

Method Path : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\  
 Method File : 8270-SIM-BN043021.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Fri Apr 30 17:46:00 2021  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BN014500.D 0.2 =BN014501.D 0.4 =BN014502.D 0.8 =BN014503.D 1.6 =BN014504.D 3.2 =BN014505.D 5 =BN014506.D

|           | Compound              | 0.1            | 0.2   | 0.4   | 0.8   | 1.6   | 3.2   | 5     | Avg   | %RSD  |
|-----------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|
| 1) I      | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |
| 2)        | 1,4-Dioxane           | 0.670          | 0.522 | 0.402 | 0.441 | 0.460 | 0.451 | 0.409 | 0.479 | 19.36 |
| 3)        | n-Nitrosodimet...     |                | 0.128 | 0.121 | 0.201 | 0.232 | 0.261 | 0.251 | 0.199 | 30.87 |
| 4) S      | 2-Fluorophenol        | 0.894          | 0.824 | 0.702 | 0.756 | 0.799 | 0.827 | 0.781 | 0.798 | 7.59  |
| 5) S      | Phenol-d6             | 0.933          | 0.890 | 0.796 | 0.878 | 0.966 | 1.038 | 0.996 | 0.928 | 8.75  |
| 6)        | bis(2-Chloroet...     | 0.987          | 0.946 | 0.865 | 0.960 | 1.004 | 0.983 | 0.899 | 0.949 | 5.32  |
| 7) I      | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |
| 8) S      | Nitrobenzene-d5       | 0.211          | 0.212 | 0.196 | 0.223 | 0.240 | 0.262 | 0.258 | 0.229 | 10.93 |
| 9)        | Naphthalene           | 1.075          | 1.064 | 0.966 | 1.053 | 1.081 | 1.091 | 1.022 | 1.050 | 4.15  |
| 10)       | Hexachlorobuta...     | 0.225          | 0.217 | 0.192 | 0.207 | 0.209 | 0.210 | 0.195 | 0.208 | 5.51  |
| 11) SURR2 | Methylnaphth...       | 0.854          | 0.817 | 0.745 | 0.821 | 0.855 | 0.872 | 0.825 | 0.827 | 5.04  |
| 12)       | 2-Methylnaphth...     | 0.700          | 0.685 | 0.622 | 0.686 | 0.711 | 0.718 | 0.674 | 0.685 | 4.66  |
| 13) I     | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 14) S     | 2,4,6-Tribromo...     | 0.128          | 0.106 | 0.097 | 0.116 | 0.139 | 0.161 |       | 0.125 | 18.57 |
| 15) S     | 2-Fluorobiphenyl      | 1.495          | 1.558 | 1.428 | 1.561 | 1.594 | 1.602 | 1.468 | 1.529 | 4.32  |
| 16)       | Acenaphthylene        | 1.376          | 1.372 | 1.271 | 1.435 | 1.599 | 1.769 | 1.722 | 1.506 | 12.72 |
| 17)       | Acenaphthene          | 1.124          | 1.118 | 1.016 | 1.107 | 1.167 | 1.205 | 1.143 | 1.126 | 5.22  |
| 18)       | Fluorene              | 1.397          | 1.361 | 1.241 | 1.375 | 1.437 | 1.480 | 1.405 | 1.385 | 5.40  |
| 19) I     | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |
| 20)       | 4,6-Dinitro-2-...     | 0.012          | 0.008 | 0.013 | 0.025 | 0.040 | 0.053 | 0.061 | 0.030 | 69.93 |
| 21)       | 4-Bromophenyl-...     | 0.215          | 0.212 | 0.191 | 0.217 | 0.237 | 0.244 | 0.233 | 0.222 | 8.18  |
| 22)       | Hexachlorobenzene     | 0.280          | 0.287 | 0.256 | 0.284 | 0.299 | 0.296 | 0.280 | 0.283 | 4.93  |
| 23)       | Atrazine              | 0.125          | 0.115 | 0.099 | 0.115 | 0.130 | 0.145 | 0.155 | 0.126 | 15.12 |
| 24)       | Pentachlorophenol     |                | 0.035 | 0.029 | 0.047 | 0.064 | 0.078 | 0.088 | 0.057 | 41.30 |
| 25)       | Phenanthrene          | 1.150          | 1.140 | 1.018 | 1.134 | 1.227 | 1.234 | 1.178 | 1.154 | 6.25  |
| 26)       | Anthracene            | 0.839          | 0.802 | 0.735 | 0.852 | 0.965 | 1.033 | 1.035 | 0.895 | 13.12 |
| 27) SURR  | Fluoranthene-d10      | 1.382          | 1.370 | 1.231 | 1.376 | 1.515 | 1.567 | 1.539 | 1.426 | 8.40  |
| 28)       | Fluoranthene          | 1.196          | 1.180 | 1.077 | 1.225 | 1.356 | 1.380 | 1.326 | 1.249 | 8.81  |
| 29) I     | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |
| 30)       | Pyrene                | 1.543          | 1.418 | 1.216 | 1.333 | 1.367 | 1.365 | 1.301 | 1.363 | 7.44  |
| 31) S     | Terphenyl-d14         | 0.997          | 0.941 | 0.866 | 0.913 | 0.941 | 0.928 | 0.878 | 0.923 | 4.75  |
| 32)       | Benzo(a)anthra...     | 1.096          | 1.078 | 0.976 | 1.118 | 1.198 | 1.215 | 1.232 | 1.130 | 8.08  |
| 33)       | Chrysene              | 1.324          | 1.344 | 1.160 | 1.285 | 1.336 | 1.373 | 1.247 | 1.296 | 5.61  |
| 34)       | Bis(2-ethylhex...     | 0.611          | 0.513 | 0.359 | 0.362 | 0.361 | 0.388 | 0.420 | 0.430 | 22.41 |
| 35) I     | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |

Method Path : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\

Method File : 8270-SIM-BN043021.M

|       |                   |       |       |       |       |       |       |       |       |       |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 36)   | Indeno(1,2,3-c... | 1.237 | 1.355 | 1.240 | 1.480 | 1.671 | 1.754 | 1.664 | 1.486 | 14.47 |
| 37)   | Benzo(b)fluora... | 1.341 | 1.427 | 1.298 | 1.433 | 1.552 | 1.664 | 1.543 | 1.465 | 8.76  |
| 38)   | Benzo(k)fluora... | 1.654 | 1.721 | 1.504 | 1.732 | 1.817 | 1.790 | 1.702 | 1.703 | 6.05  |
| 39) C | Benzo(a)pyrene    | 1.312 | 1.305 | 1.127 | 1.281 | 1.373 | 1.449 | 1.397 | 1.321 | 7.86  |
| 40)   | Dibenzo(a,h)an... | 1.003 | 1.100 | 1.029 | 1.233 | 1.393 | 1.472 | 1.395 | 1.232 | 15.58 |
| 41)   | Benzo(g,h,i)pe... | 1.387 | 1.414 | 1.314 | 1.471 | 1.564 | 1.612 | 1.508 | 1.467 | 7.07  |

-----  
(#) = Out of Range