

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120616\  
 Data File : BG024972.D  
 Acq On : 6 Dec 2016 9:29  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02015

Quant Time: Dec 07 02:35:43 2016  
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG113016.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 02 23:46:34 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   | 107   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.504 | 0.489 | 3.0   | 107   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.387 | 0.391 | -1.0  | 109   | 0.00     |
| 4    | Benzaldehyde                | 1.299 | 1.358 | -4.5  | 107   | 0.00     |
| 5 S  | Phenol-d5                   | 1.725 | 1.681 | 2.6   | 104   | 0.00     |
| 6    | Phenol                      | 1.771 | 1.802 | -1.8  | 105   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.067 | 1.097 | -2.8  | 109   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.296 | 1.250 | 3.5   | 105   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.207 | 1.207 | 0.0   | 105   | 0.00     |
| 10   | 2-Chlorophenol              | 1.213 | 1.155 | 4.8   | 101   | 0.00     |
| 11   | 2-Methylphenol              | 1.304 | 1.284 | 1.5   | 105   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.204 | 2.219 | -0.7  | 105   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.442 | 1.374 | 4.7   | 104   | 0.00     |
| 14   | Acetophenone                | 2.194 | 2.180 | 0.6   | 107   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.267 | 1.271 | -0.3  | 105   | 0.00     |
| 16   | 4-Methylphenol              | 1.451 | 1.422 | 2.0   | 111   | 0.00     |
| 17   | Hexachloroethane            | 0.537 | 0.559 | -4.1  | 109   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.149 | -4.9  | 111   | 0.00     |
| 20   | Nitrobenzene                | 0.435 | 0.435 | 0.0   | 109   | 0.00     |
| 21   | Isophorone                  | 0.824 | 0.811 | 1.6   | 103   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.177 | 0.164 | 7.3   | 100   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.179 | 0.178 | 0.6   | 106   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.410 | 0.406 | 1.0   | 102   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.420 | 0.419 | 0.2   | 103   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.340 | 0.345 | -1.5  | 107   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.330 | 0.333 | -0.9  | 105   | 0.00     |
| 28   | Naphthalene                 | 0.930 | 0.908 | 2.4   | 103   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.334 | 0.391 | -17.1 | 123   | 0.00     |
| 30   | 4-Chloroaniline             | 0.341 | 0.375 | -10.0 | 114   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.283 | 0.280 | 1.1   | 103   | 0.00     |
| 32   | Caprolactam                 | 0.137 | 0.134 | 2.2   | 106   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.406 | 0.399 | 1.7   | 103   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.734 | 0.737 | -0.4  | 103   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 103   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.603 | 0.625 | -3.6  | 110   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.354 | 0.338 | 4.5   | 107   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.379 | 0.374 | 1.3   | 104   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.415 | 0.412 | 0.7   | 101   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.213 | 1.252 | -3.2  | 105   | 0.00     |
| 41   | 2-Chloronaphthalene         | 0.965 | 0.981 | -1.7  | 108   | 0.00     |
| 42   | 2-Nitroaniline              | 0.378 | 0.395 | -4.5  | 104   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.376 | 1.403 | -2.0  | 102   | 0.00     |
| 44   | Dimethylphthalate           | 1.352 | 1.384 | -2.4  | 105   | 0.00     |

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 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
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 LabSampleId :  
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 Quant Title : SVOA CALIBRATION  
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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                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.291 | 0.283 | 2.7   | 98    | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.507 | 1.536 | -1.9  | 104   | 0.00     |
| 47   | Acenaphthylene              | 1.465 | 1.485 | -1.4  | 102   | 0.00     |
| 48   | 3-Nitroaniline              | 0.252 | 0.283 | -12.3 | 112   | 0.00     |
| 49 C | Acenaphthene                | 1.004 | 1.019 | -1.5  | 105   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.189 | 0.181 | 4.2   | 111   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.236 | 0.249 | -5.5  | 107   | 0.00     |
| 52   | 4-Nitrophenol               | 0.284 | 0.316 | -11.3 | 117   | 0.00     |
| 53   | Dibenzofuran                | 1.587 | 1.602 | -0.9  | 103   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.438 | 0.448 | -2.3  | 104   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.439 | 0.436 | 0.7   | 102   | 0.00     |
| 56   | Diethylphthalate            | 1.432 | 1.433 | -0.1  | 101   | 0.00     |
| 57 S | Fluorene-d10                | 1.295 | 1.314 | -1.5  | 103   | 0.00     |
| 58   | Fluorene                    | 1.292 | 1.336 | -3.4  | 104   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.728 | 0.736 | -1.1  | 104   | 0.00     |
| 60   | 4-Nitroaniline              | 0.303 | 0.308 | -1.7  | 105   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 105   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.124 | 0.122 | 1.6   | 109   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.128 | 0.129 | -0.8  | 110   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.515 | 0.506 | 1.7   | 102   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.226 | 0.224 | 0.9   | 105   | 0.00     |
| 66   | Hexachlorobenzene           | 0.254 | 0.268 | -5.5  | 108   | 0.00     |
| 67   | Atrazine                    | 0.245 | 0.248 | -1.2  | 104   | 0.00     |
| 68 C | Pentachlorophenol           | 0.153 | 0.151 | 1.3   | 108   | 0.00     |
| 69   | Phenanthrene                | 0.955 | 0.993 | -4.0  | 104   | 0.00     |
| 70 S | Anthracene-d10              | 0.844 | 0.879 | -4.1  | 107   | 0.00     |
| 71   | Anthracene                  | 0.981 | 1.011 | -3.1  | 103   | 0.00     |
| 72   | Carbazole                   | 0.819 | 0.908 | -10.9 | 108   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.034 | 1.091 | -5.5  | 106   | 0.00     |
| 74 C | Fluoranthene                | 1.134 | 1.335 | -17.7 | 109   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 113   | 0.00     |
| 76 S | Pyrene-d10                  | 0.824 | 0.817 | 0.8   | 109   | 0.00     |
| 77   | Pyrene                      | 0.961 | 0.976 | -1.6  | 110   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.376 | 0.380 | -1.1  | 110   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.349 | 0.375 | -7.4  | 116   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.039 | 1.053 | -1.3  | 111   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.518 | 0.513 | 1.0   | 109   | 0.00     |
| 82   | Chrysene                    | 0.972 | 0.987 | -1.5  | 112   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 114   | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.812 | 0.851 | -4.8  | 111   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 0.992 | 1.018 | -2.6  | 113   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 0.943 | 0.950 | -0.7  | 112   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.810 | 0.833 | -2.8  | 114   | 0.00     |

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Misc :  
ALS Vial : 2 Sample Multiplier: 1

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 0.954 | 0.971 | -1.8 | 113   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.185 | 1.199 | -1.2 | 114   | 0.01     |
| 90   | Dibenzo(a,h)anthracene | 0.998 | 0.998 | 0.0  | 113   | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 0.989 | 0.995 | -0.6 | 113   | 0.00     |

(#) = Out of Range

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