

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG082218\  
 Data File : BG036348.D  
 Acq On : 22 Aug 2018 18:35  
 Operator : JU/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 LabSampleId :  
 SSTD02021

Quant Time: Aug 22 19:14:11 2018  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG082118MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Aug 22 18:34:10 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   | 109   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.566 | 0.539 | 4.8   | 111   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.537 | 0.493 | 8.2   | 98    | 0.00     |
| 4    | Benzaldehyde                | 1.247 | 1.317 | -5.6  | 102   | 0.00     |
| 5 S  | Phenol-d5                   | 1.861 | 1.859 | 0.1   | 106   | 0.00     |
| 6    | Phenol                      | 1.860 | 1.896 | -1.9  | 109   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.275 | 1.273 | 0.2   | 106   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.381 | 1.417 | -2.6  | 108   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.344 | 1.347 | -0.2  | 107   | 0.00     |
| 10   | 2-Chlorophenol              | 1.344 | 1.355 | -0.8  | 111   | 0.00     |
| 11   | 2-Methylphenol              | 1.417 | 1.397 | 1.4   | 106   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 3.040 | 2.933 | 3.5   | 102   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.420 | 1.419 | 0.1   | 108   | 0.00     |
| 14   | Acetophenone                | 2.177 | 2.214 | -1.7  | 107   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.312 | 1.246 | 5.0   | 103   | 0.00     |
| 16   | 4-Methylphenol              | 1.482 | 1.490 | -0.5  | 107   | 0.00     |
| 17   | Hexachloroethane            | 0.535 | 0.521 | 2.6   | 105   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.122 | 0.141 | -15.6 | 132   | 0.00     |
| 20   | Nitrobenzene                | 0.346 | 0.381 | -10.1 | 123   | 0.00     |
| 21   | Isophorone                  | 0.840 | 0.803 | 4.4   | 102   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.123 | 0.136 | -10.6 | 130   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.130 | 0.148 | -13.8 | 128   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.397 | 0.389 | 2.0   | 106   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.454 | 0.453 | 0.2   | 106   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.329 | 0.328 | 0.3   | 107   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.311 | 0.316 | -1.6  | 107   | 0.00     |
| 28   | Naphthalene                 | 1.025 | 1.029 | -0.4  | 107   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.363 | 0.353 | 2.8   | 99    | 0.00     |
| 30   | 4-Chloroaniline             | 0.354 | 0.348 | 1.7   | 100   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.222 | 0.225 | -1.4  | 108   | 0.00     |
| 32   | Caprolactam                 | 0.129 | 0.122 | 5.4   | 98    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.366 | 0.356 | 2.7   | 102   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.765 | 0.773 | -1.0  | 108   | 0.00     |
| 35   | 1-Methylnaphthalene         | 0.746 | 0.745 | 0.1   | 105   | 0.00     |
| 36 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 106   | 0.00     |
| 37   | 1,2,4,5-Tetrachlorobenzene  | 0.681 | 0.697 | -2.3  | 108   | 0.00     |
| 38   | Hexachlorocyclopentadiene   | 0.408 | 0.397 | 2.7   | 108   | 0.00     |
| 39 C | 2,4,6-Trichlorophenol       | 0.421 | 0.413 | 1.9   | 105   | 0.00     |
| 40   | 2,4,5-Trichlorophenol       | 0.436 | 0.455 | -4.4  | 113   | 0.00     |
| 41   | 1,1'-Biphenyl               | 1.565 | 1.582 | -1.1  | 109   | 0.00     |
| 42   | 2-Chloronaphthalene         | 1.201 | 1.220 | -1.6  | 108   | 0.00     |
| 43   | 2-Nitroaniline              | 0.314 | 0.360 | -14.6 | 134   | 0.00     |
| 44 S | Dimethylphthalate-d6        | 1.672 | 1.674 | -0.1  | 106   | 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   | Dimethylphthalate           | 1.605 | 1.640 | -2.2   | 109   | 0.00     |
| 46   | 2,6-Dinitrotoluene          | 0.224 | 0.262 | -17.0  | 132   | 0.00     |
| 47 S | Acenaphthylene-d8           | 2.016 | 2.056 | -2.0   | 108   | 0.00     |
| 48   | Acenaphthylene              | 1.901 | 1.945 | -2.3   | 108   | 0.00     |
| 49   | 3-Nitroaniline              | 0.234 | 0.272 | -16.2  | 127   | 0.00     |
| 50 C | Acenaphthene                | 1.361 | 1.384 | -1.7   | 107   | 0.00     |
| 51   | 2,4-Dinitrophenol           | 0.114 | 0.115 | -0.9   | 132   | 0.00     |
| 52 S | 4-Nitrophenol-d4            | 0.247 | 0.277 | -12.1  | 126   | 0.00     |
| 53   | 4-Nitrophenol               | 0.230 | 0.249 | -8.3   | 122   | 0.00     |
| 54   | Dibenzofuran                | 1.884 | 1.931 | -2.5   | 109   | 0.00     |
| 55   | 2,4-Dinitrotoluene          | 0.310 | 0.383 | -23.5# | 141   | 0.00     |
| 56   | 2,3,4,6-Tetrachlorophenol   | 0.408 | 0.410 | -0.5   | 109   | 0.00     |
| 57   | Diethylphthalate            | 1.629 | 1.618 | 0.7    | 105   | 0.00     |
| 58 S | Fluorene-d10                | 1.454 | 1.452 | 0.1    | 106   | 0.00     |
| 59   | Fluorene                    | 1.519 | 1.534 | -1.0   | 106   | 0.00     |
| 60   | 4-Chlorophenyl-phenylether  | 0.827 | 0.838 | -1.3   | 107   | 0.00     |
| 61   | 4-Nitroaniline              | 0.296 | 0.341 | -15.2  | 127   | 0.00     |
| 62 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 107   | 0.00     |
| 63 S | 4,6-Dinitro-2-methylphenol- | 0.072 | 0.073 | -1.4   | 135   | 0.00     |
| 64   | 4,6-Dinitro-2-methylphenol  | 0.075 | 0.077 | -2.7   | 129   | 0.00     |
| 65   | N-Nitrosodiphenylamine      | 0.560 | 0.570 | -1.8   | 107   | 0.00     |
| 66   | 4-Bromophenyl-phenylether   | 0.221 | 0.222 | -0.5   | 107   | 0.00     |
| 67   | Hexachlorobenzene           | 0.221 | 0.233 | -5.4   | 112   | 0.00     |
| 68   | Atrazine                    | 0.231 | 0.241 | -4.3   | 106   | 0.00     |
| 69 C | Pentachlorophenol           | 0.134 | 0.135 | -0.7   | 109   | 0.00     |
| 70   | Phenanthrene                | 1.039 | 1.073 | -3.3   | 107   | 0.00     |
| 71 S | Anthracene-d10              | 0.933 | 0.947 | -1.5   | 106   | 0.00     |
| 72   | Anthracene                  | 1.046 | 1.069 | -2.2   | 108   | 0.00     |
| 73   | 1,2,3,4-Tetrachlorobenzene  | 0.306 | 0.309 | -1.0   | 108   | 0.00     |
| 74   | Pentachlorobenzene          | 0.266 | 0.272 | -2.3   | 108   | 0.00     |
| 75   | Carbazole                   | 0.879 | 0.963 | -9.6   | 107   | 0.00     |
| 76   | Di-n-butylphthalate         | 1.149 | 1.164 | -1.3   | 106   | 0.00     |
| 77 C | Fluoranthene                | 1.205 | 1.340 | -11.2  | 107   | 0.00     |
| 78 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 107   | 0.00     |
| 79 S | Pyrene-d10                  | 1.025 | 1.031 | -0.6   | 106   | 0.00     |
| 80   | Pyrene                      | 1.212 | 1.221 | -0.7   | 106   | 0.00     |
| 81   | Butylbenzylphthalate        | 0.493 | 0.487 | 1.2    | 106   | 0.00     |
| 82   | 3,3'-Dichlorobenzidine      | 0.412 | 0.397 | 3.6    | 96    | 0.00     |
| 83   | Benzo(a)anthracene          | 1.264 | 1.277 | -1.0   | 107   | 0.00     |
| 84   | Bis(2-ethylhexyl)phthalate  | 0.703 | 0.692 | 1.6    | 103   | 0.00     |
| 85   | Chrysene                    | 1.159 | 1.166 | -0.6   | 105   | 0.00     |
| 86 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 103   | 0.00     |
| 87   | Di-n-octyl phthalate        | 1.124 | 1.166 | -3.7   | 101   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88   | Benzo(b)fluoranthene   | 1.184 | 1.222 | -3.2 | 106   | 0.00     |
| 89   | Benzo(k)fluoranthene   | 1.142 | 1.176 | -3.0 | 106   | 0.00     |
| 90 S | Benzo(a)pyrene-d12     | 1.085 | 1.104 | -1.8 | 104   | 0.00     |
| 91 C | Benzo(a)pyrene         | 1.121 | 1.147 | -2.3 | 104   | 0.00     |
| 92   | Indeno(1,2,3-cd)pyrene | 1.334 | 1.370 | -2.7 | 106   | -0.01    |
| 93   | Dibenzo(a,h)anthracene | 1.082 | 1.117 | -3.2 | 107   | -0.01    |
| 94   | Benzo(g,h,i)perylene   | 1.100 | 1.132 | -2.9 | 106   | 0.00     |

(#) = Out of Range

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