

Method Path : Z:\HPCHEM1\BNA M\METHODS\  
 Method File : SOM02.2-EPA-BM042815.M  
 Title : SVOA CALIBRATION  
 Last Update : Tue Apr 28 06:36:34 2015  
 Response Via : Initial Calibration

## Calibration Files

5 =BM001125.D 10 =BM001126.D 20 =BM001127.D  
 40 =BM001128.D 80 =BM001129.D 160 =BM001130.D

|       | Compound              | 5              | 10    | 20    | 40    | 80    | 160   | Avg   | %RSD  |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|
| 1) I  | 1,4-Dichlorobenzene-d | -----ISTD----- |       |       |       |       |       |       |       |
| 2)    | 1,4-Dioxane           | 0.506          | 0.510 | 0.510 | 0.458 | 0.453 |       | 0.487 | 6.01  |
| 3) S  | 1,4-Dioxane-d8        | 0.426          | 0.449 | 0.483 | 0.430 | 0.420 |       | 0.441 | 5.79  |
| 4)    | Benzaldehyde          |                | 1.019 | 1.124 | 0.950 | 0.746 | 0.362 | 0.841 | 35.79 |
| 5) S  | Phenol-d5             |                | 1.506 | 1.663 | 1.622 | 1.672 | 1.792 | 1.651 | 6.22  |
| 6)    | Phenol                |                | 1.660 | 1.820 | 1.739 | 1.769 | 1.896 | 1.777 | 4.98  |
| 7) S  | Bis-(2-Chloroethy     |                | 1.024 | 1.107 | 1.022 | 1.017 | 1.067 | 1.047 | 3.70  |
| 8)    | Bis(2-Chloroethyl     |                | 1.330 | 1.443 | 1.342 | 1.347 | 1.417 | 1.376 | 3.68  |
| 9) S  | 2-Chlorophenol-d4     | 1.193          | 1.241 | 1.353 | 1.305 | 1.321 |       | 1.283 | 5.02  |
| 10)   | 2-Chlorophenol        | 1.293          | 1.305 | 1.435 | 1.347 | 1.384 |       | 1.353 | 4.31  |
| 11)   | 2-Methylphenol        |                | 1.201 | 1.347 | 1.297 | 1.323 | 1.428 | 1.319 | 6.25  |
| 12)   | 2,2'-oxybis(1-Chl     |                | 2.580 | 2.701 | 2.534 | 2.511 | 2.583 | 2.582 | 2.85  |
| 13) S | 4-Methylphenol-d8     |                | 1.228 | 1.365 | 1.322 | 1.356 | 1.431 | 1.340 | 5.55  |
| 14)   | Acetophenone          |                | 2.018 | 2.204 | 2.131 | 2.158 | 2.301 | 2.162 | 4.78  |
| 15) P | N-Nitroso-di-n-pr     | 0.920          | 0.993 | 1.129 | 1.098 | 1.126 |       | 1.053 | 8.80  |
| 16)   | 4-Methylphenol        |                | 1.331 | 1.478 | 1.440 | 1.454 | 1.546 | 1.450 | 5.38  |
| 17)   | Hexachloroethane      | 0.544          | 0.535 | 0.586 | 0.550 | 0.560 |       | 0.555 | 3.50  |
| 18) I | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |
| 19) S | Nitrobenzene-d5       | 0.122          | 0.129 | 0.139 | 0.140 | 0.146 |       | 0.135 | 7.01  |
| 20)   | Nitrobenzene          | 0.343          | 0.371 | 0.394 | 0.388 | 0.396 |       | 0.378 | 5.85  |
| 21)   | Isophorone            | 0.554          | 0.619 | 0.685 | 0.681 | 0.711 |       | 0.650 | 9.74  |
| 22) S | 2-Nitrophenol-d4      | 0.121          | 0.137 | 0.152 | 0.156 | 0.168 |       | 0.147 | 12.53 |
| 23) C | 2-Nitrophenol         | 0.131          | 0.162 | 0.172 | 0.174 | 0.180 |       | 0.164 | 11.78 |
| 24)   | 2,4-Dimethylpheno     | 0.333          | 0.358 | 0.375 | 0.368 | 0.379 |       | 0.363 | 5.15  |
| 25)   | Bis(2-Chloroethox     | 0.400          | 0.421 | 0.434 | 0.423 | 0.429 |       | 0.421 | 3.13  |
| 26) S | 2,4-Dichloropheno     | 0.253          | 0.279 | 0.302 | 0.300 | 0.315 |       | 0.290 | 8.43  |
| 27) C | 2,4-Dichloropheno     | 0.259          | 0.287 | 0.301 | 0.297 | 0.311 |       | 0.291 | 6.84  |
| 28)   | Naphthalene           | 1.044          | 1.056 | 1.089 | 1.033 | 1.052 |       | 1.055 | 1.98  |
| 29) S | 4-Chloroaniline-d     |                | 0.346 | 0.399 | 0.372 | 0.329 | 0.265 | 0.342 | 14.88 |
| 30)   | 4-Chloroaniline       |                | 0.359 | 0.413 | 0.381 | 0.340 | 0.279 | 0.354 | 14.09 |
| 31) C | Hexachlorobutadie     | 0.178          | 0.178 | 0.183 | 0.178 | 0.178 |       | 0.179 | 1.24  |
| 32)   | Caprolactam           |                | 0.072 | 0.084 | 0.091 | 0.102 | 0.107 | 0.091 | 15.18 |
| 33) C | 4-Chloro-3-methyl     | 0.276          | 0.307 | 0.329 | 0.325 | 0.345 |       | 0.316 | 8.38  |
| 34)   | 2-Methylnaphthale     | 0.709          | 0.727 | 0.752 | 0.721 | 0.743 |       | 0.730 | 2.35  |
| 35) I | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |
| 36)   | 1,2,4,5-Tetrachlo     | 0.607          | 0.618 | 0.642 | 0.613 | 0.618 |       | 0.620 | 2.17  |
| 37)   | Hexachlorocyclope     |                | 0.290 | 0.336 | 0.343 | 0.359 | 0.410 | 0.348 | 12.43 |
| 38) C | 2,4,6-Trichloroph     | 0.318          | 0.352 | 0.387 | 0.391 | 0.406 |       | 0.371 | 9.64  |
| 39)   | 2,4,5-Trichloroph     | 0.358          | 0.402 | 0.438 | 0.435 | 0.445 |       | 0.416 | 8.70  |
| 40)   | 1,1'-Biphenyl         | 1.684          | 1.671 | 1.740 | 1.678 | 1.667 |       | 1.688 | 1.78  |
| 41)   | 2-Chloronaphthale     | 1.276          | 1.303 | 1.336 | 1.291 | 1.278 |       | 1.297 | 1.89  |
| 42)   | 2-Nitroaniline        | 0.276          | 0.329 | 0.397 | 0.422 | 0.455 |       | 0.376 | 19.27 |
| 43) S | Dimethylphthalate     | 1.264          | 1.367 | 1.450 | 1.391 | 1.443 |       | 1.383 | 5.43  |
| 44)   | Dimethylphthalate     | 1.527          | 1.548 | 1.620 | 1.555 | 1.604 |       | 1.571 | 2.49  |
| 45)   | 2,6-Dinitrotoluen     | 0.215          | 0.257 | 0.302 | 0.316 | 0.339 |       | 0.286 | 17.36 |
| 46) S | Acenaphthylene-d8     | 1.786          | 1.924 | 2.045 | 1.998 | 2.023 |       | 1.955 | 5.38  |
| 47)   | Acenaphthylene        | 1.970          | 2.105 | 2.204 | 2.137 | 2.148 |       | 2.113 | 4.13  |
| 48)   | 3-Nitroaniline        |                | 0.252 | 0.333 | 0.329 | 0.323 | 0.317 | 0.311 | 10.70 |
| 49) C | Acenaphthene          | 1.408          | 1.392 | 1.462 | 1.394 | 1.411 |       | 1.413 | 2.01  |
| 50)   | 2,4-Dinitrophenol     |                | 0.097 | 0.149 | 0.173 | 0.204 | 0.236 | 0.172 | 30.76 |
| 51) S | 4-Nitrophenol-d4      |                | 0.235 | 0.276 | 0.287 | 0.321 | 0.335 | 0.291 | 13.59 |

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|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|--------|
| 52)   | 4-Nitrophenol     |                | 0.212 | 0.254 | 0.259 | 0.273 | 0.282 | 0.256 | 10.51  |
| 53)   | Dibenzofuran      | 1.931          | 1.945 | 2.002 | 1.925 | 1.945 |       | 1.950 | 1.57   |
| 54)   | 2,4-Dinitrotoluen | 0.342          | 0.402 | 0.463 | 0.468 | 0.505 |       | 0.436 | 14.71  |
| 55)   | 2,3,4,6-Tetrachlo | 0.255          | 0.305 | 0.350 | 0.351 | 0.376 |       | 0.327 | 14.58  |
| 56)   | Diethylphthalate  | 1.454          | 1.576 | 1.664 | 1.612 | 1.675 |       | 1.596 | 5.58   |
| 57) S | Fluorene-d10      | 1.356          | 1.370 | 1.393 | 1.333 | 1.381 |       | 1.367 | 1.70   |
| 58)   | Fluorene          | 1.592          | 1.604 | 1.665 | 1.619 | 1.706 |       | 1.637 | 2.89   |
| 59)   | 4-Chlorophenyl-ph | 0.787          | 0.765 | 0.796 | 0.761 | 0.783 |       | 0.778 | 1.91   |
| 60)   | 4-Nitroaniline    |                | 0.295 | 0.358 | 0.343 | 0.393 | 0.394 | 0.357 | 11.50  |
| 61) I | Phenanthrene-d10  | -----ISTD----- |       |       |       |       |       |       |        |
| 62) S | 4,6-Dinitro-2-met |                | 0.092 | 0.113 | 0.118 | 0.129 | 0.148 | 0.120 | 17.18  |
| 63)   | 4,6-Dinitro-2-met |                | 0.104 | 0.122 | 0.127 | 0.140 | 0.156 | 0.130 | 15.09  |
| 64)   | N-Nitrosodiphenyl | 0.574          | 0.593 | 0.624 | 0.611 | 0.620 |       | 0.605 | 3.40   |
| 65)   | 4-Bromophenyl-phe | 0.178          | 0.192 | 0.203 | 0.195 | 0.202 |       | 0.194 | 5.30   |
| 66)   | Hexachlorobenzene | 0.215          | 0.216 | 0.230 | 0.217 | 0.225 |       | 0.221 | 3.03   |
| 67)   | Atrazine          |                | 0.190 | 0.211 | 0.207 | 0.218 | 0.232 | 0.212 | 7.13   |
| 68) C | Pentachlorophenol |                | 0.089 | 0.110 | 0.118 | 0.135 | 0.156 | 0.122 | 20.82# |
| 69)   | Phenanthrene      | 1.140          | 1.135 | 1.176 | 1.117 | 1.157 |       | 1.145 | 1.97   |
| 70) S | Anthracene-d10    | 0.874          | 0.921 | 0.955 | 0.924 | 0.960 |       | 0.927 | 3.69   |
| 71)   | Anthracene        | 1.115          | 1.168 | 1.204 | 1.148 | 1.196 |       | 1.166 | 3.13   |
| 72)   | Carbazole         |                | 1.028 | 1.086 | 1.049 | 1.108 | 1.151 | 1.084 | 4.46   |
| 73)   | Di-n-butylphthala | 0.917          | 1.079 | 1.227 | 1.216 | 1.331 |       | 1.154 | 13.88  |
| 74) C | Fluoranthene      |                | 1.271 | 1.336 | 1.269 | 1.352 | 1.358 | 1.317 | 3.34   |
| 75) I | Chrysene-d12      | -----ISTD----- |       |       |       |       |       |       |        |
| 76) S | Pyrene-d10        | 0.840          | 0.879 | 0.919 | 0.889 | 0.905 |       | 0.887 | 3.40   |
| 77)   | Pyrene            | 1.179          | 1.227 | 1.270 | 1.211 | 1.218 |       | 1.221 | 2.69   |
| 78)   | Butylbenzylphthal | 0.315          | 0.394 | 0.477 | 0.499 | 0.544 |       | 0.446 | 20.42  |
| 79)   | 3,3'-Dichlorobenz |                | 0.315 | 0.389 | 0.363 | 0.363 | 0.355 | 0.357 | 7.50   |
| 80)   | Benzo(a)anthracen | 1.155          | 1.179 | 1.231 | 1.185 | 1.198 |       | 1.190 | 2.34   |
| 81)   | Bis(2-ethylhexyl) | 0.535          | 0.626 | 0.745 | 0.762 | 0.835 |       | 0.701 | 17.02  |
| 82)   | Chrysene          | 1.123          | 1.133 | 1.155 | 1.125 | 1.116 |       | 1.130 | 1.34   |
| 83) I | Perylene-d12      | -----ISTD----- |       |       |       |       |       |       |        |
| 84)   | Di-n-octyl phthal |                | 1.128 | 1.323 | 1.358 | 1.467 | 1.470 | 1.349 | 10.35  |
| 85)   | Benzo(b)fluoranth | 1.185          | 1.224 | 1.288 | 1.225 | 1.291 |       | 1.243 | 3.70   |
| 86)   | Benzo(k)fluoranth | 1.076          | 1.135 | 1.209 | 1.181 | 1.201 |       | 1.160 | 4.77   |
| 87) S | Benzo(a)pyrene-d1 | 0.851          | 0.877 | 0.950 | 0.912 | 0.945 |       | 0.907 | 4.73   |
| 88) C | Benzo(a)pyrene    | 1.125          | 1.148 | 1.221 | 1.182 | 1.224 |       | 1.180 | 3.71   |
| 89)   | Indeno(1,2,3-cd)p | 1.276          | 1.289 | 1.374 | 1.351 | 1.387 |       | 1.335 | 3.78   |
| 90)   | Dibenzo(a,h)anthr | 1.051          | 1.077 | 1.159 | 1.131 | 1.165 |       | 1.117 | 4.53   |
| 91)   | Benzo(g,h,i)peryl | 1.090          | 1.087 | 1.151 | 1.120 | 1.135 |       | 1.117 | 2.50   |

(#) = Out of Range