

Data Path : Z:\HPCHEM1\BNA\_M\Data\BM061317\  
 Data File : BM010575.D  
 Acq On : 14 Jun 2017 00:09  
 Operator : SJ/MA  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02030

Quant Time: Jun 14 09:12:02 2017  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM052717.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Jun 14 01:46:30 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                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0    | 32    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.520 | 0.400 | 23.1#  | 26    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.487 | 0.412 | 15.4   | 27    | 0.00     |
| 4    | Benzaldehyde                | 1.019 | 1.134 | -11.3  | 34    | 0.00     |
| 5 S  | Phenol-d5                   | 1.516 | 1.287 | 15.1   | 28    | 0.00     |
| 6    | Phenol                      | 1.554 | 1.260 | 18.9   | 27    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.852 | 0.745 | 12.6   | 29    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.078 | 0.901 | 16.4   | 28    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.208 | 1.108 | 8.3    | 29    | 0.00     |
| 10   | 2-Chlorophenol              | 1.212 | 1.091 | 10.0   | 29    | 0.00     |
| 11   | 2-Methylphenol              | 1.134 | 1.000 | 11.8   | 29    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 0.486 | 0.346 | 28.8#  | 23    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.223 | 1.053 | 13.9   | 29    | 0.00     |
| 14   | Acetophenone                | 1.995 | 1.756 | 12.0   | 29    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 0.997 | 0.988 | 0.9    | 31    | 0.00     |
| 16   | 4-Methylphenol              | 1.244 | 1.093 | 12.1   | 29    | 0.00     |
| 17   | Hexachloroethane            | 0.550 | 0.630 | -14.5  | 38    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 30    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.147 | 0.145 | 1.4    | 29    | 0.00     |
| 20   | Nitrobenzene                | 0.418 | 0.483 | -15.6  | 35    | 0.00     |
| 21   | Isophorone                  | 0.644 | 0.683 | -6.1   | 32    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.170 | 0.191 | -12.4  | 34    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.180 | 0.192 | -6.7   | 32    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.418 | 0.433 | -3.6   | 31    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.362 | 0.350 | 3.3    | 29    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.349 | 0.411 | -17.8  | 36    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.342 | 0.406 | -18.7  | 35    | 0.00     |
| 28   | Naphthalene                 | 1.003 | 0.961 | 4.2    | 29    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.336 | 0.369 | -9.8   | 36    | 0.00     |
| 30   | 4-Chloroaniline             | 0.325 | 0.355 | -9.2   | 34    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.313 | 0.404 | -29.1# | 39    | 0.00     |
| 32   | Caprolactam                 | 0.085 | 0.095 | -11.8  | 38    | 0.01     |
| 33 C | 4-Chloro-3-methylphenol     | 0.329 | 0.368 | -11.9  | 34    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.761 | 0.819 | -7.6   | 32    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 35    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.832 | 0.919 | -10.5  | 39    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.558 | 0.501 | 10.2   | 33    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.473 | 0.524 | -10.8  | 40    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.477 | 0.587 | -23.1# | 43    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.622 | 1.629 | -0.4   | 35    | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.275 | 1.262 | 1.0    | 36    | 0.00     |
| 42   | 2-Nitroaniline              | 0.325 | 0.376 | -15.7  | 41    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.662 | 1.853 | -11.5  | 41    | 0.00     |
| 44   | Dimethylphthalate           | 1.634 | 1.743 | -6.7   | 39    | 0.00     |

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Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
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|      | Compound                    | AvgRF | CCRF  | %Dev   | Area% | Dev(min) |
|------|-----------------------------|-------|-------|--------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.300 | 0.328 | -9.3   | 39    | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.939 | 1.971 | -1.7   | 36    | 0.00     |
| 47   | Acenaphthylene              | 1.894 | 1.893 | 0.1    | 36    | 0.00     |
| 48   | 3-Nitroaniline              | 0.289 | 0.275 | 4.8    | 37    | 0.00     |
| 49 C | Acenaphthene                | 1.329 | 1.303 | 2.0    | 35    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.226 | 0.230 | -1.8   | 43    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.249 | 0.248 | 0.4    | 41    | 0.00     |
| 52   | 4-Nitrophenol               | 0.341 | 0.469 | -37.5# | 59    | 0.00     |
| 53   | Dibenzofuran                | 1.965 | 2.133 | -8.5   | 39    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.440 | 0.510 | -15.9  | 42    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.458 | 0.586 | -27.9# | 47    | 0.00     |
| 56   | Diethylphthalate            | 1.581 | 1.735 | -9.7   | 41    | 0.00     |
| 57 S | Fluorene-d10                | 1.461 | 1.603 | -9.7   | 40    | 0.00     |
| 58   | Fluorene                    | 1.611 | 1.746 | -8.4   | 40    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.921 | 1.095 | -18.9  | 43    | 0.00     |
| 60   | 4-Nitroaniline              | 0.302 | 0.289 | 4.3    | 41    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 44    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.142 | 0.137 | 3.5    | 49    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.150 | 0.138 | 8.0    | 45    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.583 | 0.547 | 6.2    | 41    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.246 | 0.250 | -1.6   | 44    | 0.00     |
| 66   | Hexachlorobenzene           | 0.258 | 0.242 | 6.2    | 41    | 0.00     |
| 67   | Atrazine                    | 0.253 | 0.271 | -7.1   | 52    | 0.00     |
| 68 C | Pentachlorophenol           | 0.164 | 0.157 | 4.3    | 47    | 0.00     |
| 69   | Phenanthrene                | 1.068 | 1.059 | 0.8    | 44    | 0.00     |
| 70 S | Anthracene-d10              | 0.973 | 0.970 | 0.3    | 44    | 0.00     |
| 71   | Anthracene                  | 1.115 | 1.096 | 1.7    | 44    | 0.00     |
| 72   | Carbazole                   | 0.943 | 0.889 | 5.7    | 44    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.065 | 1.078 | -1.2   | 46    | 0.00     |
| 74 C | Fluoranthene                | 1.313 | 1.444 | -10.0  | 50    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 48    | 0.00     |
| 76 S | Pyrene-d10                  | 0.827 | 0.860 | -4.0   | 50    | 0.00     |
| 77   | Pyrene                      | 1.031 | 1.026 | 0.5    | 48    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.356 | 0.359 | -0.8   | 51    | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.393 | 0.384 | 2.3    | 49    | 0.00     |
| 80   | Benzo(a)anthracene          | 1.128 | 1.157 | -2.6   | 51    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.530 | 0.519 | 2.1    | 50    | 0.00     |
| 82   | Chrysene                    | 1.020 | 1.035 | -1.5   | 50    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 46    | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.921 | 0.933 | -1.3   | 47    | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.164 | 1.254 | -7.7   | 51    | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.112 | 1.124 | -1.1   | 47    | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.931 | 0.932 | -0.1   | 47    | 0.00     |

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| 88 C | Benzo(a)pyrene         | 1.154 | 1.174 | -1.7 | 48    | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.457 | 1.471 | -1.0 | 48    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.255 | 1.261 | -0.5 | 48    | 0.00     |
| 91   | Benzo(g,h,i)perylene   | 1.201 | 1.173 | 2.3  | 46    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 1