

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA\_G\Data\BG012218\  
 Data File : BG032223.D  
 Acq On : 22 Jan 2018 22:38  
 Operator : SJ/JU  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02083

Quant Time: Jan 23 07:04:37 2018  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG012218.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 22 17:12:25 2018  
 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 91    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.375 | 0.342 | 8.8    | 86    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.330 | 0.315 | 4.5    | 80    | 0.00     |
| 4    | Benzaldehyde                | 0.684 | 0.858 | -25.4# | 90    | 0.00     |
| 5 S  | Phenol-d5                   | 1.374 | 1.421 | -3.4   | 93    | 0.00     |
| 6    | Phenol                      | 1.347 | 1.386 | -2.9   | 89    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.648 | 0.651 | -0.5   | 87    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 0.947 | 1.007 | -6.3   | 95    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.282 | 1.350 | -5.3   | 93    | 0.00     |
| 10   | 2-Chlorophenol              | 1.211 | 1.250 | -3.2   | 93    | 0.00     |
| 11   | 2-Methylphenol              | 1.072 | 1.101 | -2.7   | 93    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 0.943 | 0.959 | -1.7   | 89    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.248 | 1.279 | -2.5   | 92    | 0.00     |
| 14   | Acetophenone                | 1.748 | 1.824 | -4.3   | 92    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 0.816 | 0.831 | -1.8   | 90    | 0.00     |
| 16   | 4-Methylphenol              | 1.173 | 1.221 | -4.1   | 92    | 0.00     |
| 17   | Hexachloroethane            | 0.478 | 0.483 | -1.0   | 93    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 92    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.146 | 0.139 | 4.8    | 90    | 0.00     |
| 20   | Nitrobenzene                | 0.293 | 0.293 | 0.0    | 94    | 0.00     |
| 21   | Isophorone                  | 0.528 | 0.526 | 0.4    | 90    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.175 | 0.184 | -5.1   | 95    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.179 | 0.186 | -3.9   | 93    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.359 | 0.355 | 1.1    | 91    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.335 | 0.328 | 2.1    | 92    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.383 | 0.388 | -1.3   | 94    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.361 | 0.358 | 0.8    | 90    | 0.00     |
| 28   | Naphthalene                 | 0.912 | 0.933 | -2.3   | 95    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.267 | 0.314 | -17.6  | 88    | 0.00     |
| 30   | 4-Chloroaniline             | 0.259 | 0.319 | -23.2# | 92    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.326 | 0.324 | 0.6    | 92    | 0.00     |
| 32   | Caprolactam                 | 0.093 | 0.101 | -8.6   | 101   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.327 | 0.335 | -2.4   | 95    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.776 | 0.793 | -2.2   | 94    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 96    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.892 | 0.888 | 0.4    | 95    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.385 | 0.330 | 14.3   | 92    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.515 | 0.510 | 1.0    | 95    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.532 | 0.535 | -0.6   | 96    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.578 | 1.555 | 1.5    | 94    | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.266 | 1.253 | 1.0    | 94    | 0.00     |
| 42   | 2-Nitroaniline              | 0.253 | 0.254 | -0.4   | 95    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.670 | 1.655 | 0.9    | 94    | 0.00     |
| 44   | Dimethylphthalate           | 1.603 | 1.622 | -1.2   | 95    | 0.00     |

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|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.327 | 0.348 | -6.4  | 99    | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.047 | 2.047 | 0.0   | 94    | 0.00     |
| 47   | Acenaphthylene              | 1.785 | 1.790 | -0.3  | 97    | 0.00     |
| 48   | 3-Nitroaniline              | 0.239 | 0.259 | -8.4  | 96    | 0.00     |
| 49 C | Acenaphthene                | 1.303 | 1.308 | -0.4  | 95    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.169 | 0.121 | 28.4# | 100   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.223 | 0.214 | 4.0   | 94    | 0.00     |
| 52   | 4-Nitrophenol               | 0.205 | 0.201 | 2.0   | 94    | 0.00     |
| 53   | Dibenzofuran                | 1.962 | 1.959 | 0.2   | 95    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.473 | 0.478 | -1.1  | 95    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.509 | 0.520 | -2.2  | 94    | 0.00     |
| 56   | Diethylphthalate            | 1.517 | 1.537 | -1.3  | 95    | 0.00     |
| 57 S | Fluorene-d10                | 1.555 | 1.589 | -2.2  | 97    | 0.00     |
| 58   | Fluorene                    | 1.556 | 1.578 | -1.4  | 95    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.982 | 0.993 | -1.1  | 96    | 0.00     |
| 60   | 4-Nitroaniline              | 0.275 | 0.287 | -4.4  | 99    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.123 | 0.125 | -1.6  | 109   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.129 | 0.126 | 2.3   | 103   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.569 | 0.564 | 0.9   | 97    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.288 | 0.281 | 2.4   | 94    | 0.00     |
| 66   | Hexachlorobenzene           | 0.299 | 0.287 | 4.0   | 94    | 0.00     |
| 67   | Atrazine                    | 0.256 | 0.257 | -0.4  | 93    | 0.00     |
| 68 C | Pentachlorophenol           | 0.149 | 0.132 | 11.4  | 92    | 0.00     |
| 69   | Phenanthrene                | 1.052 | 1.056 | -0.4  | 96    | 0.00     |
| 70 S | Anthracene-d10              | 0.962 | 0.967 | -0.5  | 97    | 0.00     |
| 71   | Anthracene                  | 1.080 | 1.088 | -0.7  | 96    | 0.00     |
| 72   | 1,2,3,4-Tetrachlorobenzene  | 0.381 | 0.369 | 3.1   | 94    | 0.00     |
| 73   | Pentachlorobenzene          | 0.400 | 0.385 | 3.8   | 94    | 0.00     |
| 74   | Carbazole                   | 0.837 | 0.870 | -3.9  | 98    | 0.00     |
| 75   | Di-n-butylphthalate         | 0.987 | 1.002 | -1.5  | 96    | 0.00     |
| 76 C | Fluoranthene                | 1.298 | 1.418 | -9.2  | 98    | 0.00     |
| 77 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 99    | 0.00     |
| 78 S | Pyrene-d10                  | 0.928 | 0.923 | 0.5   | 97    | 0.00     |
| 79   | Pyrene                      | 1.129 | 1.120 | 0.8   | 96    | 0.00     |
| 80   | Butylbenzylphthalate        | 0.368 | 0.363 | 1.4   | 97    | 0.00     |
| 81   | 3,3'-Dichlorobenzidine      | 0.320 | 0.327 | -2.2  | 99    | 0.00     |
| 82   | Benzo(a)anthracene          | 1.209 | 1.220 | -0.9  | 99    | 0.00     |
| 83   | Bis(2-ethylhexyl)phthalate  | 0.522 | 0.519 | 0.6   | 98    | 0.00     |
| 84   | Chrysene                    | 1.097 | 1.107 | -0.9  | 99    | 0.00     |
| 85 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 100   | 0.00     |
| 86   | Di-n-octyl phthalate        | 0.834 | 0.830 | 0.5   | 97    | 0.00     |
| 87   | Benzo(b)fluoranthene        | 1.182 | 1.177 | 0.4   | 98    | 0.00     |

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| 88   | Benzo(k)fluoranthene   | 1.161 | 1.153 | 0.7  | 98    | 0.00     |
| 89 S | Benzo(a)pyrene-d12     | 0.925 | 0.904 | 2.3  | 97    | -0.01    |
| 90 C | Benzo(a)pyrene         | 1.140 | 1.116 | 2.1  | 98    | 0.00     |
| 91   | Indeno(1,2,3-cd)pyrene | 1.291 | 1.236 | 4.3  | 95    | 0.00     |
| 92   | Dibenzo(a,h)anthracene | 1.056 | 0.928 | 12.1 | 86    | -0.02    |
| 93   | Benzo(g,h,i)perylene   | 1.027 | 1.020 | 0.7  | 103   | -0.02    |

(#) = Out of Range

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