

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100317\  
 Data File : BM011853.D  
 Acq On : 03 Oct 2017 16:32  
 Operator : SJ/JU  
 Sample : SSTDC0020  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTD02016

Quant Time: Oct 03 18:26:17 2017  
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM100317.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Oct 03 15:47:54 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    | 87    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.430 | 0.459 | -6.7   | 86    | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.423 | 0.423 | 0.0    | 82    | 0.00     |
| 4    | Benzaldehyde                | 0.857 | 0.931 | -8.6   | 84    | 0.00     |
| 5 S  | Phenol-d5                   | 1.721 | 1.705 | 0.9    | 86    | 0.00     |
| 6    | Phenol                      | 1.703 | 1.710 | -0.4   | 86    | 0.00     |
| 7 S  | Bis-(2-Chloroethyl)ether-d8 | 1.018 | 1.032 | -1.4   | 84    | 0.00     |
| 8    | Bis(2-Chloroethyl)ether     | 1.279 | 1.309 | -2.3   | 85    | 0.00     |
| 9 S  | 2-Chlorophenol-d4           | 1.380 | 1.455 | -5.4   | 86    | 0.00     |
| 10   | 2-Chlorophenol              | 1.349 | 1.422 | -5.4   | 87    | 0.00     |
| 11   | 2-Methylphenol              | 1.295 | 1.318 | -1.8   | 87    | 0.00     |
| 12   | 2,2'-oxybis(1-Chloropropane | 1.635 | 1.672 | -2.3   | 83    | 0.00     |
| 13 S | 4-Methylphenol-d8           | 1.487 | 1.509 | -1.5   | 86    | 0.00     |
| 14   | Acetophenone                | 2.177 | 2.254 | -3.5   | 86    | 0.00     |
| 15 P | N-Nitroso-di-n-propylamine  | 1.114 | 1.173 | -5.3   | 84    | 0.00     |
| 16   | 4-Methylphenol              | 1.450 | 1.483 | -2.3   | 86    | 0.00     |
| 17   | Hexachloroethane            | 0.541 | 0.572 | -5.7   | 87    | 0.00     |
| 18 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0    | 87    | 0.00     |
| 19 S | Nitrobenzene-d5             | 0.142 | 0.151 | -6.3   | 86    | 0.00     |
| 20   | Nitrobenzene                | 0.344 | 0.363 | -5.5   | 85    | 0.00     |
| 21   | Isophorone                  | 0.642 | 0.671 | -4.5   | 85    | 0.00     |
| 22 S | 2-Nitrophenol-d4            | 0.160 | 0.173 | -8.1   | 89    | 0.00     |
| 23 C | 2-Nitrophenol               | 0.169 | 0.183 | -8.3   | 88    | 0.00     |
| 24   | 2,4-Dimethylphenol          | 0.362 | 0.378 | -4.4   | 87    | 0.00     |
| 25   | Bis(2-Chloroethoxy)methane  | 0.394 | 0.411 | -4.3   | 85    | 0.00     |
| 26 S | 2,4-Dichlorophenol-d3       | 0.324 | 0.348 | -7.4   | 87    | 0.00     |
| 27 C | 2,4-Dichlorophenol          | 0.313 | 0.332 | -6.1   | 87    | 0.00     |
| 28   | Naphthalene                 | 1.003 | 1.038 | -3.5   | 87    | 0.00     |
| 29 S | 4-Chloroaniline-d4          | 0.354 | 0.429 | -21.2# | 87    | 0.00     |
| 30   | 4-Chloroaniline             | 0.347 | 0.414 | -19.3  | 86    | 0.00     |
| 31 C | Hexachlorobutadiene         | 0.230 | 0.240 | -4.3   | 87    | 0.00     |
| 32   | Caprolactam                 | 0.103 | 0.101 | 1.9    | 87    | 0.00     |
| 33 C | 4-Chloro-3-methylphenol     | 0.340 | 0.361 | -6.2   | 87    | 0.00     |
| 34   | 2-Methylnaphthalene         | 0.765 | 0.802 | -4.8   | 87    | 0.00     |
| 35 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0    | 89    | 0.00     |
| 36   | 1,2,4,5-Tetrachlorobenzene  | 0.678 | 0.705 | -4.0   | 89    | 0.00     |
| 37   | Hexachlorocyclopentadiene   | 0.395 | 0.367 | 7.1    | 85    | 0.00     |
| 38 C | 2,4,6-Trichlorophenol       | 0.431 | 0.456 | -5.8   | 89    | 0.00     |
| 39   | 2,4,5-Trichlorophenol       | 0.454 | 0.482 | -6.2   | 90    | 0.00     |
| 40   | 1,1'-Biphenyl               | 1.599 | 1.623 | -1.5   | 87    | 0.00     |
| 41   | 2-Chloronaphthalene         | 1.251 | 1.297 | -3.7   | 89    | 0.00     |
| 42   | 2-Nitroaniline              | 0.318 | 0.339 | -6.6   | 86    | 0.00     |
| 43 S | Dimethylphthalate-d6        | 1.733 | 1.815 | -4.7   | 89    | 0.00     |
| 44   | Dimethylphthalate           | 1.668 | 1.740 | -4.3   | 89    | 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 45   | 2,6-Dinitrotoluene          | 0.314 | 0.341 | -8.6  | 88    | 0.00     |
| 46 S | Acenaphthylene-d8           | 2.037 | 2.136 | -4.9  | 88    | 0.00     |
| 47   | Acenaphthylene              | 1.975 | 2.086 | -5.6  | 89    | 0.00     |
| 48   | 3-Nitroaniline              | 0.292 | 0.293 | -0.3  | 87    | 0.00     |
| 49 C | Acenaphthene                | 1.367 | 1.413 | -3.4  | 89    | 0.00     |
| 50   | 2,4-Dinitrophenol           | 0.208 | 0.184 | 11.5  | 90    | 0.00     |
| 51 S | 4-Nitrophenol-d4            | 0.303 | 0.292 | 3.6   | 89    | 0.00     |
| 52   | 4-Nitrophenol               | 0.242 | 0.246 | -1.7  | 91    | 0.00     |
| 53   | Dibenzofuran                | 1.984 | 2.071 | -4.4  | 89    | 0.00     |
| 54   | 2,4-Dinitrotoluene          | 0.465 | 0.514 | -10.5 | 91    | 0.00     |
| 55   | 2,3,4,6-Tetrachlorophenol   | 0.443 | 0.473 | -6.8  | 90    | 0.00     |
| 56   | Diethylphthalate            | 1.693 | 1.778 | -5.0  | 89    | 0.00     |
| 57 S | Fluorene-d10                | 1.530 | 1.598 | -4.4  | 90    | 0.00     |
| 58   | Fluorene                    | 1.656 | 1.714 | -3.5  | 89    | 0.00     |
| 59   | 4-Chlorophenyl-phenylether  | 0.876 | 0.907 | -3.5  | 89    | 0.00     |
| 60   | 4-Nitroaniline              | 0.340 | 0.318 | 6.5   | 90    | 0.00     |
| 61 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 91    | 0.00     |
| 62 S | 4,6-Dinitro-2-methylphenol- | 0.135 | 0.129 | 4.4   | 92    | 0.00     |
| 63   | 4,6-Dinitro-2-methylphenol  | 0.137 | 0.134 | 2.2   | 93    | 0.00     |
| 64   | N-Nitrosodiphenylamine      | 0.578 | 0.602 | -4.2  | 89    | 0.00     |
| 65   | 4-Bromophenyl-phenylether   | 0.224 | 0.229 | -2.2  | 90    | 0.00     |
| 66   | Hexachlorobenzene           | 0.250 | 0.256 | -2.4  | 90    | 0.00     |
| 67   | Atrazine                    | 0.229 | 0.232 | -1.3  | 90    | 0.00     |
| 68 C | Pentachlorophenol           | 0.158 | 0.153 | 3.2   | 92    | 0.00     |
| 69   | Phenanthrene                | 1.100 | 1.143 | -3.9  | 91    | 0.00     |
| 70 S | Anthracene-d10              | 0.986 | 1.038 | -5.3  | 92    | 0.00     |
| 71   | Anthracene                  | 1.119 | 1.182 | -5.6  | 91    | 0.00     |
| 72   | Carbazole                   | 0.970 | 0.993 | -2.4  | 92    | 0.00     |
| 73   | Di-n-butylphthalate         | 1.162 | 1.228 | -5.7  | 90    | 0.00     |
| 74 C | Fluoranthene                | 1.342 | 1.448 | -7.9  | 93    | 0.00     |
| 75 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 76 S | Pyrene-d10                  | 0.902 | 0.969 | -7.4  | 93    | 0.00     |
| 77   | Pyrene                      | 1.109 | 1.200 | -8.2  | 93    | 0.00     |
| 78   | Butylbenzylphthalate        | 0.434 | 0.471 | -8.5  | 93    | 0.00     |
| 79   | 3,3'-Dichlorobenzidine      | 0.372 | 0.405 | -8.9  | 101   | 0.00     |
| 80   | Benzo(a)anthracene          | 1.146 | 1.258 | -9.8  | 98    | 0.00     |
| 81   | Bis(2-ethylhexyl)phthalate  | 0.646 | 0.694 | -7.4  | 94    | 0.00     |
| 82   | Chrysene                    | 1.089 | 1.183 | -8.6  | 97    | 0.00     |
| 83 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 104   | 0.00     |
| 84   | Di-n-octyl phthalate        | 1.168 | 1.233 | -5.6  | 98    | 0.00     |
| 85   | Benzo(b)fluoranthene        | 1.216 | 1.267 | -4.2  | 100   | 0.00     |
| 86   | Benzo(k)fluoranthene        | 1.214 | 1.307 | -7.7  | 105   | 0.00     |
| 87 S | Benzo(a)pyrene-d12          | 0.962 | 1.020 | -6.0  | 104   | 0.00     |

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Misc :  
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|      | Compound               | AvgRF | CCRF  | %Dev | Area% | Dev(min) |
|------|------------------------|-------|-------|------|-------|----------|
| 88 C | Benzo(a)pyrene         | 1.168 | 1.240 | -6.2 | 104   | 0.00     |
| 89   | Indeno(1,2,3-cd)pyrene | 1.250 | 1.294 | -3.5 | 109   | 0.00     |
| 90   | Dibenzo(a,h)anthracene | 1.027 | 1.078 | -5.0 | 110   | 0.01     |
| 91   | Benzo(a,h,i)perylene   | 1.034 | 1.063 | -2.8 | 109   | 0.01     |

(#) = Out of Range

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