

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF073024.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Tue Jul 30 17:50:01 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BF138680.D 5 =BF138681.D 10 =BF138682.D 20 =BF138683.D 40 =BF138684.D 50 =BF138685.D 60 =BF138686.D 80 =BF138687.D

	Compound	2.5	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.597	0.555	0.581	0.576	0.557	0.563	0.544	0.567	3.20	
3)	Pyridine	1.373	1.365	1.453	1.381	1.380	1.358	1.308	1.374	3.13	
4)	n-Nitrosodimet...	0.813	0.817	0.829	0.814	0.804	0.849	0.803	0.818	1.94	
5) S	2-Fluorophenol	1.404	1.367	1.341	1.274	1.244	1.256	1.183	1.296	6.00	
6)	Aniline	1.656	1.626	1.704	1.532	1.487	1.480	1.375	1.551	7.47	
7) S	Phenol-d6	1.938	1.816	1.828	1.681	1.654	1.684	1.575	1.740	7.19	
8)	2-Chlorophenol	1.471	1.418	1.443	1.325	1.308	1.325	1.252	1.363	5.94	
9)	Benzaldehyde	1.230	1.119	1.099	0.904	0.862			1.043	14.84	
10) C	Phenol	2.012	1.946	1.913	1.775	1.739	1.771	1.664	1.832	6.89	
11)	bis(2-Chloroet...	1.494	1.476	1.434	1.363	1.348	1.416	1.335	1.409	4.46	
12)	1,3-Dichlorobe...	1.687	1.609	1.608	1.475	1.445	1.468	1.389	1.526	7.12	
13) C	1,4-Dichlorobe...	1.689	1.615	1.620	1.496	1.482	1.470	1.407	1.540	6.61	
14)	1,2-Dichlorobe...	1.621	1.511	1.532	1.408	1.358	1.361	1.284	1.439	8.25	
15)	Benzyl Alcohol	1.326	1.287	1.334	1.218	1.213	1.246	1.151	1.254	5.27	
16)	2,2'-oxybis(1-...	2.695	2.576	2.592	2.350	2.311	2.317	2.138	2.426	8.17	
17)	2-Methylphenol	1.192	1.154	1.183	1.103	1.072	1.130	1.045	1.126	4.91	
18)	Hexachloroethane	0.619	0.614	0.598	0.568	0.554	0.564	0.540	0.580	5.31	
19) P	n-Nitroso-di-n...	1.109	1.130	1.087	1.123	0.992	0.970	1.025	0.970	1.051	6.56
20)	3+4-Methylphenols	1.661	1.554	1.558	1.367	1.328	1.365	1.277	1.444	10.02	
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.538	0.510	0.508	0.482	0.457	0.475	0.458	0.490	6.14	
23) S	Nitrobenzene-d5	0.426	0.416	0.424	0.407	0.393	0.405	0.392	0.409	3.33	
24)	Nitrobenzene	0.431	0.423	0.433	0.417	0.400	0.409	0.401	0.416	3.31	
25)	Isophorone	0.735	0.710	0.735	0.681	0.655	0.697	0.677	0.699	4.31	
26) C	2-Nitrophenol	0.169	0.176	0.188	0.184	0.175	0.184	0.179	0.179	3.75	
27)	2,4-Dimethylph...	0.219	0.217	0.221	0.214	0.207	0.215	0.207	0.214	2.55	
28)	bis(2-Chloroet...	0.448	0.433	0.449	0.418	0.402	0.416	0.410	0.425	4.31	
29) C	2,4-Dichloroph...	0.279	0.276	0.293	0.276	0.264	0.276	0.264	0.275	3.55	
30)	1,2,4-Trichlor...	0.344	0.324	0.329	0.316	0.304	0.308	0.299	0.318	4.96	
31)	Naphthalene	1.124	1.102	1.110	1.044	1.000	1.014	0.976	1.053	5.63	
32)	Benzoic acid		0.136	0.158	0.166	0.166	0.190	0.194	0.168	12.72	
33)	4-Chloroaniline	0.354	0.363	0.369	0.346	0.337	0.358	0.348	0.353	3.09	
34) C	Hexachlorobuta...	0.205	0.197	0.200	0.190	0.187	0.187	0.181	0.192	4.39	
35)	Caprolactam	0.078	0.084	0.088	0.078	0.076	0.087	0.085	0.082	5.59	
36) C	4-Chloro-3-met...	0.326	0.319	0.343	0.307	0.290	0.314	0.303	0.315	5.37	
37)	2-Methylnaphth...	0.723	0.699	0.709	0.649	0.619	0.640	0.615	0.665	6.71	
38)	1-Methylnaphth...	0.713	0.685	0.692	0.636	0.606	0.631	0.597	0.652	6.95	

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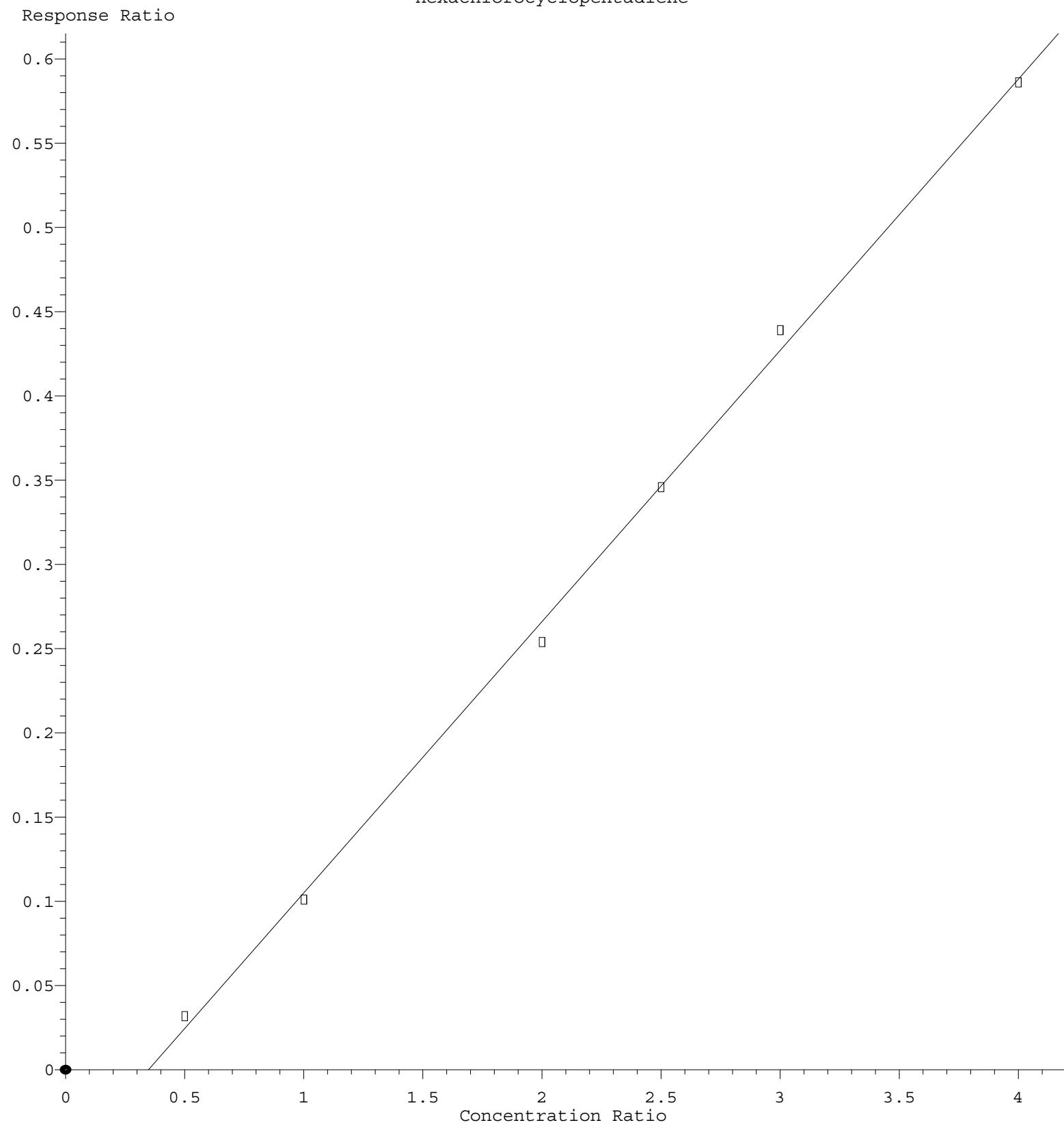
39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.606	0.560	0.581	0.550	0.542	0.534	0.515	0.556	5.48	
41) P	Hexachlorocycl...	0.064	0.101	0.127	0.138	0.146	0.147	0.120		27.08	
42) S	2,4,6-Tribromo...	0.168	0.164	0.176	0.159	0.156	0.166	0.158	0.164	4.19	
43) C	2,4,6-Trichlor...	0.338	0.338	0.353	0.336	0.330	0.341	0.334	0.339	2.10	
44)	2,4,5-Trichlor...	0.368	0.379	0.387	0.368	0.361	0.368	0.360	0.370	2.63	
45) S	2-Fluorobiphenyl	1.512	1.424	1.402	1.306	1.260	1.242	1.172	1.331	8.95	
46)	1,1'-Biphenyl	1.714	1.635	1.630	1.539	1.507	1.497	1.442	1.566	6.12	
47)	2-Chloronaphth...	1.263	1.206	1.213	1.151	1.119	1.114	1.089	1.165	5.46	
48)	2-Nitroaniline	0.394	0.403	0.414	0.389	0.380	0.395	0.390	0.395	2.74	
49)	Acenaphthylene	1.796	1.744	1.743	1.637	1.573	1.587	1.486	1.652	6.79	
50)	Dimethylphthalate	1.336	1.297	1.367	1.229	1.208	1.278	1.237	1.279	4.57	
51)	2,6-Dinitrotol...	0.286	0.288	0.304	0.288	0.277	0.300	0.277	0.289	3.57	
52) C	Acenaphthene	1.217	1.173	1.169	1.082	1.049	1.069	1.016	1.111	6.77	
53)	3-Nitroaniline	0.308	0.300	0.317	0.288	0.286	0.303	0.286	0.298	4.09	
54) P	2,4-Dinitrophenol	0.097	0.142	0.127	0.131	0.152	0.148	0.133		14.84	
55)	Dibenzofuran	1.729	1.680	1.666	1.515	1.484	1.488	1.412	1.568	7.75	
56) P	4-Nitrophenol	0.156	0.186	0.175	0.176	0.197	0.187	0.179		7.86	
57)	2,4-Dinitrotol...	0.377	0.376	0.402	0.362	0.342	0.371	0.348	0.368	5.41	
58)	Fluorene	1.374	1.326	1.351	1.206	1.166	1.189	1.127	1.249	7.97	
59)	2,3,4,6-Tetrac...	0.270	0.279	0.297	0.274	0.280	0.300	0.282	0.283	3.99	
60)	Diethylphthalate	1.268	1.259	1.321	1.155	1.113	1.216	1.156	1.213	6.16	
61)	4-Chlorophenyl...	0.662	0.664	0.662	0.589	0.580	0.591	0.551	0.614	7.67	
62)	4-Nitroaniline	0.284	0.292	0.313	0.272	0.265	0.292	0.267	0.284	6.04	
63)	Azobenzene	1.427	1.391	1.450	1.300	1.259	1.329	1.258	1.345	5.85	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.107	0.123	0.125	0.123	0.128	0.127	0.122		6.44	
66) c	n-Nitrosodiphe...	0.642	0.645	0.645	0.631	0.606	0.612	0.594	0.625	3.34	
67)	4-Bromophenyl-...	0.227	0.219	0.226	0.217	0.211	0.206	0.209	0.217	3.80	
68)	Hexachlorobenzene	0.231	0.231	0.234	0.220	0.214	0.219	0.216	0.224	3.73	
69)	Atrazine	0.171	0.172	0.176	0.162	0.153	0.153	0.141	0.161	7.84	
70) C	Pentachlorophenol	0.077	0.097	0.102	0.104	0.113	0.111	0.101		12.93	
71)	Phenanthrene	1.112	1.109	1.087	1.021	0.970	0.965	0.944	1.030	7.04	
72)	Anthracene	1.082	1.077	1.078	1.003	0.965	0.967	0.930	1.015	6.31	
73)	Carbazole	0.970	0.938	0.940	0.870	0.810	0.825	0.774	0.875	8.62	
74)	Di-n-butylphth...	0.991	0.990	1.040	0.991	0.944	0.992	0.941	0.984	3.41	
75) C	Fluoranthene	1.068	1.045	1.039	0.965	0.899	0.886	0.827	0.961	9.67	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.491	0.482	0.484	0.535	0.470	0.408	0.478		8.58	
78)	Pyrene	1.978	1.958	2.140	1.750	1.694	1.929	1.733	1.883	8.64	
79) S	Terphenyl-d14	1.264	1.265	1.362	1.106	1.072	1.210	1.083	1.195	9.25	
80)	Butylbenzylpht...	0.577	0.571	0.605	0.628	0.622	0.620	0.598	0.603	3.71	
81)	Benzo(a)anthra...	1.401	1.438	1.410	1.395	1.351	1.351	1.295	1.377	3.49	
82)	3,3'-Dichlorob...	0.388	0.364	0.377	0.360	0.346	0.319	0.314	0.352	8.00	
83)	Chrysene	1.342	1.228	1.260	1.222	1.203	1.229	1.213	1.243	3.82	
84)	Bis(2-ethylhex...	0.876	0.838	0.861	0.948	0.945	0.847	0.866	0.883	5.12	
85) c	Di-n-octyl pht...	1.637	1.524	1.614	1.744	1.742	1.570	1.605	1.634	5.06	

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86) I	Perylene-d12	-----ISTD-----									
87)	Indeno(1,2,3-c...	1.469	1.464	1.517	1.395	1.406	1.405	1.377	1.433	3.54	
88)	Benzo(b)fluora...	1.287	1.361	1.287	1.229	1.181	1.135	1.200	1.240	6.19	
89)	Benzo(k)fluora...	1.202	1.043	1.098	1.002	1.060	1.117	0.992	1.073	6.79	
90) C	Benzo(a)pyrene	1.084	1.047	1.071	1.023	1.026	1.035	1.014	1.043	2.50	
91)	Dibenzo(a,h)an...	1.226	1.231	1.253	1.134	1.140	1.147	1.106	1.177	4.94	
92)	Benzo(g,h,i)pe...	1.261	1.234	1.289	1.199	1.196	1.198	1.169	1.221	3.46	

(#) = Out of Range

Hexachlorocyclopentadiene



Response = $1.611\text{e-}001 \cdot \text{Amt} - 5.610\text{e-}002$
Coef of Det (r^2) = 0.998314 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF073024.M
Calibration Table Last Updated: Tue Jul 30 17:50:01 2024