

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020817\  
 Data File : BM008986.D  
 Acq On : 08 Feb 2017 14:39  
 Operator : SJ/MA  
 Sample : SSTDICV020  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 SICV41

Quant Time: Feb 08 15:13:39 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM020817.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Wed Feb 08 15:01:57 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   | 113   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.575 | 0.556 | 3.3   | 110   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.520 | 0.505 | 2.9   | 103   | 0.00     |
| 4    | Benzaldehyde                | 1.371 | 1.505 | -9.8  | 113   | -0.01    |
| 5 S  | Phenol-d5                   | 1.777 | 1.787 | -0.6  | 112   | -0.01    |
| 6    | Phenol                      | 1.909 | 1.900 | 0.5   | 108   | -0.01    |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.293 | 1.316 | -1.8  | 110   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.497 | 1.563 | -4.4  | 113   | -0.01    |
| 9 S  | 2-Chlorophenol-d4           | 1.194 | 1.228 | -2.8  | 112   | -0.01    |
| 10   | 2-Chlorophenol              | 1.251 | 1.291 | -3.2  | 111   | -0.01    |
| 11   | 2-Methylphenol              | 1.320 | 1.306 | 1.1   | 106   | -0.01    |
| 12   | 2,2'-oxybis(1-Chloropropane | 2.397 | 2.456 | -2.5  | 111   | -0.01    |
| 13 S | 4-Methylphenol-d8           | 1.327 | 1.342 | -1.1  | 113   | -0.01    |
| 14   | Acetophenone                | 2.307 | 2.336 | -1.3  | 110   | -0.01    |
| 15 P | N-Nitroso-di-n-propylamine  | 1.327 | 1.378 | -3.8  | 109   | -0.01    |
| 16   | 4-Methylphenol              | 1.441 | 1.448 | -0.5  | 108   | 0.00     |
| 17   | Hexachloroethane            | 0.607 | 0.612 | -0.8  | 112   | -0.01    |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 109   | -0.01    |
| 19 S | Nitrobenzene-d5             | 0.137 | 0.146 | -6.6  | 111   | 0.00     |
| 20   | Nitrobenzene                | 0.478 | 0.495 | -3.6  | 109   | -0.01    |
| 21   | Isophorone                  | 0.827 | 0.846 | -2.3  | 108   | -0.01    |
| 22 S | 2-Nitrophenol-d4            | 0.150 | 0.161 | -7.3  | 110   | -0.01    |
| 23 C | 2-Nitrophenol               | 0.165 | 0.175 | -6.1  | 111   | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.417 | 0.440 | -5.5  | 111   | -0.02    |
| 25   | Bis(2-Chloroethoxy)methane  | 0.478 | 0.500 | -4.6  | 109   | -0.01    |
| 26 S | 2,4-Dichlorophenol-d3       | 0.320 | 0.328 | -2.5  | 108   | -0.01    |
| 27 C | 2,4-Dichlorophenol          | 0.320 | 0.325 | -1.6  | 106   | 0.00     |
| 28   | Naphthalene                 | 0.997 | 1.035 | -3.8  | 110   | -0.01    |
| 29 S | 4-Chloroaniline-d4          | 0.325 | 0.360 | -10.8 | 110   | -0.01    |
| 30   | 4-Chloroaniline             | 0.341 | 0.374 | -9.7  | 110   | -0.02    |
| 31 C | Hexachlorobutadiene         | 0.263 | 0.265 | -0.8  | 107   | -0.01    |
| 32   | Caprolactam                 | 0.097 | 0.094 | 3.1   | 103   | -0.01    |
| 33 C | 4-Chloro-3-methylphenol     | 0.363 | 0.373 | -2.8  | 108   | -0.01    |
| 34   | 2-Methylnaphthalene         | 0.742 | 0.754 | -1.6  | 108   | -0.01    |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 110   | -0.01    |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.653 | 0.661 | -1.2  | 108   | -0.01    |
| 37   | Hexachlorocyclopentadiene   | 0.417 | 0.393 | 5.8   | 109   | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.372 | 0.384 | -3.2  | 110   | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.406 | 0.424 | -4.4  | 112   | -0.01    |
| 40   | 1,1'-Biphenyl               | 1.388 | 1.389 | -0.1  | 106   | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.093 | 1.106 | -1.2  | 109   | -0.01    |
| 42   | 2-Nitroaniline              | 0.371 | 0.383 | -3.2  | 107   | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.368 | 1.378 | -0.7  | 106   | -0.01    |
| 44   | Dimethylphthalate           | 1.368 | 1.376 | -0.6  | 107   | -0.01    |

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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          | 0.249 | 0.255 | -2.4 | 107   | -0.01    |
| 46 S | Acenaphthylene-d8           | 1.652 | 1.691 | -2.4 | 108   | -0.01    |
| 47   | Acenaphthylene              | 1.639 | 1.685 | -2.8 | 109   | -0.01    |
| 48   | 3-Nitroaniline              | 0.248 | 0.249 | -0.4 | 105   | 0.00     |
| 49 C | Acenaphthene                | 1.171 | 1.174 | -0.3 | 107   | -0.01    |
| 50   | 2,4-Dinitrophenol           | 0.176 | 0.156 | 11.4 | 107   | -0.01    |
| 51 S | 4-Nitrophenol-d4            | 0.241 | 0.232 | 3.7  | 108   | -0.01    |
| 52   | 4-Nitrophenol               | 0.290 | 0.278 | 4.1  | 104   | -0.01    |
| 53   | Dibenzofuran                | 1.738 | 1.719 | 1.1  | 105   | -0.01    |
| 54   | 2,4-Dinitrotoluene          | 0.379 | 0.397 | -4.7 | 110   | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.377 | 0.380 | -0.8 | 108   | 0.00     |
| 56   | Diethylphthalate            | 1.405 | 1.419 | -1.0 | 108   | 0.00     |
| 57 S | Fluorene-d10                | 1.271 | 1.269 | 0.2  | 106   | 0.00     |
| 58   | Fluorene                    | 1.424 | 1.410 | 1.0  | 105   | -0.01    |
| 59   | 4-Chlorophenyl-phenylether  | 0.767 | 0.769 | -0.3 | 106   | 0.00     |
| 60   | 4-Nitroaniline              | 0.287 | 0.291 | -1.4 | 109   | -0.01    |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0  | 105   | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.119 | 0.117 | 1.7  | 108   | -0.01    |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.128 | 0.121 | 5.5  | 105   | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.568 | 0.579 | -1.9 | 105   | -0.01    |
| 65   | 4-Bromophenyl-phenylether   | 0.224 | 0.229 | -2.2 | 106   | 0.00     |
| 66   | Hexachlorobenzene           | 0.252 | 0.250 | 0.8  | 105   | 0.00     |
| 67   | Atrazine                    | 0.225 | 0.224 | 0.4  | 105   | -0.01    |
| 68 C | Pentachlorophenol           | 0.151 | 0.144 | 4.6  | 106   | -0.01    |
| 69   | Phenanthrene                | 1.085 | 1.103 | -1.7 | 105   | -0.01    |
| 70 S | Anthracene-d10              | 0.941 | 0.942 | -0.1 | 104   | 0.00     |
| 71   | Anthracene                  | 1.109 | 1.116 | -0.6 | 103   | 0.00     |
| 72   | Carbazole                   | 0.987 | 0.985 | 0.2  | 105   | 0.00     |
| 73   | Di-n-butylphthalate         | 1.106 | 1.151 | -4.1 | 107   | -0.01    |
| 74 C | Fluoranthene                | 1.323 | 1.354 | -2.3 | 104   | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0  | 104   | -0.01    |
| 76 S | Pyrene-d10                  | 0.857 | 0.870 | -1.5 | 102   | 0.00     |
| 77   | Pyrene                      | 1.099 | 1.112 | -1.2 | 104   | 0.00     |
| 78   | Butylbenzylphthalate        | 0.388 | 0.405 | -4.4 | 107   | -0.01    |
| 79   | 3,3'-Dichlorobenzidine      | 0.379 | 0.399 | -5.3 | 105   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.133 | 1.173 | -3.5 | 105   | -0.01    |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.570 | 0.593 | -4.0 | 107   | -0.01    |
| 82   | Chrysene                    | 1.088 | 1.101 | -1.2 | 103   | -0.01    |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0  | 101   | -0.01    |
| 84   | Di-n-octyl phthalate        | 1.102 | 1.128 | -2.4 | 106   | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.196 | 1.233 | -3.1 | 101   | -0.01    |
| 86   | Benzo(k)fluoranthene        | 1.137 | 1.194 | -5.0 | 103   | -0.01    |
| 87 S | Benzo(a)pyrene-d12          | 0.901 | 0.928 | -3.0 | 101   | -0.01    |

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|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.136 | 1.174 | -3.3 | 100   | -0.01    |
| 89   | Indeno(1,2,3-cd)pyrene | 1.251 | 1.276 | -2.0 | 100   | -0.01    |
| 90   | Dibenzo(a,h)anthracene | 1.078 | 1.085 | -0.6 | 100   | -0.02    |
| 91   | Benzo(a,h,i)perylene   | 1.051 | 1.061 | -1.0 | 100   | -0.02    |

(#) = Out of Range

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