

Method Path : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\  
 Method File : 8270E-BP110724.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Thu Nov 07 23:36:11 2024  
 Response Via : Initial Calibration

## Calibration Files

2.5 =BP022914.D 5 =BP022915.D 10 =BP022916.D 20 =BP022917.D 40 =BP022918.D 50 =BP022919.D 60 =BP022920.D 80 =BP022921.D

|       | Compound              | 2.5            | 5     | 10    | 20    | 40    | 50    | 60    | 80    | Avg   | %RSD  |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 1) I  | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 2)    | 1,4-Dioxane           | 0.499          | 0.417 | 0.417 | 0.422 | 0.393 | 0.404 | 0.393 | 0.421 |       | 8.62  |
| 3)    | Pyridine              | 1.147          | 1.097 | 1.199 | 1.164 | 1.138 | 1.179 | 1.163 | 1.155 |       | 2.82  |
| 4)    | n-Nitrosodimet...     | 0.549          | 0.566 | 0.567 | 0.558 | 0.550 | 0.562 | 0.547 | 0.557 |       | 1.49  |
| 5) S  | 2-Fluorophenol        | 1.048          | 1.021 | 1.062 | 1.067 | 1.032 | 1.079 | 1.068 | 1.054 |       | 2.01  |
| 6)    | Aniline               | 1.309          | 1.273 | 1.298 | 1.239 | 1.173 | 1.192 | 1.092 | 1.225 |       | 6.37  |
| 7) S  | Phenol-d6             | 1.454          | 1.427 | 1.504 | 1.473 | 1.444 | 1.507 | 1.479 | 1.470 |       | 2.04  |
| 8)    | 2-Chlorophenol        | 1.110          | 1.082 | 1.126 | 1.115 | 1.071 | 1.129 | 1.100 | 1.105 |       | 1.98  |
| 9)    | Benzaldehyde          | 0.962          | 0.978 | 0.987 | 0.845 | 0.767 | 0.754 | 0.653 | 0.849 |       | 15.40 |
| 10) C | Phenol                | 1.393          | 1.399 | 1.501 | 1.475 | 1.451 | 1.501 | 1.467 | 1.456 |       | 3.05  |
| 11)   | bis(2-Chloroet...     | 1.144          | 1.066 | 1.106 | 1.097 | 1.073 | 1.122 | 1.090 | 1.100 |       | 2.48  |
| 12)   | 1,3-Dichlorobe...     | 1.492          | 1.411 | 1.434 | 1.396 | 1.372 | 1.404 | 1.359 | 1.410 |       | 3.12  |
| 13) C | 1,4-Dichlorobe...     | 1.521          | 1.441 | 1.484 | 1.409 | 1.389 | 1.442 | 1.392 | 1.440 |       | 3.40  |
| 14)   | 1,2-Dichlorobe...     | 1.491          | 1.422 | 1.426 | 1.373 | 1.333 | 1.364 | 1.333 | 1.392 |       | 4.14  |
| 15)   | Benzyl Alcohol        | 1.269          | 0.961 | 1.087 | 1.108 | 1.091 | 1.159 | 1.117 | 1.113 |       | 8.28  |
| 16)   | 2,2'-oxybis(1-...     | 1.266          | 1.216 | 1.256 | 1.173 | 1.140 | 1.176 | 1.125 | 1.193 |       | 4.59  |
| 17)   | 2-Methylphenol        | 0.965          | 0.918 | 1.007 | 1.010 | 0.975 | 1.043 | 0.998 | 0.988 |       | 4.04  |
| 18)   | Hexachloroethane      | 0.577          | 0.535 | 0.543 | 0.524 | 0.507 | 0.528 | 0.513 | 0.533 |       | 4.36  |
| 19) P | n-Nitroso-di-n...     | 0.981          | 1.061 | 1.020 | 1.078 | 1.010 | 0.956 | 0.993 | 0.949 | 1.006 | 4.58  |
| 20)   | 3+4-Methylphenols     | 1.338          | 1.322 | 1.421 | 1.392 | 1.366 | 1.435 | 1.393 | 1.381 |       | 3.00  |
| 21) I | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |       |
| 22)   | Acetophenone          | 0.579          | 0.520 | 0.528 | 0.526 | 0.526 | 0.532 | 0.516 | 0.532 |       | 3.98  |
| 23) S | Nitrobenzene-d5       | 0.432          | 0.404 | 0.432 | 0.427 | 0.424 | 0.430 | 0.415 | 0.423 |       | 2.47  |
| 24)   | Nitrobenzene          | 0.429          | 0.427 | 0.431 | 0.435 | 0.428 | 0.438 | 0.423 | 0.430 |       | 1.15  |
| 25)   | Isophorone            | 0.729          | 0.705 | 0.715 | 0.717 | 0.714 | 0.726 | 0.709 | 0.717 |       | 1.22  |
| 26) C | 2-Nitrophenol         | 0.172          | 0.163 | 0.170 | 0.179 | 0.182 | 0.187 | 0.183 | 0.177 |       | 4.70  |
| 27)   | 2,4-Dimethylph...     | 0.206          | 0.192 | 0.200 | 0.200 | 0.202 | 0.209 | 0.203 | 0.202 |       | 2.65  |
| 28)   | bis(2-Chloroet...     | 0.429          | 0.399 | 0.407 | 0.406 | 0.400 | 0.410 | 0.398 | 0.407 |       | 2.65  |
| 29) C | 2,4-Dichloroph...     | 0.333          | 0.328 | 0.343 | 0.358 | 0.355 | 0.365 | 0.362 | 0.349 |       | 4.23  |
| 30)   | 1,2,4-Trichlor...     | 0.481          | 0.440 | 0.442 | 0.439 | 0.436 | 0.449 | 0.436 | 0.446 |       | 3.56  |
| 31)   | Naphthalene           | 1.030          | 0.990 | 0.995 | 1.000 | 1.003 | 1.012 | 0.985 | 1.002 |       | 1.50  |
| 32)   | Benzoic acid          | 0.198          | 0.215 | 0.237 | 0.254 | 0.264 | 0.263 | 0.265 | 0.242 |       | 11.03 |
| 33)   | 4-Chloroaniline       | 0.357          | 0.343 | 0.366 | 0.365 | 0.359 | 0.366 | 0.340 | 0.357 |       | 3.04  |
| 34) C | Hexachlorobuta...     | 0.350          | 0.326 | 0.323 | 0.328 | 0.320 | 0.327 | 0.316 | 0.327 |       | 3.30  |
| 35)   | Caprolactam           | 0.106          | 0.095 | 0.105 | 0.106 | 0.106 | 0.109 | 0.105 | 0.104 |       | 4.28  |
| 36) C | 4-Chloro-3-met...     | 0.375          | 0.372 | 0.397 | 0.397 | 0.391 | 0.401 | 0.391 | 0.389 |       | 2.86  |
| 37)   | 2-Methylnaphth...     | 0.796          | 0.724 | 0.740 | 0.747 | 0.743 | 0.764 | 0.739 | 0.751 |       | 3.13  |
| 38)   | 1-Methylnaphth...     | 0.762          | 0.732 | 0.751 | 0.750 | 0.740 | 0.750 | 0.730 | 0.745 |       | 1.54  |

8270E-BP110724.M Fri Nov 08 19:01:38 2024

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|       |                   |                |       |       |       |       |       |       |       |      |  |
|-------|-------------------|----------------|-------|-------|-------|-------|-------|-------|-------|------|--|
| 86) I | Perylene-d12      | -----ISTD----- |       |       |       |       |       |       |       |      |  |
| 87)   | Indeno(1,2,3-c... | 1.405          | 1.350 | 1.352 | 1.325 | 1.314 | 1.364 | 1.334 | 1.349 | 2.21 |  |
| 88)   | Benzo(b)fluora... | 1.160          | 1.188 | 1.136 | 1.197 | 1.160 | 1.206 | 1.178 | 1.175 | 2.09 |  |
| 89)   | Benzo(k)fluora... | 1.129          | 1.118 | 1.129 | 1.077 | 1.092 | 1.141 | 1.081 | 1.110 | 2.34 |  |
| 90) C | Benzo(a)pyrene    | 1.003          | 0.979 | 1.009 | 1.003 | 1.000 | 1.020 | 1.022 | 1.005 | 1.41 |  |
| 91)   | Dibenzo(a,h)an... | 1.181          | 1.106 | 1.104 | 1.074 | 1.075 | 1.110 | 1.093 | 1.106 | 3.28 |  |
| 92)   | Benzo(g,h,i)pe... | 1.189          | 1.119 | 1.128 | 1.124 | 1.093 | 1.146 | 1.135 | 1.133 | 2.60 |  |

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(#) = Out of Range