

Data Path : Z:\HPCHEM1\BNA G\Data\BG111715\  
 Data File : BG019659.D  
 Acq On : 17 Nov 2015 9:19  
 Operator : UM/NP  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02029

Quant Time: Nov 18 00:32:40 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG111615.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Nov 17 01:14:52 2015  
 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   | 86    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.534 | 0.613 | -14.8 | 127   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.468 | 0.444 | 5.1   | 90    | 0.00     |
| 4    | Benzaldehyde                | 1.493 | 1.644 | -10.1 | 85    | 0.00     |
| 5 S  | Phenol-d5                   | 2.049 | 1.994 | 2.7   | 90    | 0.00     |
| 6    | Phenol                      | 2.212 | 2.262 | -2.3  | 94    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.342 | 1.369 | -2.0  | 89    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.540 | 1.543 | -0.2  | 87    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.314 | 1.308 | 0.5   | 91    | 0.00     |
| 10   | 2-Chlorophenol              | 1.284 | 1.231 | 4.1   | 90    | 0.00     |
| 11   | 2-Methylphenol              | 1.629 | 1.569 | 3.7   | 86    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.345 | 1.311 | 2.5   | 85    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.696 | 1.635 | 3.6   | 87    | 0.00     |
| 14   | Acetophenone                | 2.803 | 2.840 | -1.3  | 91    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.894 | 1.781 | 6.0   | 87    | 0.00     |
| 16   | 4-Methylphenol              | 1.835 | 1.790 | 2.5   | 89    | 0.00     |
| 17   | Hexachloroethane            | 0.610 | 0.600 | 1.6   | 91    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 88    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.164 | 0.154 | 6.1   | 88    | 0.00     |
| 20   | Nitrobenzene                | 0.591 | 0.566 | 4.2   | 88    | 0.00     |
| 21   | Isophorone                  | 1.141 | 1.095 | 4.0   | 89    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.185 | 0.169 | 8.6   | 79    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.200 | 0.190 | 5.0   | 89    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.533 | 0.538 | -0.9  | 95    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.565 | 0.530 | 6.2   | 87    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.400 | 0.392 | 2.0   | 91    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.381 | 0.361 | 5.2   | 88    | 0.00     |
| 28   | Naphthalene                 | 1.080 | 1.032 | 4.4   | 89    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.379 | 0.407 | -7.4  | 88    | 0.00     |
| 30   | 4-Chloroaniline             | 0.395 | 0.419 | -6.1  | 86    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.343 | 0.342 | 0.3   | 91    | 0.00     |
| 32   | Caprolactam                 | 0.155 | 0.147 | 5.2   | 85    | -0.01    |
| 33 C | 4-Chloro-3-methylphenol     | 0.531 | 0.514 | 3.2   | 91    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.862 | 0.846 | 1.9   | 88    | 0.00     |
| 35   | 1-Methylnaphthalene         | 0.858 | 0.829 | 3.4   | 91    | 0.00     |
| 36 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 93    | 0.00     |
| 37   | 1,2,4,5-Tetrachlorobenzene  | 0.802 | 0.755 | 5.9   | 91    | 0.00     |
| 38   | Hexachlorocyclopentadiene   | 0.467 | 0.418 | 10.5  | 88    | 0.00     |
| 39 C | 2,4,6-Trichlorophenol       | 0.489 | 0.447 | 8.6   | 89    | 0.00     |
| 40   | 2,4,5-Trichlorophenol       | 0.512 | 0.511 | 0.2   | 97    | 0.00     |
| 41   | 1,1'-Biphenyl               | 1.488 | 1.407 | 5.4   | 92    | 0.00     |
| 42   | 2-Chloronaphthalene         | 1.156 | 1.068 | 7.6   | 88    | 0.00     |
| 43   | 2-Nitroaniline              | 0.489 | 0.457 | 6.5   | 88    | 0.00     |
| 44 S | Dimethylphthalate-d6        | 1.809 | 1.698 | 6.1   | 90    | 0.00     |

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 ALS Vial : 2 Sample Multiplier: 1

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 LabSampleId :  
 SSTD02029

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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   | Dimethylphthalate           | 1.794 | 1.725 | 3.8  | 91    | 0.00     |
| 46   | 2,6-Dinitrotoluene          | 0.359 | 0.350 | 2.5  | 95    | 0.00     |
| 47 S | Acenaphthylene-d8           | 1.961 | 1.840 | 6.2  | 90    | 0.00     |
| 48   | Acenaphthylene              | 1.957 | 1.855 | 5.2  | 90    | 0.00     |
| 49   | 3-Nitroaniline              | 0.299 | 0.309 | -3.3 | 87    | 0.00     |
| 50 C | Acenaphthene                | 1.340 | 1.277 | 4.7  | 92    | 0.00     |
| 51   | 2,4-Dinitrophenol           | 0.236 | 0.200 | 15.3 | 85    | 0.00     |
| 52 S | 4-Nitrophenol-d4            | 0.285 | 0.283 | 0.7  | 93    | 0.00     |
| 53   | 4-Nitrophenol               | 0.347 | 0.350 | -0.9 | 93    | 0.00     |
| 54   | Dibenzofuran                | 1.964 | 1.897 | 3.4  | 94    | 0.00     |
| 55   | 2,4-Dinitrotoluene          | 0.558 | 0.541 | 3.0  | 91    | 0.00     |
| 56   | 2,3,4,6-Tetrachlorophenol   | 0.543 | 0.518 | 4.6  | 90    | 0.00     |
| 57   | Diethylphthalate            | 1.792 | 1.758 | 1.9  | 94    | 0.00     |
| 58 S | Fluorene-d10                | 1.619 | 1.518 | 6.2  | 90    | 0.00     |
| 59   | Fluorene                    | 1.717 | 1.624 | 5.4  | 93    | 0.00     |
| 60   | 4-Chlorophenyl-phenylether  | 0.999 | 0.943 | 5.6  | 92    | 0.00     |
| 61   | 4-Nitroaniline              | 0.360 | 0.380 | -5.6 | 91    | 0.00     |
| 62 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0  | 92    | 0.00     |
| 63 S | 4,6-Dinitro-2-methylphenol- | 0.129 | 0.118 | 8.5  | 87    | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol  | 0.137 | 0.126 | 8.0  | 85    | 0.00     |
| 65   | N-Nitrosodiphenylamine      | 0.539 | 0.517 | 4.1  | 92    | 0.00     |
| 66   | 4-Bromophenyl-phenylether   | 0.240 | 0.225 | 6.2  | 89    | 0.00     |
| 67   | Hexachlorobenzene           | 0.254 | 0.242 | 4.7  | 90    | 0.00     |
| 68   | Atrazine                    | 0.251 | 0.251 | 0.0  | 92    | 0.00     |
| 69 C | Pentachlorophenol           | 0.142 | 0.133 | 6.3  | 92    | 0.00     |
| 70   | Phenanthrene                | 1.070 | 1.051 | 1.8  | 94    | 0.00     |
| 71 S | Anthracene-d10              | 0.971 | 0.946 | 2.6  | 92    | 0.00     |
| 72   | Anthracene                  | 1.079 | 1.056 | 2.1  | 93    | 0.00     |
| 73   | 1,2,3,4-Tetrachlorobenzene  | 0.276 | 0.269 | 2.5  | 93    | 0.00     |
| 74   | Pentachlorobenzene          | 0.288 | 0.280 | 2.8  | 92    | 0.00     |
| 75   | Carbazole                   | 0.867 | 0.897 | -3.5 | 92    | 0.00     |
| 76   | Di-n-butylphthalate         | 1.124 | 1.090 | 3.0  | 90    | 0.00     |
| 77 C | Fluoranthene                | 1.364 | 1.455 | -6.7 | 92    | 0.00     |
| 78 I | Chrysene-d12                | 1.000 | 1.000 | 0.0  | 91    | 0.00     |
| 79 S | Pyrene-d10                  | 0.911 | 0.898 | 1.4  | 94    | 0.00     |
| 80   | Pyrene                      | 1.168 | 1.135 | 2.8  | 92    | 0.00     |
| 81   | Butylbenzylphthalate        | 0.384 | 0.379 | 1.3  | 91    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine      | 0.366 | 0.388 | -6.0 | 89    | 0.00     |
| 83   | Benzo(a)anthracene          | 1.212 | 1.163 | 4.0  | 89    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate  | 0.563 | 0.541 | 3.9  | 89    | 0.00     |
| 85   | Chrysene                    | 1.141 | 1.091 | 4.4  | 90    | 0.00     |
| 86 I | Perylene-d12                | 1.000 | 1.000 | 0.0  | 90    | 0.00     |
| 87   | Di-n-octyl phthalate        | 0.971 | 0.994 | -2.4 | 90    | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.232 | 1.150 | 6.7  | 86    | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.189 | 1.182 | 0.6  | 93    | 0.00     |
| 90 S | Benzo(a)pyrene-d12     | 1.044 | 1.013 | 3.0  | 90    | -0.01    |
| 91 C | Benzo(a)pyrene         | 1.186 | 1.152 | 2.9  | 91    | 0.00     |
| 92   | Indeno(1,2,3-cd)pyrene | 1.328 | 1.273 | 4.1  | 90    | -0.02    |
| 93   | Dibenzo(a,h)anthracene | 1.091 | 1.054 | 3.4  | 89    | -0.01    |
| 94   | Benzo(a,h,i)perylene   | 1.102 | 1.066 | 3.3  | 90    | -0.02    |

(#) = Out of Range

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