

Method Path : Z:\HPCHEM1\BNA\_E\METHODS\  
 Method File : 8270-SIM-BE090115.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 Last Update : Wed Sep 02 11:40:50 2015  
 Response Via : Initial Calibration

## Calibration Files

0.1 =BE090669.D 0.2 =BE090670.D 0.5 =BE090671.D 0.8 =BE090672.D 1 =BE090673.D 2 =BE090674.D 5 =BE090675.D  
 10 =BE090676.D

|       | Compound              |                | 0.1   | 0.2   | 0.5   | 0.8   | 1     | 2     | 5     | 10    | Avg      | %RSD |
|-------|-----------------------|----------------|-------|-------|-------|-------|-------|-------|-------|-------|----------|------|
| 1) I  | 1,4-Dichlorobenzen... | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |
| 2)    | 1,4-Dioxane           | 1.042          | 1.133 | 0.707 | 0.848 | 0.792 | 0.675 | 0.682 | 0.630 | 0.814 | 22.65    |      |
| 3)    | n-Nitrosodimet...     | 1.279          | 1.507 | 0.876 | 0.975 | 0.978 | 0.862 | 0.917 | 0.871 | 1.033 | 22.71    |      |
| 4) S  | 2-Fluorophenol        | 1.153          | 1.416 | 0.948 | 1.013 | 0.976 | 0.865 | 0.959 | 0.982 | 1.039 | 16.60    |      |
| 5) S  | Phenol-d6             | 1.379          | 1.718 | 1.201 | 1.358 | 1.322 | 1.237 | 1.438 | 1.483 | 1.392 | 11.62    |      |
| 6) I  | Naphthalene-d8        | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |
| 7) S  | Nitrobenzene-d5       | 0.288          | 0.399 | 0.280 | 0.342 | 0.330 | 0.305 | 0.348 | 0.351 | 0.330 | 11.82    |      |
| 8)    | Nitrobenzene          | 0.284          | 0.449 | 0.299 | 0.365 | 0.351 | 0.352 | 0.409 | 0.412 | 0.365 | 15.54    |      |
| 9)    | Naphthalene           | 1.144          | 1.445 | 0.979 | 1.131 | 1.086 | 0.977 | 1.009 | 0.954 | 1.091 | 14.74    |      |
| 10)   | Hexachlorobuta...     | 0.173          | 0.224 | 0.155 | 0.182 | 0.176 | 0.154 | 0.158 | 0.143 | 0.171 | 14.81    |      |
| 11)   | 2-Methylnaphth...     | 0.557          | 0.732 | 0.506 | 0.572 | 0.572 | 0.541 | 0.621 | 0.617 | 0.590 | 11.65    |      |
| 12) I | Acenaphthene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |
| 13) S | 2,4,6-Tribromo...     | 0.176          | 0.204 | 0.117 | 0.130 | 0.121 | 0.107 | 0.137 | 0.150 | 0.143 | 22.98    |      |
| 14) S | 2-Fluorobiphenyl      | 1.366          | 1.786 | 1.203 | 1.497 | 1.364 | 1.246 | 1.326 | 1.289 | 1.385 | 13.36    |      |
| 15)   | Acenaphthylene        | 0.782          | 1.008 | 0.670 | 0.820 | 0.767 | 0.718 | 0.816 | 0.813 | 0.799 | E1 12.43 |      |
| 16)   | Acenaphthene          | 1.297          | 1.734 | 1.172 | 1.430 | 1.350 | 1.208 | 1.268 | 1.222 | 1.335 | 13.57    |      |
| 17)   | Fluorene              | 1.761          | 2.110 | 1.410 | 1.654 | 1.650 | 1.520 | 1.662 | 1.558 | 1.666 | 12.53    |      |
| 18) I | Phenanthrene-d10      | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |
| 19)   | 4-Bromophenyl-...     | 0.167          | 0.220 | 0.150 | 0.175 | 0.164 | 0.158 | 0.170 | 0.161 | 0.170 | 12.51    |      |
| 20)   | Hexachlorobenzene     | 0.923          | 1.239 | 0.844 | 1.015 | 0.915 | 0.845 | 0.846 | 0.794 | 0.928 | 15.41    |      |
| 21)   | Pentachlorophenol     |                |       | 0.048 | 0.053 | 0.047 | 0.044 | 0.058 | 0.073 | 0.054 | 19.58    |      |
| 22)   | Phenanthrene          | 1.219          | 1.615 | 1.102 | 1.301 | 1.234 | 1.144 | 1.174 | 1.111 | 1.237 | 13.44    |      |
| 23)   | Anthracene            | 0.975          | 1.248 | 0.855 | 1.027 | 0.995 | 0.993 | 1.111 | 1.075 | 1.035 | 11.10    |      |
| 24)   | Fluoranthene          | 1.406          | 1.714 | 1.092 | 1.247 | 1.209 | 1.162 | 1.320 | 1.286 | 1.304 | 14.66    |      |
| 25) I | Chrysene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |
| 26)   | Pyrene                | 1.100          | 1.246 | 0.877 | 0.980 | 0.981 | 0.934 | 1.084 | 1.077 | 1.035 | 11.20    |      |
| 27) S | Terphenyl-d14         | 0.582          | 0.675 | 0.474 | 0.523 | 0.528 | 0.503 | 0.573 | 0.567 | 0.553 | 11.12    |      |
| 28)   | Benzo(a)anthra...     | 1.241          | 1.547 | 0.978 | 1.126 | 1.094 | 1.035 | 1.141 | 1.086 | 1.156 | 15.23    |      |
| 29)   | Chrysene              | 1.294          | 1.745 | 1.124 | 1.341 | 1.264 | 1.135 | 1.188 | 1.138 | 1.279 | 16.05    |      |
| 30)   | Bis(2-ethylhex...     | 0.366          | 0.491 | 0.233 | 0.293 | 0.254 | 0.254 | 0.395 | 0.477 | 0.345 | 29.70    |      |
| 31)   | Indeno(1,2,3-c...     | 1.120          | 1.398 | 0.945 | 1.335 | 1.090 | 1.058 | 1.236 | 1.268 | 1.181 | 12.96    |      |
| 32) I | Perylene-d12          | -----ISTD----- |       |       |       |       |       |       |       |       |          |      |

Method Path : Z:\HPCHEM1\BNA\_E\METHODS\

Method File : 8270-SIM-BE090115.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|       |                   |       |       |       |       |       |       |       |       |       |       |
|-------|-------------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
| 33)   | Benzo(b)fluora... | 1.009 | 1.253 | 0.876 | 1.078 | 1.056 | 1.033 | 1.204 | 1.163 | 1.084 | 11.11 |
| 34)   | Benzo(k)fluora... | 1.134 | 1.530 | 1.061 | 1.231 | 1.244 | 1.245 | 1.342 | 1.313 | 1.263 | 11.18 |
| 35) C | Benzo(a)pyrene    | 0.924 | 1.156 | 0.770 | 0.947 | 0.906 | 0.901 | 1.069 | 1.085 | 0.970 | 12.87 |
| 36)   | Dibenzo(a,h)an... | 0.810 | 1.139 | 0.791 | 1.065 | 0.980 | 0.981 | 1.191 | 1.204 | 1.020 | 15.66 |
| 37)   | Benzo(g,h,i)pe... | 1.024 | 1.342 | 0.905 | 1.198 | 1.056 | 1.019 | 1.190 | 1.192 | 1.116 | 12.45 |

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