

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011017\  
 Data File : BM008744.D  
 Acq On : 09 Jan 2017 18:02  
 Operator : UM/SJ  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02062

Quant Time: Jan 10 00:38:07 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010617.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Jan 06 16:08:18 2017  
 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   | 103   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.515 | 0.534 | -3.7  | 101   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.433 | 0.431 | 0.5   | 94    | 0.00     |
| 4    | Benzaldehyde                | 1.197 | 1.420 | -18.6 | 107   | 0.00     |
| 5 S  | Phenol-d5                   | 1.643 | 1.681 | -2.3  | 109   | 0.00     |
| 6    | Phenol                      | 1.766 | 1.760 | 0.3   | 103   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.115 | 1.151 | -3.2  | 107   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.295 | 1.294 | 0.1   | 104   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.144 | 1.184 | -3.5  | 104   | 0.00     |
| 10   | 2-Chlorophenol              | 1.172 | 1.199 | -2.3  | 108   | 0.00     |
| 11   | 2-Methylphenol              | 1.297 | 1.268 | 2.2   | 102   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.003 | 2.162 | -7.9  | 108   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.274 | 1.310 | -2.8  | 107   | 0.00     |
| 14   | Acetophenone                | 2.239 | 2.240 | -0.0  | 104   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.271 | 1.323 | -4.1  | 105   | 0.00     |
| 16   | 4-Methylphenol              | 1.401 | 1.382 | 1.4   | 101   | 0.00     |
| 17   | Hexachloroethane            | 0.621 | 0.628 | -1.1  | 104   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.123 | 0.127 | -3.3  | 105   | 0.00     |
| 20   | Nitrobenzene                | 0.438 | 0.468 | -6.8  | 110   | 0.00     |
| 21   | Isophorone                  | 0.775 | 0.763 | 1.5   | 103   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.134 | 0.142 | -6.0  | 119   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.144 | 0.145 | -0.7  | 107   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.406 | 0.411 | -1.2  | 106   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.410 | 0.401 | 2.2   | 101   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.309 | 0.303 | 1.9   | 103   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.306 | 0.307 | -0.3  | 105   | 0.00     |
| 28   | Naphthalene                 | 0.914 | 0.916 | -0.2  | 106   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.342 | 0.360 | -5.3  | 105   | 0.00     |
| 30   | 4-Chloroaniline             | 0.351 | 0.386 | -10.0 | 109   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.274 | 0.272 | 0.7   | 103   | 0.01     |
| 32   | Caprolactam                 | 0.094 | 0.087 | 7.4   | 102   | 0.01     |
| 33 C | 4-Chloro-3-methylphenol     | 0.367 | 0.359 | 2.2   | 102   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.709 | 0.700 | 1.3   | 104   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 107   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.599 | 0.600 | -0.2  | 105   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.426 | 0.391 | 8.2   | 101   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.362 | 0.348 | 3.9   | 103   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.390 | 0.383 | 1.8   | 104   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.254 | 1.232 | 1.8   | 105   | 0.00     |
| 41   | 2-Chloronaphthalene         | 0.984 | 0.971 | 1.3   | 105   | 0.00     |
| 42   | 2-Nitroaniline              | 0.343 | 0.360 | -5.0  | 115   | 0.01     |
| 43 S | Dimethylphthalate-d6        | 1.319 | 1.274 | 3.4   | 103   | 0.00     |
| 44   | Dimethylphthalate           | 1.319 | 1.300 | 1.4   | 104   | 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                    | AvqRF | CCRF  | %Dev | Area% | Dev(min) |
|------|-----------------------------|-------|-------|------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.230 | 0.226 | 1.7  | 104   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.556 | 1.542 | 0.9  | 105   | 0.00     |
| 47   | Acenaphthylene              | 1.527 | 1.463 | 4.2  | 101   | 0.00     |
| 48   | 3-Nitroaniline              | 0.233 | 0.239 | -2.6 | 110   | 0.00     |
| 49 C | Acenaphthene                | 1.056 | 1.025 | 2.9  | 104   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.160 | 0.147 | 8.1  | 108   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.216 | 0.198 | 8.3  | 104   | 0.00     |
| 52   | 4-Nitrophenol               | 0.359 | 0.344 | 4.2  | 109   | 0.00     |
| 53   | Dibenzofuran                | 1.606 | 1.594 | 0.7  | 105   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.349 | 0.363 | -4.0 | 113   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.369 | 0.373 | -1.1 | 107   | 0.00     |
| 56   | Diethylphthalate            | 1.388 | 1.357 | 2.2  | 103   | 0.00     |
| 57 S | Fluorene-d10                | 1.217 | 1.178 | 3.2  | 103   | 0.00     |
| 58   | Fluorene                    | 1.316 | 1.297 | 1.4  | 106   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.747 | 0.740 | 0.9  | 105   | 0.00     |
| 60   | 4-Nitroaniline              | 0.255 | 0.260 | -2.0 | 111   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0  | 108   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.103 | 0.096 | 6.8  | 111   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.109 | 0.101 | 7.3  | 112   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.502 | 0.486 | 3.2  | 105   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.207 | 0.203 | 1.9  | 108   | 0.00     |
| 66   | Hexachlorobenzene           | 0.232 | 0.231 | 0.4  | 107   | 0.00     |
| 67   | Atrazine                    | 0.223 | 0.205 | 8.1  | 101   | 0.00     |
| 68 C | Pentachlorophenol           | 0.146 | 0.136 | 6.8  | 107   | 0.00     |
| 69   | Phenanthrene                | 0.963 | 0.944 | 2.0  | 105   | 0.00     |
| 70 S | Anthracene-d10              | 0.862 | 0.848 | 1.6  | 107   | 0.00     |
| 71   | Anthracene                  | 0.990 | 0.976 | 1.4  | 107   | 0.00     |
| 72   | Carbazole                   | 0.881 | 0.855 | 3.0  | 107   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.035 | 0.984 | 4.9  | 104   | 0.00     |
| 74 C | Fluoranthene                | 1.238 | 1.209 | 2.3  | 108   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0  | 117   | 0.00     |
| 76 S | Pyrene-d10                  | 0.881 | 0.812 | 7.8  | 108   | 0.00     |
| 77   | Pyrene                      | 1.076 | 1.001 | 7.0  | 109   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.396 | 0.356 | 10.1 | 106   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.381 | 0.377 | 1.0  | 112   | 0.01     |
| 80   | Benzo(a)anthracene          | 1.060 | 1.031 | 2.7  | 113   | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.554 | 0.513 | 7.4  | 113   | 0.00     |
| 82   | Chrysene                    | 0.984 | 0.993 | -0.9 | 118   | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0  | 115   | 0.01     |
| 84   | Di-n-octyl phthalate        | 1.041 | 0.935 | 10.2 | 113   | 0.01     |
| 85   | Benzo(b)fluoranthene        | 1.066 | 1.057 | 0.8  | 117   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.031 | 1.041 | -1.0 | 119   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.816 | 0.811 | 0.6  | 117   | 0.01     |

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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.017 | 1.024 | -0.7 | 119   | 0.01     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.153 | 1.132 | 1.8  | 115   | 0.01     |
| 90   | Dibenzo(a,h)anthracene | 0.985 | 0.971 | 1.4  | 116   | 0.01     |
| 91   | Benzo(a,h,i)perylene   | 0.978 | 0.941 | 3.8  | 112   | 0.01     |

(#) = Out of Range

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