

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM070225\  
 Data File : BM050344.D  
 Acq On : 02 Jul 2025 15:14  
 Operator : RC/JU  
 Sample : SSTDCCC020EC  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 LabSampleId :  
 SSTDCCC020EC

Quant Time: Jul 02 15:50:39 2025  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM070125.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Tue Jul 01 15:14:38 2025  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 1 I  | 1,4-Dichlorobenzene-d4      | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 2    | 1,4-Dioxane                 | 0.503 | 0.545 | -8.3  | 103   | 0.00     |
| 3 S  | 1,4-Dioxane-d8              | 0.485 | 0.521 | -7.4  | 100   | 0.00     |
| 4 S  | Pyridine-d5                 | 1.312 | 1.434 | -9.3  | 102   | 0.00     |
| 5    | Pyridine                    | 1.369 | 1.507 | -10.1 | 101   | 0.00     |
| 6    | Benzaldehyde                | 0.800 | 0.986 | -23.2 | 103   | 0.00     |
| 7 S  | Phenol-d5                   | 1.450 | 1.544 | -6.5  | 100   | 0.00     |
| 8    | Phenol                      | 1.519 | 1.667 | -9.7  | 102   | 0.00     |
| 9 S  | Bis-(2-Chloroethyl)ether-d8 | 0.919 | 1.036 | -12.7 | 104   | 0.00     |
| 10   | Bis(2-Chloroethyl)ether     | 1.193 | 1.327 | -11.2 | 103   | 0.00     |
| 11 S | 2-Chlorophenol-d4           | 1.181 | 1.230 | -4.1  | 98    | 0.00     |
| 12   | 2-Chlorophenol              | 1.239 | 1.292 | -4.3  | 99    | 0.00     |
| 13   | 2-Methylphenol              | 1.096 | 1.157 | -5.6  | 100   | 0.00     |
| 14   | 2,2'-oxybis(1-Chloropropane | 1.864 | 2.124 | -13.9 | 104   | 0.00     |
| 15 S | 4-Methylphenol-d8           | 1.105 | 1.158 | -4.8  | 99    | 0.00     |
| 16   | Acetophenone                | 1.771 | 1.966 | -11.0 | 103   | 0.00     |
| 17 P | N-Nitroso-di-n-propylamine  | 0.919 | 0.962 | -4.7  | 100   | 0.00     |
| 18   | 4-Methylphenol              | 1.191 | 1.259 | -5.7  | 99    | 0.00     |
| 19   | Hexachloroethane            | 0.511 | 0.540 | -5.7  | 101   | 0.00     |
| 20 I | Naphthalene-d8              | 1.000 | 1.000 | 0.0   | 100   | 0.00     |
| 21 S | Nitrobenzene-d5             | 0.138 | 0.138 | 0.0   | 97    | 0.00     |
| 22   | Nitrobenzene                | 0.357 | 0.376 | -5.3  | 102   | 0.00     |
| 23   | Isophorone                  | 0.653 | 0.650 | 0.5   | 98    | 0.00     |
| 24 S | 2-Nitrophenol-d4            | 0.129 | 0.117 | 9.3   | 91    | 0.00     |
| 25 C | 2-Nitrophenol               | 0.153 | 0.148 | 3.3   | 95    | 0.00     |
| 26   | 2,4-Dimethylphenol          | 0.342 | 0.349 | -2.0  | 100   | 0.00     |
| 27   | Bis(2-Chloroethoxy)methane  | 0.421 | 0.429 | -1.9  | 100   | 0.00     |
| 28 S | 2,4-Dichlorophenol-d3       | 0.313 | 0.302 | 3.5   | 95    | 0.00     |
| 29 C | 2,4-Dichlorophenol          | 0.312 | 0.303 | 2.9   | 95    | 0.00     |
| 30   | Naphthalene                 | 1.029 | 1.040 | -1.1  | 99    | 0.00     |
| 31 S | 4-Chloroaniline-d4          | 0.416 | 0.429 | -3.1  | 99    | 0.00     |
| 32   | 4-Chloroaniline             | 0.421 | 0.434 | -3.1  | 99    | 0.00     |
| 33 C | Hexachlorobutadiene         | 0.219 | 0.206 | 5.9   | 94    | 0.00     |
| 34   | Caprolactam                 | 0.091 | 0.082 | 9.9   | 93    | 0.00     |
| 35 C | 4-Chloro-3-methylphenol     | 0.285 | 0.286 | -0.4  | 98    | 0.00     |
| 36   | 2-Methylnaphthalene         | 0.715 | 0.714 | 0.1   | 99    | 0.00     |
| 37   | 1-Methylnaphthalene         | 0.708 | 0.702 | 0.8   | 98    | 0.00     |
| 38 I | Acenaphthene-d10            | 1.000 | 1.000 | 0.0   | 97    | 0.00     |
| 39   | 1,2,4,5-Tetrachlorobenzene  | 0.651 | 0.622 | 4.5   | 94    | 0.00     |
| 40   | Hexachlorocyclopentadiene   | 0.339 | 0.303 | 10.6  | 91    | 0.00     |
| 41 C | 2,4,6-Trichlorophenol       | 0.373 | 0.359 | 3.8   | 93    | 0.00     |
| 42   | 2,4,5-Trichlorophenol       | 0.415 | 0.401 | 3.4   | 93    | 0.00     |
| 43   | 1,1'-Biphenyl               | 1.473 | 1.490 | -1.2  | 97    | 0.00     |
| 44   | 2-Chloronaphthalene         | 1.180 | 1.199 | -1.6  | 98    | 0.00     |
| 45   | 2-Nitroaniline              | 0.272 | 0.271 | 0.4   | 95    | 0.00     |
| 46 S | Dimethylphthalate-d6        | 1.362 | 1.336 | 1.9   | 95    | 0.00     |
| 47   | Dimethylphthalate           | 1.355 | 1.346 | 0.7   | 96    | 0.00     |

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 25% Max. Rel. Area : 150%

|      | Compound                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|------|-----------------------------|-------|-------|-------|-------|----------|
| 48   | 2,6-Dinitrotoluene          | 0.221 | 0.213 | 3.6   | 92    | 0.00     |
| 49 S | Acenaphthylene-d8           | 1.668 | 1.678 | -0.6  | 97    | 0.00     |
| 50   | Acenaphthylene              | 1.858 | 1.892 | -1.8  | 97    | 0.00     |
| 51   | 3-Nitroaniline              | 0.240 | 0.245 | -2.1  | 96    | 0.00     |
| 52 C | Acenaphthene                | 1.202 | 1.227 | -2.1  | 99    | 0.00     |
| 53   | 2,4-Dinitrophenol           | 0.099 | 0.069 | 30.3# | 89    | 0.00     |
| 54 S | 4-Nitrophenol-d4            | 0.245 | 0.232 | 5.3   | 94    | 0.00     |
| 55   | 4-Nitrophenol               | 0.192 | 0.198 | -3.1  | 97    | 0.00     |
| 56   | Dibenzofuran                | 1.731 | 1.743 | -0.7  | 97    | 0.00     |
| 57   | 2,4-Dinitrotoluene          | 0.330 | 0.329 | 0.3   | 94    | 0.00     |
| 58   | 2,3,4,6-Tetrachlorophenol   | 0.320 | 0.302 | 5.6   | 90    | 0.00     |
| 59   | Diethylphthalate            | 1.274 | 1.285 | -0.9  | 96    | 0.00     |
| 60 S | Fluorene-d10                | 1.237 | 1.233 | 0.3   | 96    | 0.00     |
| 61   | Fluorene                    | 1.438 | 1.460 | -1.5  | 98    | 0.00     |
| 62   | 4-Chlorophenyl-phenylether  | 0.735 | 0.729 | 0.8   | 97    | 0.00     |
| 63   | 4-Nitroaniline              | 0.237 | 0.258 | -8.9  | 96    | 0.00     |
| 64 I | Phenanthrene-d10            | 1.000 | 1.000 | 0.0   | 94    | 0.00     |
| 65 S | 4,6-Dinitro-2-methylphenol- | 0.086 | 0.067 | 22.1  | 90    | 0.00     |
| 66   | 4,6-Dinitro-2-methylphenol  | 0.097 | 0.077 | 20.6  | 87    | 0.00     |
| 67   | N-Nitrosodiphenylamine      | 0.615 | 0.616 | -0.2  | 95    | 0.00     |
| 68   | 4-Bromophenyl-phenylether   | 0.218 | 0.205 | 6.0   | 90    | 0.00     |
| 69   | Hexachlorobenzene           | 0.262 | 0.252 | 3.8   | 92    | 0.00     |
| 70   | Atrazine                    | 0.210 | 0.194 | 7.6   | 88    | 0.00     |
| 71 C | Pentachlorophenol           | 0.149 | 0.131 | 12.1  | 89    | 0.00     |
| 72   | Phenanthrene                | 1.138 | 1.139 | -0.1  | 94    | 0.00     |
| 73 S | Anthracene-d10              | 0.992 | 0.991 | 0.1   | 93    | 0.00     |
| 74   | Anthracene                  | 1.152 | 1.169 | -1.5  | 95    | 0.00     |
| 75   | 1,2,3,4-Tetrachlorobenzene  | 0.339 | 0.328 | 3.2   | 94    | 0.00     |
| 76   | Pentachlorobenzene          | 0.319 | 0.314 | 1.6   | 95    | 0.00     |
| 77   | Carbazole                   | 1.008 | 1.024 | -1.6  | 93    | 0.00     |
| 78   | Di-n-butylphthalate         | 1.018 | 1.009 | 0.9   | 90    | 0.00     |
| 79 I | Chrysene-d12                | 1.000 | 1.000 | 0.0   | 91    | 0.00     |
| 80   | Fluoranthene                | 1.229 | 1.184 | 3.7   | 90    | 0.00     |
| 81 S | Pyrene-d10                  | 1.068 | 1.039 | 2.7   | 90    | 0.00     |
| 82   | Pyrene                      | 1.333 | 1.320 | 1.0   | 92    | 0.00     |
| 83   | Butylbenzylphthalate        | 0.376 | 0.348 | 7.4   | 82    | 0.00     |
| 84   | 3,3'-Dichlorobenzidine      | 0.363 | 0.325 | 10.5  | 79    | 0.00     |
| 85   | Benzo(a)anthracene          | 1.305 | 1.300 | 0.4   | 90    | 0.00     |
| 86   | Bis(2-ethylhexyl)phthalate  | 0.588 | 0.564 | 4.1   | 84    | 0.00     |
| 87   | Chrysene                    | 1.234 | 1.241 | -0.6  | 91    | 0.00     |
| 88 I | Perylene-d12                | 1.000 | 1.000 | 0.0   | 89    | 0.00     |
| 89   | Di-n-octyl phthalate        | 1.001 | 0.818 | 18.3  | 81    | 0.00     |
| 90   | Benzo(b)fluoranthene        | 1.206 | 1.177 | 2.4   | 87    | 0.00     |
| 91   | Benzo(k)fluoranthene        | 1.229 | 1.256 | -2.2  | 91    | 0.00     |
| 92 S | Benzo(a)pyrene-d12          | 1.014 | 0.989 | 2.5   | 87    | 0.00     |
| 93 C | Benzo(a)pyrene              | 1.155 | 1.165 | -0.9  | 89    | 0.00     |

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|----|------------------------|-------|-------|------|-------|----------|
| 94 | Indeno(1,2,3-cd)pyrene | 1.428 | 1.418 | 0.7  | 90    | 0.00     |
| 95 | Dibenzo(a,h)anthracene | 1.186 | 1.182 | 0.3  | 90    | 0.00     |
| 96 | Benzo(g,h,i)perylene   | 1.148 | 1.154 | -0.5 | 91    | 0.00     |

(#) = Out of Range

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