

Data Path : Z:\HPCHEM1\BNA G\DATA\BG123115\  
 Data File : BG020652.D  
 Acq On : 31 Dec 2015 21:49  
 Operator : UM/SJ  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02011

Quant Time: Jan 01 04:56:53 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG123015.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 01 03:15:46 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    | 128   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.422 | 0.476 | -12.8  | 143   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.334 | 0.402 | -20.4# | 129   | 0.00     |
| 4    | Benzaldehyde                | 1.035 | 1.182 | -14.2  | 125   | 0.00     |
| 5 S  | Phenol-d5                   | 1.689 | 1.716 | -1.6   | 123   | 0.00     |
| 6    | Phenol                      | 1.820 | 1.816 | 0.2    | 119   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.022 | 1.076 | -5.3   | 126   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.298 | 1.385 | -6.7   | 126   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.292 | 1.402 | -8.5   | 128   | 0.00     |
| 10   | 2-Chlorophenol              | 1.350 | 1.432 | -6.1   | 127   | 0.00     |
| 11   | 2-Methylphenol              | 1.402 | 1.435 | -2.4   | 124   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.697 | 1.835 | -8.1   | 126   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.449 | 1.465 | -1.1   | 121   | 0.00     |
| 14   | Acetophenone                | 2.375 | 2.443 | -2.9   | 122   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.199 | 1.256 | -4.8   | 121   | 0.00     |
| 16   | 4-Methylphenol              | 1.551 | 1.558 | -0.5   | 120   | 0.00     |
| 17   | Hexachloroethane            | 0.567 | 0.603 | -6.3   | 127   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 120   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.155 | 0.172 | -11.0  | 124   | 0.00     |
| 20   | Nitrobenzene                | 0.386 | 0.426 | -10.4  | 123   | 0.00     |
| 21   | Isophorone                  | 0.751 | 0.817 | -8.8   | 118   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.179 | 0.190 | -6.1   | 113   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.189 | 0.203 | -7.4   | 122   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.398 | 0.440 | -10.6  | 121   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.415 | 0.454 | -9.4   | 120   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.341 | 0.373 | -9.4   | 118   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.352 | 0.387 | -9.9   | 121   | 0.00     |
| 28   | Naphthalene                 | 1.059 | 1.133 | -7.0   | 118   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.405 | 0.455 | -12.3  | 119   | 0.00     |
| 30   | 4-Chloroaniline             | 0.416 | 0.465 | -11.8  | 117   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.284 | 0.309 | -8.8   | 123   | 0.00     |
| 32   | Caprolactam                 | 0.121 | 0.120 | 0.8    | 113   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.381 | 0.415 | -8.9   | 119   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.793 | 0.845 | -6.6   | 118   | 0.00     |
| 35   | 1-Methylnaphthalene         | 0.765 | 0.804 | -5.1   | 116   | 0.00     |
| 36 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 112   | 0.00     |
| 37   | 1,2,4,5-Tetrachlorobenzene  | 0.770 | 0.851 | -10.5  | 117   | 0.00     |
| 38   | Hexachlorocyclopentadiene   | 0.277 | 0.201 | 27.4#  | 99    | 0.00     |
| 39 C | 2,4,6-Trichlorophenol       | 0.431 | 0.460 | -6.7   | 110   | 0.00     |
| 40   | 2,4,5-Trichlorophenol       | 0.473 | 0.529 | -11.8  | 114   | 0.00     |
| 41   | 1,1'-Biphenyl               | 1.505 | 1.654 | -9.9   | 116   | 0.00     |
| 42   | 2-Chloronaphthalene         | 1.211 | 1.329 | -9.7   | 116   | 0.00     |
| 43   | 2-Nitroaniline              | 0.348 | 0.387 | -11.2  | 113   | 0.00     |
| 44 S | Dimethylphthalate-d6        | 1.637 | 1.741 | -6.4   | 111   | 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   | Dimethylphthalate           | 1.672 | 1.783 | -6.6  | 112   | 0.00     |
| 46   | 2,6-Dinitrotoluene          | 0.347 | 0.371 | -6.9  | 111   | 0.00     |
| 47 S | Acenaphthylene-d8           | 1.859 | 2.042 | -9.8  | 115   | 0.00     |
| 48   | Acenaphthylene              | 1.980 | 2.158 | -9.0  | 115   | 0.00     |
| 49   | 3-Nitroaniline              | 0.316 | 0.363 | -14.9 | 117   | 0.00     |
| 50 C | Acenaphthene                | 1.340 | 1.446 | -7.9  | 114   | 0.00     |
| 51   | 2,4-Dinitrophenol           | 0.146 | 0.111 | 24.0# | 100   | 0.00     |
| 52 S | 4-Nitrophenol-d4            | 0.257 | 0.260 | -1.2  | 107   | 0.00     |
| 53   | 4-Nitrophenol               | 0.227 | 0.240 | -5.7  | 114   | 0.00     |
| 54   | Dibenzofuran                | 2.025 | 2.152 | -6.3  | 112   | 0.00     |
| 55   | 2,4-Dinitrotoluene          | 0.506 | 0.546 | -7.9  | 110   | 0.00     |
| 56   | 2,3,4,6-Tetrachlorophenol   | 0.441 | 0.468 | -6.1  | 107   | 0.00     |
| 57   | Diethylphthalate            | 1.723 | 1.830 | -6.2  | 112   | 0.00     |
| 58 S | Fluorene-d10                | 1.484 | 1.566 | -5.5  | 112   | 0.00     |
| 59   | Fluorene                    | 1.629 | 1.729 | -6.1  | 111   | 0.00     |
| 60   | 4-Chlorophenyl-phenylether  | 0.941 | 1.002 | -6.5  | 111   | 0.00     |
| 61   | 4-Nitroaniline              | 0.345 | 0.390 | -13.0 | 117   | 0.00     |
| 62 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 110   | 0.00     |
| 63 S | 4,6-Dinitro-2-methylphenol- | 0.121 | 0.120 | 0.8   | 110   | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol  | 0.130 | 0.128 | 1.5   | 106   | 0.00     |
| 65   | N-Nitrosodiphenylamine      | 0.551 | 0.604 | -9.6  | 112   | 0.00     |
| 66   | 4-Bromophenyl-phenylether   | 0.243 | 0.259 | -6.6  | 110   | 0.00     |
| 67   | Hexachlorobenzene           | 0.267 | 0.280 | -4.9  | 109   | 0.00     |
| 68   | Atrazine                    | 0.243 | 0.265 | -9.1  | 111   | 0.00     |
| 69 C | Pentachlorophenol           | 0.107 | 0.096 | 10.3  | 103   | 0.00     |
| 70   | Phenanthrene                | 1.087 | 1.167 | -7.4  | 110   | 0.00     |
| 71 S | Anthracene-d10              | 0.932 | 0.997 | -7.0  | 110   | 0.00     |
| 72   | Anthracene                  | 1.112 | 1.189 | -6.9  | 110   | 0.00     |
| 73   | 1,2,3,4-Tetrachlorobenzene  | 0.318 | 0.359 | -12.9 | 120   | 0.00     |
| 74   | Pentachlorobenzene          | 0.303 | 0.329 | -8.6  | 112   | 0.00     |
| 75   | Carbazole                   | 0.938 | 1.020 | -8.7  | 109   | 0.00     |
| 76   | Di-n-butylphthalate         | 1.101 | 1.212 | -10.1 | 111   | 0.00     |
| 77 C | Fluoranthene                | 1.328 | 1.499 | -12.9 | 111   | 0.00     |
| 78 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 79 S | Pyrene-d10                  | 0.897 | 0.970 | -8.1  | 110   | 0.00     |
| 80   | Pyrene                      | 1.158 | 1.252 | -8.1  | 110   | 0.00     |
| 81   | Butylbenzylphthalate        | 0.408 | 0.451 | -10.5 | 114   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine      | 0.381 | 0.426 | -11.8 | 113   | 0.00     |
| 83   | Benzo(a)anthracene          | 1.170 | 1.268 | -8.4  | 112   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate  | 0.586 | 0.642 | -9.6  | 113   | 0.00     |
| 85   | Chrysene                    | 1.101 | 1.177 | -6.9  | 111   | 0.00     |
| 86 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 110   | 0.00     |
| 87   | Di-n-octyl phthalate        | 1.050 | 1.160 | -10.5 | 115   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.188 | 1.297 | -9.2 | 111   | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.138 | 1.238 | -8.8 | 112   | 0.00     |
| 90 S | Benzo(a)pyrene-d12     | 0.998 | 1.072 | -7.4 | 111   | 0.00     |
| 91 C | Benzo(a)pyrene         | 1.141 | 1.239 | -8.6 | 112   | 0.00     |
| 92   | Indeno(1,2,3-cd)pyrene | 1.361 | 1.453 | -6.8 | 110   | 0.00     |
| 93   | Dibenzo(a,h)anthracene | 1.135 | 1.217 | -7.2 | 110   | 0.00     |
| 94   | Benzo(a,h,i)perylene   | 1.118 | 1.183 | -5.8 | 109   | 0.00     |

(#) = Out of Range

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