

Data Path : U:\HPCHEM1\BNA M\Data\BM012418\  
 Data File : BM013783.D  
 Acq On : 25 Jan 2018 00:12  
 Operator : SJ/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02040

Quant Time: Jan 25 03:19:21 2018  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jan 25 00:32:09 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                    | AvqRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 112   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.554 | 0.444 | 19.9   | 94    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.506 | 0.392 | 22.5#  | 88    | 0.00     |
| 4    | Benzaldehyde                | 1.063 | 1.038 | 2.4    | 101   | 0.00     |
| 5 S  | Phenol-d5                   | 1.535 | 1.480 | 3.6    | 113   | 0.00     |
| 6    | Phenol                      | 1.540 | 1.500 | 2.6    | 111   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.930 | 0.869 | 6.6    | 105   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.205 | 1.124 | 6.7    | 109   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.251 | 1.289 | -3.0   | 114   | 0.00     |
| 10   | 2-Chlorophenol              | 1.247 | 1.241 | 0.5    | 114   | 0.00     |
| 11   | 2-Methylphenol              | 1.135 | 1.129 | 0.5    | 116   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.226 | 1.156 | 5.7    | 107   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.180 | 1.234 | -4.6   | 122   | 0.00     |
| 14   | Acetophenone                | 1.953 | 1.901 | 2.7    | 113   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 0.953 | 0.961 | -0.8   | 118   | 0.00     |
| 16   | 4-Methylphenol              | 1.214 | 1.225 | -0.9   | 116   | 0.00     |
| 17   | Hexachloroethane            | 0.607 | 0.571 | 5.9    | 109   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 117   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.155 | 0.150 | 3.2    | 115   | 0.00     |
| 20   | Nitrobenzene                | 0.401 | 0.369 | 8.0    | 109   | 0.00     |
| 21   | Isophorone                  | 0.648 | 0.625 | 3.5    | 114   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.164 | 0.169 | -3.0   | 123   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.171 | 0.172 | -0.6   | 120   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.360 | 0.358 | 0.6    | 117   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.395 | 0.373 | 5.6    | 109   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.317 | 0.332 | -4.7   | 122   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.305 | 0.304 | 0.3    | 117   | 0.00     |
| 28   | Naphthalene                 | 0.995 | 1.000 | -0.5   | 120   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.275 | 0.355 | -29.1# | 168#  | 0.00     |
| 30   | 4-Chloroaniline             | 0.273 | 0.360 | -31.9# | 177#  | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.251 | 0.238 | 5.2    | 113   | 0.00     |
| 32   | Caprolactam                 | 0.085 | 0.094 | -10.6  | 135   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.309 | 0.329 | -6.5   | 129   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.742 | 0.723 | 2.6    | 118   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 121   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.741 | 0.742 | -0.1   | 122   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.308 | 0.202 | 34.4#  | 91    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.429 | 0.459 | -7.0   | 127   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.446 | 0.477 | -7.0   | 128   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.665 | 1.643 | 1.3    | 119   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.301 | 1.289 | 0.9    | 119   | 0.00     |
| 42   | 2-Nitroaniline              | 0.359 | 0.380 | -5.8   | 122   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.648 | 1.718 | -4.2   | 127   | 0.00     |
| 44   | Dimethylphthalate           | 1.621 | 1.675 | -3.3   | 127   | 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          | 0.314 | 0.337 | -7.3  | 123   | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.102 | 2.112 | -0.5  | 122   | 0.00     |
| 47   | Acenaphthylene              | 1.923 | 1.939 | -0.8  | 120   | 0.00     |
| 48   | 3-Nitroaniline              | 0.247 | 0.296 | -19.8 | 156#  | 0.00     |
| 49 C | Acenaphthene                | 1.332 | 1.317 | 1.1   | 120   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.180 | 0.149 | 17.2  | 116   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.268 | 0.281 | -4.9  | 141   | 0.00     |
| 52   | 4-Nitrophenol               | 0.272 | 0.271 | 0.4   | 120   | 0.00     |
| 53   | Dibenzofuran                | 2.004 | 1.996 | 0.4   | 119   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.458 | 0.507 | -10.7 | 130   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.420 | 0.446 | -6.2  | 131   | 0.00     |
| 56   | Diethylphthalate            | 1.631 | 1.661 | -1.8  | 126   | 0.00     |
| 57 S | Fluorene-d10                | 1.450 | 1.472 | -1.5  | 122   | 0.00     |
| 58   | Fluorene                    | 1.641 | 1.686 | -2.7  | 125   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.887 | 0.872 | 1.7   | 119   | 0.00     |
| 60   | 4-Nitroaniline              | 0.289 | 0.313 | -8.3  | 126   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 124   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.120 | 0.108 | 10.0  | 124   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.127 | 0.119 | 6.3   | 125   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.588 | 0.594 | -1.0  | 128   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.236 | 0.232 | 1.7   | 124   | 0.00     |
| 66   | Hexachlorobenzene           | 0.254 | 0.253 | 0.4   | 126   | 0.00     |
| 67   | Atrazine                    | 0.235 | 0.235 | 0.0   | 130   | 0.00     |
| 68 C | Pentachlorophenol           | 0.135 | 0.131 | 3.0   | 137   | 0.00     |
| 69   | Phenanthrene                | 1.112 | 1.113 | -0.1  | 125   | 0.00     |
| 70 S | Anthracene-d10              | 0.974 | 0.971 | 0.3   | 123   | 0.00     |
| 71   | Anthracene                  | 1.147 | 1.138 | 0.8   | 126   | 0.00     |
| 72   | Carbazole                   | 0.947 | 0.975 | -3.0  | 131   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.113 | 1.144 | -2.8  | 129   | 0.00     |
| 74 C | Fluoranthene                | 1.327 | 1.353 | -2.0  | 128   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 119   | 0.00     |
| 76 S | Pyrene-d10                  | 0.931 | 1.007 | -8.2  | 129   | 0.00     |
| 77   | Pyrene                      | 1.216 | 1.292 | -6.3  | 128   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.428 | 0.478 | -11.7 | 134   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.331 | 0.395 | -19.3 | 149   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.209 | 1.261 | -4.3  | 126   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.643 | 0.714 | -11.0 | 132   | 0.00     |
| 82   | Chrysene                    | 1.109 | 1.136 | -2.4  | 123   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 121   | 0.00     |
| 84   | Di-n-octyl phthalate        | 1.205 | 1.196 | 0.7   | 129   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.228 | 1.320 | -7.5  | 127   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.187 | 1.173 | 1.2   | 117   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.919 | 0.947 | -3.0  | 124   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.160 | 1.172 | -1.0 | 121   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.370 | 1.364 | 0.4  | 121   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.157 | 1.154 | 0.3  | 121   | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 1.127 | 1.125 | 0.2  | 121   | 0.00     |

(#) = Out of Range

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