 | -20.4# | 72    | 0.01     |
| 66   | Hexachlorobenzene           | 0.254 | 0.292 | -15.0  | 67    | 0.01     |
| 67   | Atrazine                    | 0.245 | 0.266 | -8.6   | 63    | 0.01     |
| 68 C | Pentachlorophenol           | 0.153 | 0.165 | -7.8   | 67    | 0.01     |
| 69   | Phenanthrene                | 0.955 | 1.099 | -15.1  | 65    | 0.01     |
| 70 S | Anthracene-d10              | 0.844 | 0.962 | -14.0  | 66    | 0.00     |
| 71   | Anthracene                  | 0.981 | 1.142 | -16.4  | 66    | 0.01     |
| 72   | Carbazole                   | 0.819 | 1.006 | -22.8# | 68    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.034 | 1.167 | -12.9  | 64    | 0.00     |
| 74 C | Fluoranthene                | 1.134 | 1.408 | -24.2  | 65    | 0.01     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 60    | 0.01     |
| 76 S | Pyrene-d10                  | 0.824 | 0.912 | -10.7  | 65    | 0.00     |
| 77   | Pyrene                      | 0.961 | 1.059 | -10.2  | 63    | 0.01     |
| 78   | Butylbenzylphthalate        | 0.376 | 0.434 | -15.4  | 66    | 0.01     |
| 79   | 3,3'-Dichlorobenzidine      | 0.349 | 0.470 | -34.7# | 77    | 0.01     |
| 80   | Benzo(a)anthracene          | 1.039 | 1.207 | -16.2  | 67    | 0.01     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.518 | 0.578 | -11.6  | 65    | 0.01     |
| 82   | Chrysene                    | 0.972 | 1.132 | -16.5  | 68    | 0.01     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 59    | 0.02     |
| 84   | Di-n-octyl phthalate        | 0.812 | 0.976 | -20.2# | 66    | 0.02     |
| 85   | Benzo(b)fluoranthene        | 0.992 | 1.123 | -13.2  | 65    | 0.01     |
| 86   | Benzo(k)fluoranthene        | 0.943 | 1.122 | -19.0  | 69    | 0.02     |
| 87 S | Benzo(a)pyrene-d12          | 0.810 | 0.930 | -14.8  | 66    | 0.02     |

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|      | Compound               | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|------------------------|-------|-------|-------|-------|----------|
| 88 C | Benzo(a)pyrene         | 0.954 | 1.100 | -15.3 | 67    | 0.02     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.185 | 1.366 | -15.3 | 67    | 0.04     |
| 90   | Dibenzo(a,h)anthracene | 0.998 | 1.152 | -15.4 | 68    | 0.02     |
| 91   | Benzo(a,h,i)perylene   | 0.989 | 1.050 | -6.2  | 62    | 0.04     |

(#) = Out of Range

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