

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM092016\  
 Data File : BM007491.D  
 Acq On : 20 Sep 2016 09:56  
 Operator : UM/SJ  
 Sample : SSTDCCC020  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02059

Quant Time: Sep 20 16:25:55 2016  
 Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM082316.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Sep 20 01:35:12 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                    | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|-----------------------------|-------|-------|------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0  | 103   | 0.00     |
| 2    | 1,4-Dioxane                 | 0.436 | 0.471 | -8.0 | 113   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.391 | 0.424 | -8.4 | 113   | 0.00     |
| 4    | Benzaldehyde                | 1.076 | 1.163 | -8.1 | 98    | 0.00     |
| 5 S  | Phenol-d5                   | 1.889 | 1.810 | 4.2  | 99    | 0.00     |
| 6    | Phenol                      | 1.960 | 1.894 | 3.4  | 99    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 0.993 | 1.007 | -1.4 | 101   | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.365 | 1.376 | -0.8 | 101   | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.396 | 1.371 | 1.8  | 101   | 0.00     |
| 10   | 2-Chlorophenol              | 1.430 | 1.434 | -0.3 | 101   | 0.00     |
| 11   | 2-Methylphenol              | 1.532 | 1.461 | 4.6  | 98    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.536 | 1.554 | -1.2 | 100   | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.648 | 1.556 | 5.6  | 97    | 0.00     |
| 14   | Acetophenone                | 2.527 | 2.475 | 2.1  | 97    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.343 | 1.313 | 2.2  | 97    | 0.00     |
| 16   | 4-Methylphenol              | 1.718 | 1.646 | 4.2  | 97    | 0.00     |
| 17   | Hexachloroethane            | 0.556 | 0.571 | -2.7 | 104   | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0  | 99    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.147 | 0.148 | -0.7 | 98    | 0.00     |
| 20   | Nitrobenzene                | 0.377 | 0.385 | -2.1 | 98    | 0.00     |
| 21   | Isophorone                  | 0.768 | 0.761 | 0.9  | 94    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.171 | 0.172 | -0.6 | 98    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.186 | 0.187 | -0.5 | 95    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.411 | 0.423 | -2.9 | 99    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.433 | 0.433 | 0.0  | 96    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.341 | 0.340 | 0.3  | 96    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.339 | 0.344 | -1.5 | 97    | 0.00     |
| 28   | Naphthalene                 | 0.995 | 1.019 | -2.4 | 98    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.426 | 0.431 | -1.2 | 93    | 0.00     |
| 30   | 4-Chloroaniline             | 0.438 | 0.441 | -0.7 | 92    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.229 | 0.241 | -5.2 | 101   | 0.00     |
| 32   | Caprolactam                 | 0.137 | 0.121 | 11.7 | 82    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.420 | 0.417 | 0.7  | 94    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.814 | 0.816 | -0.2 | 96    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0  | 95    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.666 | 0.704 | -5.7 | 97    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.402 | 0.340 | 15.4 | 81    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.471 | 0.460 | 2.3  | 89    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.491 | 0.482 | 1.8  | 91    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.551 | 1.592 | -2.6 | 94    | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.216 | 1.234 | -1.5 | 93    | 0.00     |
| 42   | 2-Nitroaniline              | 0.397 | 0.405 | -2.0 | 93    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.725 | 1.709 | 0.9  | 92    | 0.00     |
| 44   | Dimethylphthalate           | 1.719 | 1.695 | 1.4  | 90    | 0.00     |

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 LabSampleId :  
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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          | 0.351 | 0.347 | 1.1  | 89    | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.000 | 2.001 | -0.0 | 92    | 0.00     |
| 47   | Acenaphthylene              | 2.058 | 2.099 | -2.0 | 93    | 0.00     |
| 48   | 3-Nitroaniline              | 0.359 | 0.347 | 3.3  | 84    | 0.00     |
| 49 C | Acenaphthene                | 1.339 | 1.366 | -2.0 | 94    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.238 | 0.205 | 13.9 | 86    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.326 | 0.294 | 9.8  | 82    | 0.00     |
| 52   | 4-Nitrophenol               | 0.347 | 0.322 | 7.2  | 85    | 0.00     |
| 53   | Dibenzofuran                | 2.004 | 2.012 | -0.4 | 92    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.538 | 0.525 | 2.4  | 89    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.504 | 0.497 | 1.4  | 90    | 0.00     |
| 56   | Diethylphthalate            | 1.793 | 1.740 | 3.0  | 89    | 0.00     |
| 57 S | Fluorene-d10                | 1.492 | 1.495 | -0.2 | 91    | 0.00     |
| 58   | Fluorene                    | 1.650 | 1.668 | -1.1 | 93    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.863 | 0.866 | -0.3 | 92    | 0.00     |
| 60   | 4-Nitroaniline              | 0.407 | 0.380 | 6.6  | 82    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0  | 91    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.126 | 0.122 | 3.2  | 88    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.134 | 0.136 | -1.5 | 91    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.558 | 0.583 | -4.5 | 91    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.226 | 0.234 | -3.5 | 90    | 0.00     |
| 66   | Hexachlorobenzene           | 0.253 | 0.255 | -0.8 | 89    | 0.00     |
| 67   | Atrazine                    | 0.236 | 0.242 | -2.5 | 85    | 0.00     |
| 68 C | Pentachlorophenol           | 0.163 | 0.155 | 4.9  | 84    | 0.00     |
| 69   | Phenanthrene                | 1.074 | 1.110 | -3.4 | 90    | 0.00     |
| 70 S | Anthracene-d10              | 0.934 | 0.961 | -2.9 | 90    | 0.00     |
| 71   | Anthracene                  | 1.108 | 1.154 | -4.2 | 90    | 0.00     |
| 72   | Carbazole                   | 0.963 | 1.021 | -6.0 | 88    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.212 | 1.202 | 0.8  | 86    | 0.00     |
| 74 C | Fluoranthene                | 1.347 | 1.464 | -8.7 | 89    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0  | 91    | 0.00     |
| 76 S | Pyrene-d10                  | 0.836 | 0.846 | -1.2 | 90    | 0.00     |
| 77   | Pyrene                      | 1.073 | 1.094 | -2.0 | 90    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.446 | 0.434 | 2.7  | 86    | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.397 | 0.421 | -6.0 | 84    | 0.00     |
| 80   | Benzo(a)anthracene          | 1.179 | 1.196 | -1.4 | 90    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.648 | 0.642 | 0.9  | 87    | 0.00     |
| 82   | Chrysene                    | 1.068 | 1.089 | -2.0 | 89    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0  | 88    | 0.00     |
| 84   | Di-n-octyl phthalate        | 0.996 | 1.086 | -9.0 | 86    | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.183 | 1.193 | -0.8 | 84    | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.093 | 1.150 | -5.2 | 90    | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.921 | 0.931 | -1.1 | 87    | 0.00     |

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| 88 C | Benzo(a)pyrene         | 1.132 | 1.162 | -2.7 | 87    | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.389 | 1.328 | 4.4  | 81    | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.155 | 1.096 | 5.1  | 80    | 0.00     |
| 91   | Benzo(g,h,i)perylene   | 1.179 | 1.112 | 5.7  | 80    | 0.00     |

(#) = Out of Range

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