

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG061617\  
 Data File : BG027490.D  
 Acq On : 17 Jun 2017 3:19  
 Operator : SJ/MA  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02086

Quant Time: Jun 17 03:51:48 2017  
 Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG061617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jun 16 22:56:07 2017  
 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    | 99    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.579 | 0.543 | 6.2    | 96    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.500 | 0.498 | 0.4    | 102   | 0.00     |
| 4    | Benzaldehyde                | 1.268 | 1.541 | -21.5# | 103   | 0.00     |
| 5 S  | Phenol-d5                   | 2.169 | 2.160 | 0.4    | 103   | 0.00     |
| 6    | Phenol                      | 2.273 | 2.330 | -2.5   | 106   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.478 | 1.558 | -5.4   | 104   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.671 | 1.722 | -3.1   | 102   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.433 | 1.471 | -2.7   | 101   | 0.00     |
| 10   | 2-Chlorophenol              | 1.443 | 1.504 | -4.2   | 101   | 0.00     |
| 11   | 2-Methylphenol              | 1.643 | 1.626 | 1.0    | 102   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.782 | 3.100 | -11.4  | 109   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.696 | 1.765 | -4.1   | 108   | 0.00     |
| 14   | Acetophenone                | 2.621 | 2.667 | -1.8   | 102   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.481 | 1.661 | -12.2  | 112   | 0.00     |
| 16   | 4-Methylphenol              | 1.804 | 1.853 | -2.7   | 105   | 0.00     |
| 17   | Hexachloroethane            | 0.553 | 0.571 | -3.3   | 100   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 102   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.145 | 0.151 | -4.1   | 104   | 0.00     |
| 20   | Nitrobenzene                | 0.409 | 0.439 | -7.3   | 108   | 0.00     |
| 21   | Isophorone                  | 0.807 | 0.844 | -4.6   | 105   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.154 | 0.169 | -9.7   | 112   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.169 | 0.182 | -7.7   | 108   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.404 | 0.439 | -8.7   | 110   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.496 | 0.514 | -3.6   | 104   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.292 | 0.305 | -4.5   | 107   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.295 | 0.309 | -4.7   | 105   | 0.00     |
| 28   | Naphthalene                 | 1.028 | 1.050 | -2.1   | 103   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.369 | 0.446 | -20.9# | 103   | 0.00     |
| 30   | 4-Chloroaniline             | 0.364 | 0.430 | -18.1  | 102   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.181 | 0.180 | 0.6    | 103   | 0.00     |
| 32   | Caprolactam                 | 0.130 | 0.128 | 1.5    | 100   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.388 | 0.419 | -8.0   | 107   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.762 | 0.794 | -4.2   | 104   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 104   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.592 | 0.579 | 2.2    | 102   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.332 | 0.309 | 6.9    | 104   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.392 | 0.405 | -3.3   | 106   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.408 | 0.423 | -3.7   | 106   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.582 | 1.575 | 0.4    | 103   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.193 | 1.176 | 1.4    | 103   | 0.00     |
| 42   | 2-Nitroaniline              | 0.449 | 0.503 | -12.0  | 116   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.675 | 1.643 | 1.9    | 102   | 0.00     |
| 44   | Dimethylphthalate           | 1.661 | 1.621 | 2.4    | 100   | 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          | 0.309 | 0.315 | -1.9  | 105   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.929 | 1.965 | -1.9  | 104   | 0.00     |
| 47   | Acenaphthylene              | 1.933 | 1.950 | -0.9  | 104   | 0.00     |
| 48   | 3-Nitroaniline              | 0.336 | 0.338 | -0.6  | 102   | 0.00     |
| 49 C | Acenaphthene                | 1.348 | 1.328 | 1.5   | 104   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.214 | 0.191 | 10.7  | 106   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.328 | 0.306 | 6.7   | 98    | 0.00     |
| 52   | 4-Nitrophenol               | 0.261 | 0.260 | 0.4   | 103   | 0.00     |
| 53   | Dibenzofuran                | 1.965 | 1.934 | 1.6   | 103   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.473 | 0.475 | -0.4  | 101   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.397 | 0.403 | -1.5  | 107   | 0.00     |
| 56   | Diethylphthalate            | 1.755 | 1.714 | 2.3   | 101   | 0.00     |
| 57 S | Fluorene-d10                | 1.571 | 1.540 | 2.0   | 103   | 0.00     |
| 58   | Fluorene                    | 1.630 | 1.630 | 0.0   | 104   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.820 | 0.797 | 2.8   | 102   | 0.00     |
| 60   | 4-Nitroaniline              | 0.411 | 0.389 | 5.4   | 99    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 98    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.124 | 0.117 | 5.6   | 104   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.131 | 0.127 | 3.1   | 104   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.575 | 0.597 | -3.8  | 101   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.203 | 0.206 | -1.5  | 102   | 0.00     |
| 66   | Hexachlorobenzene           | 0.213 | 0.219 | -2.8  | 100   | 0.00     |
| 67   | Atrazine                    | 0.220 | 0.223 | -1.4  | 96    | 0.00     |
| 68 C | Pentachlorophenol           | 0.142 | 0.132 | 7.0   | 98    | 0.00     |
| 69   | Phenanthrene                | 1.075 | 1.079 | -0.4  | 98    | 0.00     |
| 70 S | Anthracene-d10              | 0.933 | 0.954 | -2.3  | 99    | 0.00     |
| 71   | Anthracene                  | 1.093 | 1.116 | -2.1  | 98    | 0.00     |
| 72   | Carbazole                   | 0.940 | 0.944 | -0.4  | 95    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.162 | 1.172 | -0.9  | 94    | 0.00     |
| 74 C | Fluoranthene                | 1.209 | 1.221 | -1.0  | 93    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 96    | 0.00     |
| 76 S | Pyrene-d10                  | 0.977 | 0.985 | -0.8  | 92    | 0.00     |
| 77   | Pyrene                      | 1.182 | 1.196 | -1.2  | 92    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.446 | 0.492 | -10.3 | 104   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.368 | 0.381 | -3.5  | 97    | 0.00     |
| 80   | Benzo(a)anthracene          | 1.114 | 1.139 | -2.2  | 94    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.640 | 0.719 | -12.3 | 103   | 0.00     |
| 82   | Chrysene                    | 1.006 | 1.040 | -3.4  | 97    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 84   | Di-n-octyl phthalate        | 1.253 | 1.303 | -4.0  | 104   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.191 | 1.167 | 2.0   | 96    | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.145 | 1.173 | -2.4  | 99    | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.901 | 0.907 | -0.7  | 98    | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.144 | 1.138 | 0.5  | 97    | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.306 | 1.321 | -1.1 | 99    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.121 | 1.126 | -0.4 | 98    | 0.00     |
| 91   | Benzo(g,h,i)perylene   | 1.069 | 1.070 | -0.1 | 97    | 0.00     |

(#) = Out of Range

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