

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121816\  
 Data File : BM008403.D  
 Acq On : 17 Dec 2016 09:51  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02062

Quant Time: Dec 18 02:32:29 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121416.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Sun Dec 18 00:00:12 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    | 117   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.499 | 0.500 | -0.2   | 110   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.426 | 0.466 | -9.4   | 119   | 0.00     |
| 4    | Benzaldehyde                | 1.039 | 1.204 | -15.9  | 115   | 0.00     |
| 5 S  | Phenol-d5                   | 1.565 | 1.614 | -3.1   | 116   | 0.00     |
| 6    | Phenol                      | 1.623 | 1.690 | -4.1   | 117   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.901 | 0.954 | -5.9   | 116   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.212 | 1.254 | -3.5   | 115   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.174 | 1.268 | -8.0   | 119   | 0.00     |
| 10   | 2-Chlorophenol              | 1.188 | 1.296 | -9.1   | 118   | 0.00     |
| 11   | 2-Methylphenol              | 1.186 | 1.223 | -3.1   | 116   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.387 | 1.484 | -7.0   | 116   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.207 | 1.236 | -2.4   | 115   | 0.00     |
| 14   | Acetophenone                | 1.978 | 2.041 | -3.2   | 116   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.013 | 1.081 | -6.7   | 116   | 0.00     |
| 16   | 4-Methylphenol              | 1.281 | 1.320 | -3.0   | 117   | 0.00     |
| 17   | Hexachloroethane            | 0.542 | 0.586 | -8.1   | 117   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 117   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.144 | 0.154 | -6.9   | 118   | 0.00     |
| 20   | Nitrobenzene                | 0.370 | 0.406 | -9.7   | 120   | 0.00     |
| 21   | Isophorone                  | 0.671 | 0.721 | -7.5   | 116   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.159 | 0.167 | -5.0   | 116   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.163 | 0.175 | -7.4   | 118   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.367 | 0.391 | -6.5   | 116   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.383 | 0.414 | -8.1   | 114   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.293 | 0.317 | -8.2   | 115   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.295 | 0.315 | -6.8   | 114   | 0.00     |
| 28   | Naphthalene                 | 0.947 | 1.017 | -7.4   | 116   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.335 | 0.366 | -9.3   | 113   | 0.00     |
| 30   | 4-Chloroaniline             | 0.340 | 0.384 | -12.9  | 117   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.240 | 0.261 | -8.8   | 120   | 0.00     |
| 32   | Caprolactam                 | 0.092 | 0.089 | 3.3    | 109   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.319 | 0.341 | -6.9   | 114   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.684 | 0.731 | -6.9   | 116   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 116   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.627 | 0.673 | -7.3   | 115   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.250 | 0.349 | -39.6# | 201#  | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.356 | 0.383 | -7.6   | 116   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.385 | 0.405 | -5.2   | 116   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.336 | 1.437 | -7.6   | 115   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.025 | 1.095 | -6.8   | 114   | 0.00     |
| 42   | 2-Nitroaniline              | 0.302 | 0.329 | -8.9   | 116   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.332 | 1.403 | -5.3   | 112   | 0.00     |
| 44   | Dimethylphthalate           | 1.305 | 1.379 | -5.7   | 112   | 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          | 0.252 | 0.275 | -9.1  | 116   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.571 | 1.698 | -8.1  | 115   | 0.00     |
| 47   | Acenaphthylene              | 1.606 | 1.719 | -7.0  | 113   | 0.00     |
| 48   | 3-Nitroaniline              | 0.253 | 0.269 | -6.3  | 116   | 0.00     |
| 49 C | Acenaphthene                | 1.106 | 1.181 | -6.8  | 113   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.148 | 0.133 | 10.1  | 119   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.196 | 0.157 | 19.9  | 97    | 0.00     |
| 52   | 4-Nitrophenol               | 0.189 | 0.212 | -12.2 | 134   | 0.00     |
| 53   | Dibenzofuran                | 1.607 | 1.696 | -5.5  | 112   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.379 | 0.412 | -8.7  | 114   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.358 | 0.379 | -5.9  | 112   | 0.00     |
| 56   | Diethylphthalate            | 1.298 | 1.369 | -5.5  | 113   | 0.00     |
| 57 S | Fluorene-d10                | 1.159 | 1.238 | -6.8  | 114   | 0.00     |
| 58   | Fluorene                    | 1.316 | 1.409 | -7.1  | 115   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.703 | 0.752 | -7.0  | 113   | 0.00     |
| 60   | 4-Nitroaniline              | 0.244 | 0.253 | -3.7  | 116   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 111   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.119 | 0.105 | 11.8  | 101   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.122 | 0.110 | 9.8   | 102   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.543 | 0.591 | -8.8  | 112   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.219 | 0.240 | -9.6  | 112   | 0.00     |
| 66   | Hexachlorobenzene           | 0.249 | 0.263 | -5.6  | 108   | 0.00     |
| 67   | Atrazine                    | 0.222 | 0.227 | -2.3  | 110   | 0.00     |
| 68 C | Pentachlorophenol           | 0.134 | 0.126 | 6.0   | 107   | 0.00     |
| 69   | Phenanthrene                | 1.027 | 1.106 | -7.7  | 111   | 0.00     |
| 70 S | Anthracene-d10              | 0.871 | 0.953 | -9.4  | 113   | 0.00     |
| 71   | Anthracene                  | 1.041 | 1.138 | -9.3  | 112   | 0.00     |
| 72   | Carbazole                   | 0.914 | 0.973 | -6.5  | 112   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.022 | 1.104 | -8.0  | 111   | 0.00     |
| 74 C | Fluoranthene                | 1.258 | 1.328 | -5.6  | 110   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 109   | 0.00     |
| 76 S | Pyrene-d10                  | 0.772 | 0.854 | -10.6 | 111   | 0.00     |
| 77   | Pyrene                      | 0.981 | 1.072 | -9.3  | 110   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.351 | 0.393 | -12.0 | 113   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.350 | 0.379 | -8.3  | 107   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.030 | 1.132 | -9.9  | 111   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.513 | 0.571 | -11.3 | 111   | 0.00     |
| 82   | Chrysene                    | 0.996 | 1.083 | -8.7  | 108   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 110   | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.957 | 1.001 | -4.6  | 111   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.076 | 1.183 | -9.9  | 113   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.031 | 1.109 | -7.6  | 106   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.834 | 0.920 | -10.3 | 112   | 0.00     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.042 | 1.135 | -8.9 | 109   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.262 | 1.366 | -8.2 | 109   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.059 | 1.146 | -8.2 | 109   | 0.00     |
| 91   | Benzo(a,h,i)perylene   | 1.052 | 1.147 | -9.0 | 110   | 0.01     |

(#) = Out of Range

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