

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042816\  
 Data File : BM005133.D  
 Acq On : 29 Apr 2016 02:54  
 Operator : UM/SJ  
 Sample : SSTDC0020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02051

Quant Time: Apr 29 05:40:20 2016  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM042516.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Apr 29 02:20:37 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    | 113   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.730 | 0.759 | -4.0   | 115   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.394 | 0.415 | -5.3   | 116   | 0.00     |
| 4    | Benzaldehyde                | 0.821 | 1.068 | -30.1# | 116   | 0.00     |
| 5 S  | Phenol-d5                   | 1.654 | 1.779 | -7.6   | 118   | 0.00     |
| 6    | Phenol                      | 1.725 | 1.855 | -7.5   | 117   | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.941 | 1.021 | -8.5   | 117   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.322 | 1.411 | -6.7   | 116   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.309 | 1.385 | -5.8   | 116   | 0.00     |
| 10   | 2-Chlorophenol              | 1.345 | 1.430 | -6.3   | 117   | 0.00     |
| 11   | 2-Methylphenol              | 1.346 | 1.424 | -5.8   | 117   | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.710 | 1.839 | -7.5   | 116   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.403 | 1.467 | -4.6   | 116   | 0.00     |
| 14   | Acetophenone                | 2.111 | 2.309 | -9.4   | 114   | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.057 | 1.161 | -9.8   | 117   | 0.00     |
| 16   | 4-Methylphenol              | 1.498 | 1.603 | -7.0   | 117   | 0.00     |
| 17   | Hexachloroethane            | 0.536 | 0.542 | -1.1   | 112   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 111   | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.139 | 0.148 | -6.5   | 115   | 0.00     |
| 20   | Nitrobenzene                | 0.341 | 0.357 | -4.7   | 113   | 0.00     |
| 21   | Isophorone                  | 0.647 | 0.708 | -9.4   | 117   | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.155 | 0.171 | -10.3  | 119   | 0.00     |
| 23 C | 2-Nitrophenol               | 0.168 | 0.184 | -9.5   | 116   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.362 | 0.378 | -4.4   | 113   | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.402 | 0.431 | -7.2   | 117   | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.294 | 0.321 | -9.2   | 117   | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.302 | 0.327 | -8.3   | 115   | 0.00     |
| 28   | Naphthalene                 | 1.009 | 1.037 | -2.8   | 112   | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.344 | 0.427 | -24.1# | 115   | 0.00     |
| 30   | 4-Chloroaniline             | 0.350 | 0.432 | -23.4# | 115   | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.199 | 0.191 | 4.0    | 106   | 0.00     |
| 32   | Caprolactam                 | 0.104 | 0.108 | -3.8   | 115   | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.345 | 0.385 | -11.6  | 118   | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.747 | 0.778 | -4.1   | 112   | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 109   | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.625 | 0.629 | -0.6   | 107   | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.371 | 0.338 | 8.9    | 103   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.397 | 0.431 | -8.6   | 114   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.445 | 0.481 | -8.1   | 113   | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.582 | 1.651 | -4.4   | 110   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.203 | 1.248 | -3.7   | 110   | 0.00     |
| 42   | 2-Nitroaniline              | 0.326 | 0.378 | -16.0  | 117   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.580 | 1.636 | -3.5   | 109   | 0.00     |
| 44   | Dimethylphthalate           | 1.584 | 1.642 | -3.7   | 109   | 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          | 0.306 | 0.335 | -9.5   | 113   | 0.00     |
| 46 S | Acenaphthylene-d8           | 1.881 | 1.960 | -4.2   | 109   | 0.00     |
| 47   | Acenaphthylene              | 1.990 | 2.093 | -5.2   | 109   | 0.00     |
| 48   | 3-Nitroaniline              | 0.311 | 0.364 | -17.0  | 115   | 0.00     |
| 49 C | Acenaphthene                | 1.309 | 1.360 | -3.9   | 109   | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.179 | 0.142 | 20.7#  | 101   | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.276 | 0.293 | -6.2   | 113   | 0.00     |
| 52   | 4-Nitrophenol               | 0.247 | 0.243 | 1.6    | 107   | 0.00     |
| 53   | Dibenzofuran                | 1.931 | 1.979 | -2.5   | 107   | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.465 | 0.498 | -7.1   | 110   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.387 | 0.409 | -5.7   | 110   | 0.00     |
| 56   | Diethylphthalate            | 1.599 | 1.679 | -5.0   | 110   | 0.00     |
| 57 S | Fluorene-d10                | 1.388 | 1.425 | -2.7   | 108   | 0.00     |
| 58   | Fluorene                    | 1.499 | 1.563 | -4.3   | 108   | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.756 | 0.781 | -3.3   | 107   | 0.00     |
| 60   | 4-Nitroaniline              | 0.336 | 0.373 | -11.0  | 115   | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0    | 105   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.114 | 0.108 | 5.3    | 100   | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.120 | 0.115 | 4.2    | 99    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.550 | 0.595 | -8.2   | 109   | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.195 | 0.202 | -3.6   | 106   | 0.00     |
| 66   | Hexachlorobenzene           | 0.222 | 0.226 | -1.8   | 104   | 0.00     |
| 67   | Atrazine                    | 0.203 | 0.223 | -9.9   | 109   | 0.00     |
| 68 C | Pentachlorophenol           | 0.128 | 0.130 | -1.6   | 109   | 0.00     |
| 69   | Phenanthrene                | 1.064 | 1.098 | -3.2   | 106   | 0.00     |
| 70 S | Anthracene-d10              | 0.897 | 0.939 | -4.7   | 106   | 0.00     |
| 71   | Anthracene                  | 1.073 | 1.119 | -4.3   | 105   | 0.00     |
| 72   | Carbazole                   | 0.920 | 1.022 | -11.1  | 107   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.077 | 1.227 | -13.9  | 113   | 0.00     |
| 74 C | Fluoranthene                | 1.171 | 1.261 | -7.7   | 101   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0    | 90    | -0.01    |
| 76 S | Pyrene-d10                  | 0.910 | 1.030 | -13.2  | 100   | 0.00     |
| 77   | Pyrene                      | 1.155 | 1.304 | -12.9  | 99    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.444 | 0.554 | -24.8# | 107   | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.358 | 0.371 | -3.6   | 82    | -0.01    |
| 80   | Benzo(a)anthracene          | 1.145 | 1.191 | -4.0   | 90    | -0.01    |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.624 | 0.764 | -22.4# | 104   | -0.01    |
| 82   | Chrysene                    | 1.095 | 1.136 | -3.7   | 91    | -0.01    |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0    | 78    | -0.01    |
| 84   | Di-n-octyl phthalate        | 1.243 | 1.501 | -20.8# | 93    | -0.01    |
| 85   | Benzo(b)fluoranthene        | 1.205 | 1.308 | -8.5   | 83    | -0.01    |
| 86   | Benzo(k)fluoranthene        | 1.171 | 1.175 | -0.3   | 76    | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.896 | 0.928 | -3.6   | 78    | -0.01    |

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|      | Compound               | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|------------------------|-------|-------|-------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.125 | 1.170 | -4.0  | 78    | -0.01    |
| 89   | Indeno(1,2,3-cd)pyrene | 1.073 | 1.173 | -9.3  | 80    | -0.02    |
| 90   | Dibenzo(a,h)anthracene | 0.897 | 0.989 | -10.3 | 80    | -0.02    |
| 91   | Benzo(a,h,i)perylene   | 0.878 | 0.965 | -9.9  | 81    | -0.02    |

(#) = Out of Range

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