

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

Method File : 8270-BF120518.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Last Update : Wed Dec 05 17:30:55 2018

Response Via : Initial Calibration

Calibration Files

2.5 =BF111105.D 10 =BF111106.D 25 =BF111107.D 40 =BF111108.D 50 =BF111109.D 60 =BF111110.D 80 =BF111111.D

Compound	2.5	10	25	40	50	60	80	Avg	%RSD
----------	-----	----	----	----	----	----	----	-----	------

1) I	1,4-Dichlorobenzen...	-----ISTD-----							
2)	1,4-Dioxane	0.777	0.777	0.746	0.731	0.706	0.713	0.710	0.737
3)	Pyridine	1.619	1.690	1.674	1.628	1.599	1.609	1.626	1.635
4)	n-Nitrosodimet...	0.791	0.819	0.801	0.809	0.792	0.806	0.817	0.805
5) S	2-Fluorophenol	1.498	1.425	1.328	1.252	1.183	1.149	1.137	1.282
6)	Aniline	2.434	2.352	2.242	2.200	2.087	2.062	2.064	2.206
7) S	Phenol-d6	1.916	1.844	1.687	1.618	1.521	1.499	1.473	1.651
8)	2-Chlorophenol	1.718	1.617	1.506	1.462	1.362	1.351	1.323	1.477
9)	Benzaldehyde		1.229	1.014	0.915	0.828	0.776	0.720	0.914
10) C	Phenol	2.037	1.958	1.832	1.745	1.675	1.680	1.675	1.800
11)	bis(2-Chloroet...	1.685	1.595	1.494	1.447	1.362	1.382	1.335	1.471
12)	1,3-Dichlorobe...	1.866	1.707	1.604	1.520	1.459	1.439	1.403	1.571
13) C	1,4-Dichlorobe...	1.811	1.763	1.611	1.547	1.478	1.452	1.405	1.581
14)	1,2-Dichlorobe...	1.756	1.648	1.509	1.451	1.344	1.326	1.261	1.471
15)	Benzyl Alcohol	1.436	1.425	1.335	1.299	1.243	1.224	1.195	1.308
16)	2,2'-oxvibis(1-...	3.221	3.122	2.906	2.852	2.688	2.655	2.590	2.862
17)	2-Methylphenol	1.314	1.292	1.219	1.197	1.158	1.161	1.148	1.213
18)	Hexachloroethane	0.642	0.624	0.596	0.600	0.566	0.564	0.560	0.593
19) P	n-Nitroso-di-n...	1.226	1.201	1.107	1.080	1.047	1.056	1.050	1.110
20)	3+4-Methylphenols	1.773	1.681	1.530	1.445	1.334	1.322	1.268	1.479
									12.98
21) I	Naphthalene-d8	-----ISTD-----							
22)	Acetophenone	0.590	0.533	0.494	0.467	0.440	0.441	0.418	0.483
23) S	Nitrobenzene-d5	0.320	0.357	0.354	0.345	0.335	0.331	0.323	0.338
24)	Nitrobenzene	0.360	0.378	0.383	0.371	0.359	0.350	0.342	0.363
25)	Isophorone	0.676	0.665	0.625	0.606	0.588	0.594	0.586	0.620
26) C	2-Nitrophenol	0.109	0.142	0.160	0.164	0.164	0.169	0.167	0.154
27)	2,4-Dimethylph...	0.324	0.311	0.293	0.283	0.273	0.272	0.268	0.289
28)	bis(2-Chloroet...	0.476	0.457	0.429	0.414	0.393	0.396	0.375	0.420
29) C	2,4-Dichloroph...	0.298	0.302	0.292	0.285	0.276	0.275	0.265	0.285
30)	1,2,4-Trichlor...	0.323	0.305	0.294	0.285	0.271	0.272	0.262	0.287
31)	Naphthalene		1.067	0.969	0.878	0.821	0.776	0.713	0.871
32)	Benzoic acid		0.177	0.196	0.207	0.174	0.192	0.197	0.191
33)	4-Chloroaniline	0.473	0.474	0.435	0.422	0.389	0.379	0.374	0.421
34) C	Hexachlorobuta...	0.169	0.167	0.162	0.156	0.149	0.147	0.142	0.156
35)	Caprolactam	0.111	0.108	0.106	0.102	0.100	0.102	0.102	0.105
36) C	4-Chloro-3-met...	0.336	0.332	0.317	0.314	0.298	0.297	0.289	0.312
37)	2-Methylnaphth...	0.730	0.673	0.624	0.594	0.547	0.535	0.501	0.601
									13.52

Response Factor Report Instrumen

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

Method File : 8270-BF120518.M

38)	1-Methylnaphth...	0.682	0.656	0.598	0.566	0.533	0.520	0.488	0.577	12.42
<hr/>										
39) I	Acenaphthene-d10	-----ISTD-----								
40)	1,2,4,5-Tetrac...	0.631	0.593	0.547	0.521	0.497	0.488	0.454	0.533	11.68
41) P	Hexachlorocycl...	0.275	0.284	0.293	0.291	0.284	0.278	0.262	0.281	3.72
42) S	2,4,6-Tribromo...	0.170	0.169	0.170	0.160	0.151	0.150	0.143	0.159	7.13
43) C	2,4,6-Trichlor...	0.415	0.404	0.401	0.394	0.384	0.368	0.360	0.389	5.08
44)	2,4,5-Trichlor...	0.395	0.428	0.417	0.401	0.381	0.382	0.359	0.395	5.91
45) S	2-Fluorobiphenyl	1.476	1.286	1.134	1.049	1.003		1.190		16.21
46)	1,1'-Biphenyl	1.921	1.758	1.610	1.490	1.408		1.637		12.60
47)	2-Chloronaphth...	1.498	1.434	1.329	1.252	1.181	1.143	1.069	1.272	12.29
48)	2-Nitroaniline	0.378	0.417	0.426	0.421	0.417	0.415	0.412		4.23
49)	Acenaphthylene	2.127	1.963	1.808	1.694	1.612	1.465	1.778		13.52
50)	Dimethylphthalate	1.639	1.576	1.451	1.376	1.299	1.240	1.164	1.392	12.52
51)	2,6-Dinitrotol...	0.215	0.303	0.312	0.315	0.310	0.307	0.303	0.295	12.08
52) C	Acenaphthene	1.420	1.299	1.225	1.143	1.082	1.058	0.996	1.175	12.69
53)	3-Nitroaniline	0.297	0.377	0.398	0.390	0.381	0.383	0.370	0.371	9.12
54) P	2,4-Dinitrophenol	0.053	0.071	0.082	0.087	0.091	0.095	0.080		19.70
55)	Dibenzofuran	1.950	1.784	1.641	1.536	1.457	1.341	1.618		13.77
56) P	4-Nitrophenol	0.216	0.284	0.298	0.300	0.292	0.287	0.279	0.279	10.34
57)	2,4-Dinitrotol...	0.343	0.384	0.394	0.393	0.392	0.380	0.381		5.08
58)	Fluorene	1.388	1.239	1.154	1.072	1.037	0.956	1.141		13.62
59)	2,3,4,6-Tetrac...	0.296	0.319	0.316	0.306	0.290	0.277	0.259	0.295	7.29
60)	Diethylphthalate	1.611	1.547	1.450	1.384	1.312	1.263	1.170	1.391	11.27
61)	4-Chlorophenyl...	0.661	0.604	0.574	0.537	0.517	0.481	0.562		11.49
62)	4-Nitroaniline	0.277	0.330	0.346	0.347	0.350	0.349	0.342	0.335	7.87
63)	Azobenzene	1.772	1.668	1.548	1.470	1.364	1.328	1.217	1.481	13.24
<hr/>										
64) I	Phenanthrene-d10	-----ISTD-----								
65)	4,6-Dinitro-2-...	0.061	0.077	0.084	0.088	0.091	0.094	0.083		14.58
66) c	n-Nitrosodiphe...	0.821	0.782	0.723	0.688	0.646	0.644	0.598	0.700	11.44
67)	4-Bromophenyl-	0.239	0.226	0.220	0.209	0.199	0.201	0.189	0.212	8.17
68)	Hexachlorobenzene	0.246	0.230	0.224	0.212	0.203	0.206	0.194	0.216	8.20
69)	Atrazine	0.248	0.229	0.204	0.180	0.148	0.109		0.186	27.76
70) C	Pentachlorophenol	0.129	0.152	0.157	0.149	0.145	0.138	0.127	0.142	8.14
71)	Phenanthrene	1.168	1.067	0.970	0.895	0.869	0.803	0.962		14.10
72)	Anthracene	1.181	1.080	0.981	0.905	0.872	0.800	0.970		14.56
73)	Carbazole	1.237	1.100	1.019	0.937	0.916	0.844	1.009		14.11
74)	Di-n-butylphth...	1.425	1.275	1.167	1.033	1.005		1.181		14.75
75) C	Fluoranthene	1.152	1.021	0.964	0.874	0.850	0.788	0.941		14.04
<hr/>										
76) I	Chrysene-d12	-----ISTD-----								
77)	Benzidine	0.927	0.895	0.704	0.554	0.522	0.398		0.667	31.96
78)	Pyrene	1.609	1.586	1.577	1.559	1.525	1.553	1.499	1.558	2.38
79) S	Terphenyl-d14	1.116	1.066	0.992	0.940	0.925	0.920	0.877	0.977	8.84
80)	Butylbenzylphth...	0.640	0.713	0.774	0.790	0.773	0.797	0.781	0.753	7.53
81)	Benzo(a)anthra...	1.265	1.261	1.209	1.180	1.110	1.105	1.073	1.172	6.61
82)	3,3'-Dichlorob...	0.464	0.482	0.479	0.460	0.430	0.404	0.377	0.442	9.05

Response Factor Report Instrumen

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

Method File : 8270-BF120518.M

83)	Chrysene	1.251	1.235	1.209	1.165	1.113	1.089	1.082	1.163	6.05
84)	Bis(2-ethylhex...	0.878	0.914	0.944	0.951	0.894	0.901	0.883	0.909	3.16
85) C	Di-n-octyl phth...	1.416	1.485	1.496	1.472	1.389	1.388	1.359	1.429	3.83
86)	Indeno(1,2,3-c...	0.920	0.816	0.874	0.914	0.964	1.004	1.045	0.934	8.32

87) I Perylene-d12 -----ISTD-----

88)	Benzo(b)fluora...	1.171	1.236	1.138	1.165	1.056	1.114	1.054	1.133	5.77
89)	Benzo(k)fluora...	1.241	1.192	1.181	1.095	1.053	0.932	0.976	1.096	10.58
90) C	Benzo(a)pyrene	1.091	1.089	1.067	1.052	1.005	0.992	0.981	1.040	4.47
91)	Dibenzo(a,h)an...	0.850	0.805	0.864	0.890	0.914	0.890	0.894	0.872	4.19
92)	Benzo(g,h,i)pe...	0.846	0.770	0.868	0.895	0.919	0.905	0.934	0.877	6.37

(#) = Out of Range