

Data Path : Z:\HPCHEM1\BNA G\DATA\BG113016\  
 Data File : BG024955.D  
 Acq On : 30 Nov 2016 20:57  
 Operator : UM/SJ  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02012

Quant Time: Dec 01 01:13:57 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Dec 01 01:03:36 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  | 99    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.504 | 0.478 | 5.2  | 97    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.387 | 0.346 | 10.6 | 90    | 0.00     |
| 4    | Benzaldehyde                | 1.299 | 1.315 | -1.2 | 96    | 0.00     |
| 5 S  | Phenol-d5                   | 1.725 | 1.704 | 1.2  | 98    | 0.00     |
| 6    | Phenol                      | 1.771 | 1.781 | -0.6 | 97    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.067 | 1.092 | -2.3 | 100   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.296 | 1.318 | -1.7 | 103   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.207 | 1.273 | -5.5 | 103   | 0.00     |
| 10   | 2-Chlorophenol              | 1.213 | 1.185 | 2.3  | 96    | 0.00     |
| 11   | 2-Methylphenol              | 1.304 | 1.296 | 0.6  | 99    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.204 | 2.361 | -7.1 | 104   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.442 | 1.426 | 1.1  | 100   | 0.00     |
| 14   | Acetophenone                | 2.194 | 2.202 | -0.4 | 101   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.267 | 1.334 | -5.3 | 102   | 0.00     |
| 16   | 4-Methylphenol              | 1.451 | 1.484 | -2.3 | 108   | 0.00     |
| 17   | Hexachloroethane            | 0.537 | 0.569 | -6.0 | 102   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 103   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.144 | -1.4 | 105   | 0.00     |
| 20   | Nitrobenzene                | 0.435 | 0.427 | 1.8  | 104   | 0.00     |
| 21   | Isophorone                  | 0.824 | 0.805 | 2.3  | 100   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.177 | 0.175 | 1.1  | 105   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.179 | 0.172 | 3.9  | 100   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.410 | 0.407 | 0.7  | 100   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.420 | 0.425 | -1.2 | 102   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.340 | 0.335 | 1.5  | 101   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.330 | 0.332 | -0.6 | 103   | 0.00     |
| 28   | Naphthalene                 | 0.930 | 0.928 | 0.2  | 103   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.334 | 0.320 | 4.2  | 99    | 0.00     |
| 30   | 4-Chloroaniline             | 0.341 | 0.348 | -2.1 | 103   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.283 | 0.271 | 4.2  | 98    | 0.00     |
| 32   | Caprolactam                 | 0.137 | 0.141 | -2.9 | 109   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.406 | 0.410 | -1.0 | 103   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.734 | 0.722 | 1.6  | 98    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 103   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.603 | 0.590 | 2.2  | 104   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.354 | 0.324 | 8.5  | 103   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.379 | 0.384 | -1.3 | 106   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.415 | 0.426 | -2.7 | 104   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.213 | 1.247 | -2.8 | 104   | 0.00     |
| 41   | 2-Chloronaphthalene         | 0.965 | 0.962 | 0.3  | 106   | 0.00     |
| 42   | 2-Nitroaniline              | 0.378 | 0.385 | -1.9 | 101   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.376 | 1.385 | -0.7 | 101   | 0.00     |
| 44   | Dimethylphthalate           | 1.352 | 1.370 | -1.3 | 103   | 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 | 0.291 | 0.0   | 100   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.507 | 1.508 | -0.1  | 102   | 0.00     |
| 47   | Acenaphthylene              | 1.465 | 1.472 | -0.5  | 101   | 0.00     |
| 48   | 3-Nitroaniline              | 0.252 | 0.250 | 0.8   | 99    | 0.00     |
| 49 C | Acenaphthene                | 1.004 | 1.053 | -4.9  | 108   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.189 | 0.180 | 4.8   | 110   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.236 | 0.225 | 4.7   | 96    | 0.00     |
| 52   | 4-Nitrophenol               | 0.284 | 0.292 | -2.8  | 108   | 0.00     |
| 53   | Dibenzofuran                | 1.587 | 1.597 | -0.6  | 103   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.438 | 0.455 | -3.9  | 106   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.439 | 0.452 | -3.0  | 105   | 0.00     |
| 56   | Diethylphthalate            | 1.432 | 1.453 | -1.5  | 102   | 0.00     |
| 57 S | Fluorene-d10                | 1.295 | 1.288 | 0.5   | 101   | 0.00     |
| 58   | Fluorene                    | 1.292 | 1.306 | -1.1  | 102   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.728 | 0.738 | -1.4  | 104   | 0.00     |
| 60   | 4-Nitroaniline              | 0.303 | 0.298 | 1.7   | 102   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.124 | 0.120 | 3.2   | 106   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.128 | 0.133 | -3.9  | 111   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.515 | 0.512 | 0.6   | 101   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.226 | 0.232 | -2.7  | 106   | 0.00     |
| 66   | Hexachlorobenzene           | 0.254 | 0.262 | -3.1  | 104   | 0.00     |
| 67   | Atrazine                    | 0.245 | 0.252 | -2.9  | 103   | 0.00     |
| 68 C | Pentachlorophenol           | 0.153 | 0.145 | 5.2   | 102   | 0.00     |
| 69   | Phenanthrene                | 0.955 | 0.994 | -4.1  | 102   | 0.00     |
| 70 S | Anthracene-d10              | 0.844 | 0.878 | -4.0  | 105   | 0.00     |
| 71   | Anthracene                  | 0.981 | 1.008 | -2.8  | 100   | 0.00     |
| 72   | Carbazole                   | 0.819 | 0.886 | -8.2  | 103   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.034 | 1.071 | -3.6  | 102   | 0.00     |
| 74 C | Fluoranthene                | 1.134 | 1.303 | -14.9 | 105   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 105   | 0.00     |
| 76 S | Pyrene-d10                  | 0.824 | 0.836 | -1.5  | 104   | 0.00     |
| 77   | Pyrene                      | 0.961 | 0.981 | -2.1  | 102   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.376 | 0.383 | -1.9  | 103   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.349 | 0.374 | -7.2  | 108   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.039 | 1.062 | -2.2  | 104   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.518 | 0.509 | 1.7   | 101   | 0.00     |
| 82   | Chrysene                    | 0.972 | 0.992 | -2.1  | 104   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 103   | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.812 | 0.891 | -9.7  | 105   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 0.992 | 1.021 | -2.9  | 103   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 0.943 | 0.975 | -3.4  | 104   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.810 | 0.844 | -4.2  | 105   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 0.954 | 0.991 | -3.9 | 104   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.185 | 1.221 | -3.0 | 105   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 0.998 | 1.020 | -2.2 | 105   | -0.01    |
| 91   | Benzo(a,h,i)perylene   | 0.989 | 0.991 | -0.2 | 102   | 0.00     |

(#) = Out of Range

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