

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122815\  
 Data File : BG020567.D  
 Acq On : 28 Dec 2015 16:32  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02001

Quant Time: Dec 29 02:49:11 2015  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG122415.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Dec 28 04:59:10 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    | 71    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.416 | 0.495 | -19.0  | 76    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.363 | 0.400 | -10.2  | 71    | 0.00     |
| 4    | Benzaldehyde                | 0.973 | 1.107 | -13.8  | 68    | 0.00     |
| 5 S  | Phenol-d5                   | 1.605 | 1.758 | -9.5   | 72    | 0.00     |
| 6    | Phenol                      | 1.671 | 1.926 | -15.3  | 77    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.951 | 1.068 | -12.3  | 73    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.218 | 1.370 | -12.5  | 76    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.231 | 1.396 | -13.4  | 72    | 0.00     |
| 10   | 2-Chlorophenol              | 1.285 | 1.409 | -9.6   | 69    | 0.00     |
| 11   | 2-Methylphenol              | 1.297 | 1.461 | -12.6  | 74    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.579 | 1.820 | -15.3  | 73    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.350 | 1.509 | -11.8  | 72    | 0.00     |
| 14   | Acetophenone                | 2.164 | 2.506 | -15.8  | 75    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.094 | 1.295 | -18.4  | 75    | 0.00     |
| 16   | 4-Methylphenol              | 1.407 | 1.680 | -19.4  | 78    | 0.00     |
| 17   | Hexachloroethane            | 0.543 | 0.591 | -8.8   | 70    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 74    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.157 | 0.173 | -10.2  | 74    | 0.00     |
| 20   | Nitrobenzene                | 0.391 | 0.428 | -9.5   | 73    | 0.00     |
| 21   | Isophorone                  | 0.738 | 0.860 | -16.5  | 78    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.178 | 0.199 | -11.8  | 75    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.188 | 0.209 | -11.2  | 77    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.400 | 0.440 | -10.0  | 73    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.410 | 0.453 | -10.5  | 75    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.345 | 0.391 | -13.3  | 75    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.354 | 0.388 | -9.6   | 73    | 0.00     |
| 28   | Naphthalene                 | 1.063 | 1.145 | -7.7   | 73    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.394 | 0.478 | -21.3# | 79    | 0.00     |
| 30   | 4-Chloroaniline             | 0.397 | 0.487 | -22.7# | 79    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.296 | 0.295 | 0.3    | 69    | 0.00     |
| 32   | Caprolactam                 | 0.114 | 0.158 | -38.6# | 96    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.371 | 0.467 | -25.9# | 82    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.787 | 0.881 | -11.9  | 76    | 0.00     |
| 35   | 1-Methylnaphthalene         | 0.747 | 0.855 | -14.5  | 76    | 0.00     |
| 36 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 84    | 0.00     |
| 37   | 1,2,4,5-Tetrachlorobenzene  | 0.810 | 0.763 | 5.8    | 74    | 0.00     |
| 38   | Hexachlorocyclopentadiene   | 0.441 | 0.245 | 44.4#  | 48    | 0.00     |
| 39 C | 2,4,6-Trichlorophenol       | 0.481 | 0.487 | -1.2   | 77    | 0.00     |
| 40   | 2,4,5-Trichlorophenol       | 0.526 | 0.569 | -8.2   | 84    | 0.00     |
| 41   | 1,1'-Biphenyl               | 1.544 | 1.553 | -0.6   | 78    | 0.00     |
| 42   | 2-Chloronaphthalene         | 1.245 | 1.260 | -1.2   | 80    | 0.00     |
| 43   | 2-Nitroaniline              | 0.349 | 0.405 | -16.0  | 89    | 0.00     |
| 44 S | Dimethylphthalate-d6        | 1.648 | 1.839 | -11.6  | 86    | 0.00     |

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
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 LabSampleId :  
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 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG122415.M  
 Quant Title : SVOA CALIBRATION  
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 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) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 45   | Dimethylphthalate           | 1.676 | 1.860 | -11.0  | 85    | 0.00     |
| 46   | 2,6-Dinitrotoluene          | 0.341 | 0.395 | -15.8  | 88    | 0.00     |
| 47 S | Acenaphthylene-d8           | 1.910 | 1.989 | -4.1   | 81    | 0.00     |
| 48   | Acenaphthylene              | 2.008 | 2.123 | -5.7   | 81    | 0.00     |
| 49   | 3-Nitroaniline              | 0.315 | 0.389 | -23.5# | 93    | 0.00     |
| 50 C | Acenaphthene                | 1.357 | 1.442 | -6.3   | 83    | 0.00     |
| 51   | 2,4-Dinitrophenol           | 0.218 | 0.143 | 34.4#  | 57    | 0.00     |
| 52 S | 4-Nitrophenol-d4            | 0.275 | 0.313 | -13.8  | 88    | 0.00     |
| 53   | 4-Nitrophenol               | 0.249 | 0.271 | -8.8   | 86    | 0.00     |
| 54   | Dibenzofuran                | 2.019 | 2.236 | -10.7  | 84    | 0.00     |
| 55   | 2,4-Dinitrotoluene          | 0.510 | 0.609 | -19.4  | 90    | 0.00     |
| 56   | 2,3,4,6-Tetrachlorophenol   | 0.518 | 0.564 | -8.9   | 84    | 0.00     |
| 57   | Diethylphthalate            | 1.720 | 1.982 | -15.2  | 88    | 0.00     |
| 58 S | Fluorene-d10                | 1.491 | 1.641 | -10.1  | 85    | 0.00     |
| 59   | Fluorene                    | 1.636 | 1.786 | -9.2   | 83    | 0.00     |
| 60   | 4-Chlorophenyl-phenylether  | 0.947 | 0.999 | -5.5   | 82    | 0.00     |
| 61   | 4-Nitroaniline              | 0.353 | 0.426 | -20.7# | 93    | 0.00     |
| 62 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 91    | 0.00     |
| 63 S | 4,6-Dinitro-2-methylphenol- | 0.133 | 0.124 | 6.8    | 83    | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol  | 0.141 | 0.131 | 7.1    | 81    | 0.00     |
| 65   | N-Nitrosodiphenylamine      | 0.559 | 0.585 | -4.7   | 87    | 0.00     |
| 66   | 4-Bromophenyl-phenylether   | 0.249 | 0.251 | -0.8   | 85    | 0.00     |
| 67   | Hexachlorobenzene           | 0.278 | 0.278 | 0.0    | 84    | 0.00     |
| 68   | Atrazine                    | 0.248 | 0.271 | -9.3   | 92    | 0.00     |
| 69 C | Pentachlorophenol           | 0.150 | 0.118 | 21.3   | 71    | 0.00     |
| 70   | Phenanthrene                | 1.113 | 1.173 | -5.4   | 88    | 0.00     |
| 71 S | Anthracene-d10              | 0.937 | 1.000 | -6.7   | 90    | 0.00     |
| 72   | Anthracene                  | 1.138 | 1.207 | -6.1   | 89    | 0.00     |
| 73   | 1,2,3,4-Tetrachlorobenzene  | 0.334 | 0.295 | 11.7   | 75    | 0.00     |
| 74   | Pentachlorobenzene          | 0.310 | 0.294 | 5.2    | 81    | 0.00     |
| 75   | Carbazole                   | 0.959 | 1.069 | -11.5  | 93    | 0.00     |
| 76   | Di-n-butylphthalate         | 1.143 | 1.252 | -9.5   | 92    | 0.00     |
| 77 C | Fluoranthene                | 1.399 | 1.544 | -10.4  | 89    | 0.00     |
| 78 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 89    | 0.00     |
| 79 S | Pyrene-d10                  | 0.866 | 0.954 | -10.2  | 90    | 0.00     |
| 80   | Pyrene                      | 1.131 | 1.235 | -9.2   | 89    | 0.00     |
| 81   | Butylbenzylphthalate        | 0.402 | 0.444 | -10.4  | 90    | 0.00     |
| 82   | 3,3'-Dichlorobenzidine      | 0.392 | 0.418 | -6.6   | 85    | 0.00     |
| 83   | Benzo(a)anthracene          | 1.175 | 1.265 | -7.7   | 88    | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate  | 0.588 | 0.638 | -8.5   | 89    | 0.00     |
| 85   | Chrysene                    | 1.119 | 1.191 | -6.4   | 87    | 0.00     |
| 86 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 91    | 0.00     |
| 87   | Di-n-octyl phthalate        | 1.035 | 1.132 | -9.4   | 90    | 0.00     |

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Misc :  
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BNA\_G  
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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.200 | 1.257 | -4.7 | 86    | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.160 | 1.223 | -5.4 | 89    | 0.00     |
| 90 S | Benzo(a)pyrene-d12     | 1.011 | 1.065 | -5.3 | 87    | 0.00     |
| 91 C | Benzo(a)pyrene         | 1.154 | 1.219 | -5.6 | 88    | 0.00     |
| 92   | Indeno(1,2,3-cd)pyrene | 1.382 | 1.466 | -6.1 | 90    | 0.00     |
| 93   | Dibenzo(a,h)anthracene | 1.147 | 1.202 | -4.8 | 88    | 0.00     |
| 94   | Benzo(a,h,i)perylene   | 1.135 | 1.210 | -6.6 | 90    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 1