

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF101524.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Tue Oct 15 16:41:54 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BF139805.D 5 =BF139806.D 10 =BF139807.D 20 =BF139808.D 40 =BF139809.D 50 =BF139810.D 60 =BF139811.D 80 =BF139812.D

	Compound	2.5	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.584	0.523	0.508	0.495	0.501	0.509	0.492	0.516	6.15	
3)	Pyridine	1.350	1.312	1.285	1.231	1.223	1.240	1.195	1.262	4.38	
4)	n-Nitrosodimet...	0.813	0.750	0.743	0.737	0.732	0.752	0.719	0.750	4.05	
5) S	2-Fluorophenol	1.316	1.245	1.230	1.150	1.141	1.134	1.082	1.185	6.82	
6)	Aniline	1.535	1.483	1.427	1.343	1.291	1.266	1.162	1.358	9.64	
7) S	Phenol-d6	1.806	1.674	1.636	1.527	1.492	1.489	1.411	1.576	8.60	
8)	2-Chlorophenol	1.409	1.301	1.309	1.227	1.202	1.190	1.122	1.251	7.60	
9)	Benzaldehyde	1.068	1.012	0.972	0.874	0.800	0.776	0.695	0.885	15.50	
10) C	Phenol	1.863	1.765	1.717	1.621	1.577	1.569	1.471	1.655	8.09	
11)	bis(2-Chloroet...	1.360	1.312	1.255	1.209	1.182	1.185	1.135	1.234	6.46	
12)	1,3-Dichlorobe...	1.611	1.509	1.452	1.369	1.344	1.356	1.264	1.415	8.27	
13) C	1,4-Dichlorobe...	1.602	1.485	1.475	1.378	1.360	1.357	1.280	1.419	7.58	
14)	1,2-Dichlorobe...	1.492	1.398	1.380	1.278	1.250	1.249	1.172	1.317	8.36	
15)	Benzyl Alcohol	1.230	1.221	1.186	1.122	1.091	1.082	1.009	1.135	7.16	
16)	2,2'-oxybis(1-...	2.452	2.297	2.257	2.106	2.059	2.043	1.923	2.163	8.37	
17)	2-Methylphenol	1.101	1.060	1.051	0.993	0.980	0.974	0.928	1.012	5.92	
18)	Hexachloroethane	0.553	0.543	0.536	0.504	0.496	0.496	0.472	0.514	5.78	
19) P	n-Nitroso-di-n...	0.988	1.018	0.980	0.950	0.894	0.855	0.857	0.800	0.918	8.41
20)	3+4-Methylphenols	1.494	1.400	1.329	1.228	1.185	1.180	1.096	1.273	11.01	
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.522	0.510	0.481	0.447	0.443	0.450	0.433	0.469	7.48	
23) S	Nitrobenzene-d5	0.417	0.402	0.400	0.374	0.377	0.381	0.369	0.389	4.58	
24)	Nitrobenzene	0.431	0.418	0.421	0.387	0.395	0.398	0.387	0.405	4.35	
25)	Isophorone	0.720	0.699	0.686	0.640	0.635	0.640	0.622	0.663	5.70	
26) C	2-Nitrophenol	0.170	0.178	0.179	0.175	0.177	0.179	0.176	0.176	1.70	
27)	2,4-Dimethylph...	0.233	0.222	0.223	0.206	0.206	0.209	0.201	0.214	5.50	
28)	bis(2-Chloroet...	0.449	0.426	0.423	0.388	0.389	0.397	0.378	0.407	6.31	
29) C	2,4-Dichloroph...	0.295	0.296	0.288	0.268	0.269	0.272	0.258	0.278	5.26	
30)	1,2,4-Trichlor...	0.348	0.328	0.328	0.303	0.306	0.309	0.293	0.316	5.98	
31)	Naphthalene	1.145	1.092	1.078	0.984	0.990	0.989	0.938	1.031	7.25	
32)	Benzoic acid	0.174	0.170	0.192	0.210	0.210	0.218	0.212	0.198	9.96	
33)	4-Chloroaniline	0.381	0.366	0.355	0.321	0.321	0.319	0.303	0.338	8.60	
34) C	Hexachlorobuta...	0.220	0.208	0.207	0.192	0.195	0.198	0.186	0.201	5.84	
35)	Caprolactam	0.086	0.086	0.081	0.078	0.075	0.074	0.070	0.078	7.82	
36) C	4-Chloro-3-met...	0.319	0.317	0.305	0.288	0.284	0.285	0.269	0.295	6.42	
37)	2-Methylnaphth...	0.724	0.693	0.680	0.618	0.615	0.615	0.579	0.646	8.15	
38)	1-Methylnaphth...	0.706	0.681	0.657	0.603	0.596	0.595	0.566	0.629	8.28	

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF101524.M

39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.634	0.597	0.602	0.551	0.559	0.573	0.550	0.581	5.42	
41) P	Hexachlorocycl...	0.053	0.087	0.123	0.130	0.143	0.143	0.113		31.67	
42) S	2,4,6-Tribromo...	0.187	0.183	0.184	0.166	0.156	0.158	0.147	0.169	9.42	
43) C	2,4,6-Trichlor...	0.389	0.366	0.379	0.363	0.360	0.372	0.355	0.369	3.19	
44)	2,4,5-Trichlor...	0.412	0.408	0.418	0.380	0.382	0.392	0.378	0.396	4.22	
45) S	2-Fluorobiphenyl	1.562	1.442	1.402	1.252	1.248	1.255	1.199	1.337	9.98	
46)	1,1'-Biphenyl	1.681	1.574	1.562	1.441	1.445	1.476	1.406	1.512	6.45	
47)	2-Chloronaphth...	1.255	1.202	1.199	1.104	1.103	1.143	1.081	1.155	5.59	
48)	2-Nitroaniline	0.377	0.396	0.393	0.372	0.368	0.375	0.359	0.377	3.51	
49)	Acenaphthylene	1.870	1.786	1.762	1.642	1.620	1.647	1.552	1.697	6.56	
50)	Dimethylphthalate	1.374	1.316	1.278	1.192	1.142	1.172	1.096	1.224	8.23	
51)	2,6-Dinitrotol...	0.288	0.294	0.295	0.278	0.269	0.274	0.258	0.279	4.90	
52) C	Acenaphthene	1.201	1.144	1.124	1.024	1.008	1.044	0.981	1.075	7.60	
53)	3-Nitroaniline	0.307	0.299	0.293	0.278	0.264	0.261	0.249	0.278	7.88	
54) P	2,4-Dinitrophenol	0.084	0.111	0.122	0.118	0.125	0.121	0.114		13.46	
55)	Dibenzofuran	1.807	1.743	1.699	1.541	1.528	1.530	1.434	1.612	8.53	
56) P	4-Nitrophenol	0.156	0.167	0.175	0.179	0.168	0.172	0.157	0.168	5.24	
57)	2,4-Dinitrotol...	0.372	0.376	0.369	0.345	0.324	0.326	0.302	0.345	8.36	
58)	Fluorene	1.468	1.398	1.333	1.183	1.154	1.169	1.075	1.254	11.61	
59)	2,3,4,6-Tetrac...	0.321	0.306	0.314	0.289	0.282	0.289	0.259	0.294	7.22	
60)	Diethylphthalate	1.388	1.321	1.264	1.149	1.091	1.105	1.020	1.191	11.35	
61)	4-Chlorophenyl...	0.714	0.686	0.651	0.589	0.583	0.582	0.535	0.620	10.46	
62)	4-Nitroaniline	0.284	0.280	0.274	0.251	0.236	0.238	0.219	0.255	9.85	
63)	Azobenzene	1.397	1.362	1.296	1.209	1.165	1.177	1.086	1.242	9.13	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.089	0.100	0.108	0.108	0.113	0.111	0.105		8.70	
66) c	n-Nitrosodiphe...	0.660	0.645	0.637	0.603	0.623	0.640	0.621	0.633	2.95	
67)	4-Bromophenyl-...	0.217	0.216	0.219	0.206	0.215	0.220	0.214	0.215	2.14	
68)	Hexachlorobenzene	0.258	0.245	0.243	0.232	0.238	0.240	0.231	0.241	3.69	
69)	Atrazine	0.178	0.164	0.140	0.093	0.094	0.087	0.071	0.118	35.35	
70) C	Pentachlorophenol	0.082	0.092	0.103	0.109	0.110	0.114	0.112	0.103	11.49	
71)	Phenanthrene	1.078	1.011	0.977	0.897	0.910	0.931	0.884	0.955	7.37	
72)	Anthracene	1.042	0.990	0.984	0.887	0.894	0.912	0.863	0.939	7.07	
73)	Carbazole	0.954	0.895	0.863	0.786	0.776	0.786	0.754	0.831	8.96	
74)	Di-n-butylphth...	1.058	0.973	0.927	0.869	0.834	0.844	0.812	0.902	9.80	
75) C	Fluoranthene	1.086	1.014	0.909	0.833	0.817	0.839	0.812	0.901	12.01	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.407	0.337	0.233	0.366	0.354	0.346	0.297	0.334	16.61	
78)	Pyrene	2.326	2.258	2.089	1.608	1.549		1.966		18.55	
79) S	Terphenyl-d14	1.633	1.512	1.351	1.029	0.977	0.942	0.823	1.181	26.62	
80)	Butylbenzylpht...	0.573	0.550	0.559	0.526	0.504	0.519	0.500	0.533	5.30	
81)	Benzo(a)anthra...	1.479	1.395	1.352	1.239	1.278	1.281	1.275	1.328	6.40	
82)	3,3'-Dichlorob...	0.418	0.415	0.443	0.451	0.454	0.465	0.452	0.443	4.33	
83)	Chrysene	1.288	1.219	1.241	1.190	1.190	1.219	1.130	1.211	4.04	
84)	Bis(2-ethylhex...	0.605	0.657	0.709	0.721	0.703	0.720	0.697	0.688	6.15	
85) c	Di-n-octyl pht...	1.022	1.115	1.277	1.361	1.380	1.416	1.383	1.279	11.93	

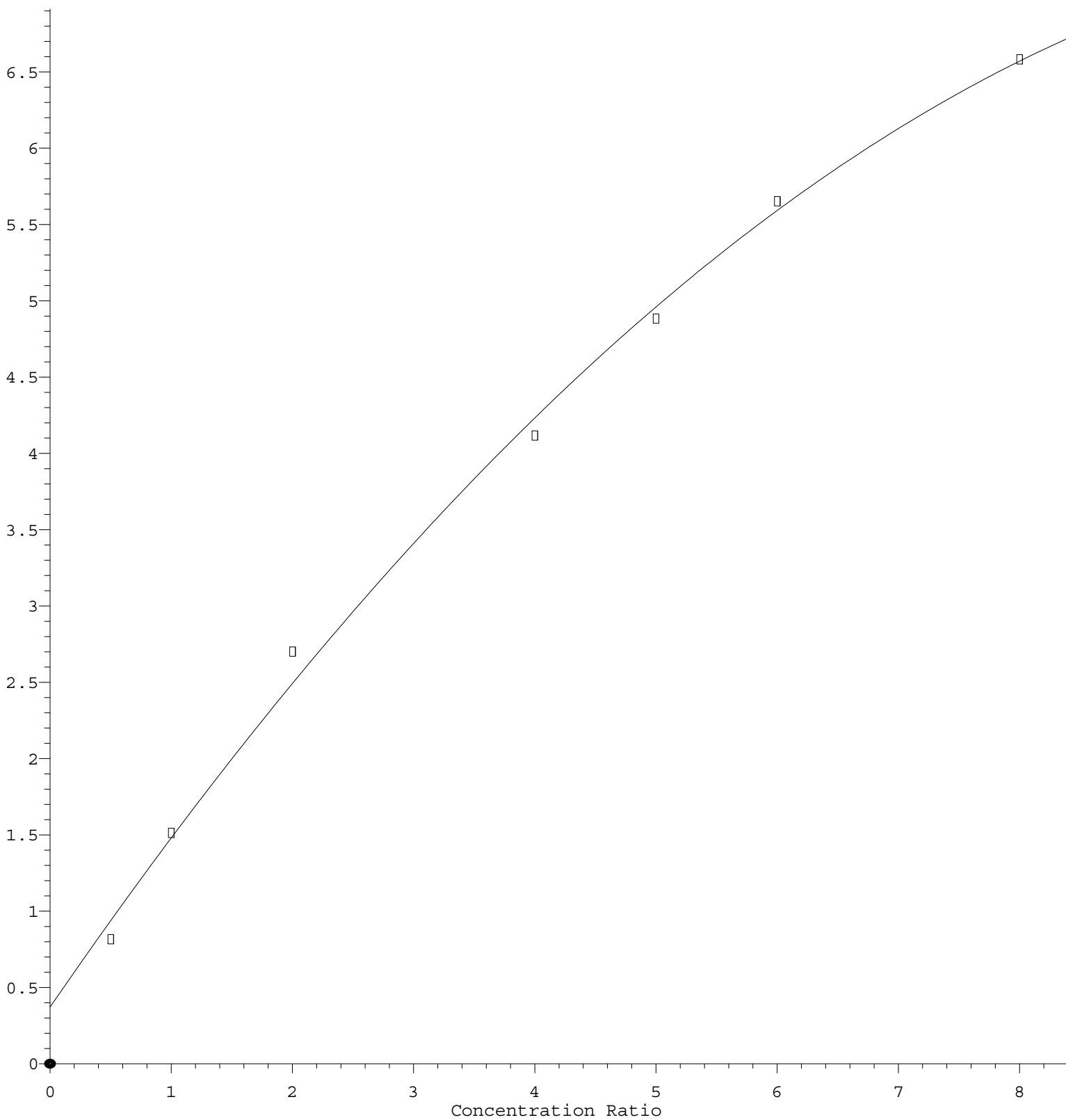
Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
Method File : 8270-BF101524.M

86) I	Perylene-d12	-----ISTD-----								
87)	Indeno(1,2,3-c...	1.466	1.428	1.426	1.246	1.273	1.299	1.237	1.339	7.26
88)	Benzo(b)fluora...	1.227	1.152	1.118	1.135	1.122	1.118	1.158	1.147	3.37
89)	Benzo(k)fluora...	1.106	1.046	1.083	1.006	0.977	0.999	0.909	1.018	6.55
90) C	Benzo(a)pyrene	1.052	1.011	1.032	0.985	0.980	0.991	0.963	1.002	3.13
91)	Dibenzo(a,h)an...	1.229	1.161	1.164	1.025	1.053	1.049	1.007	1.098	7.77
92)	Benzo(g,h,i)pe...	1.266	1.253	1.214	1.072	1.092	1.101	1.066	1.152	7.69

(#) = Out of Range

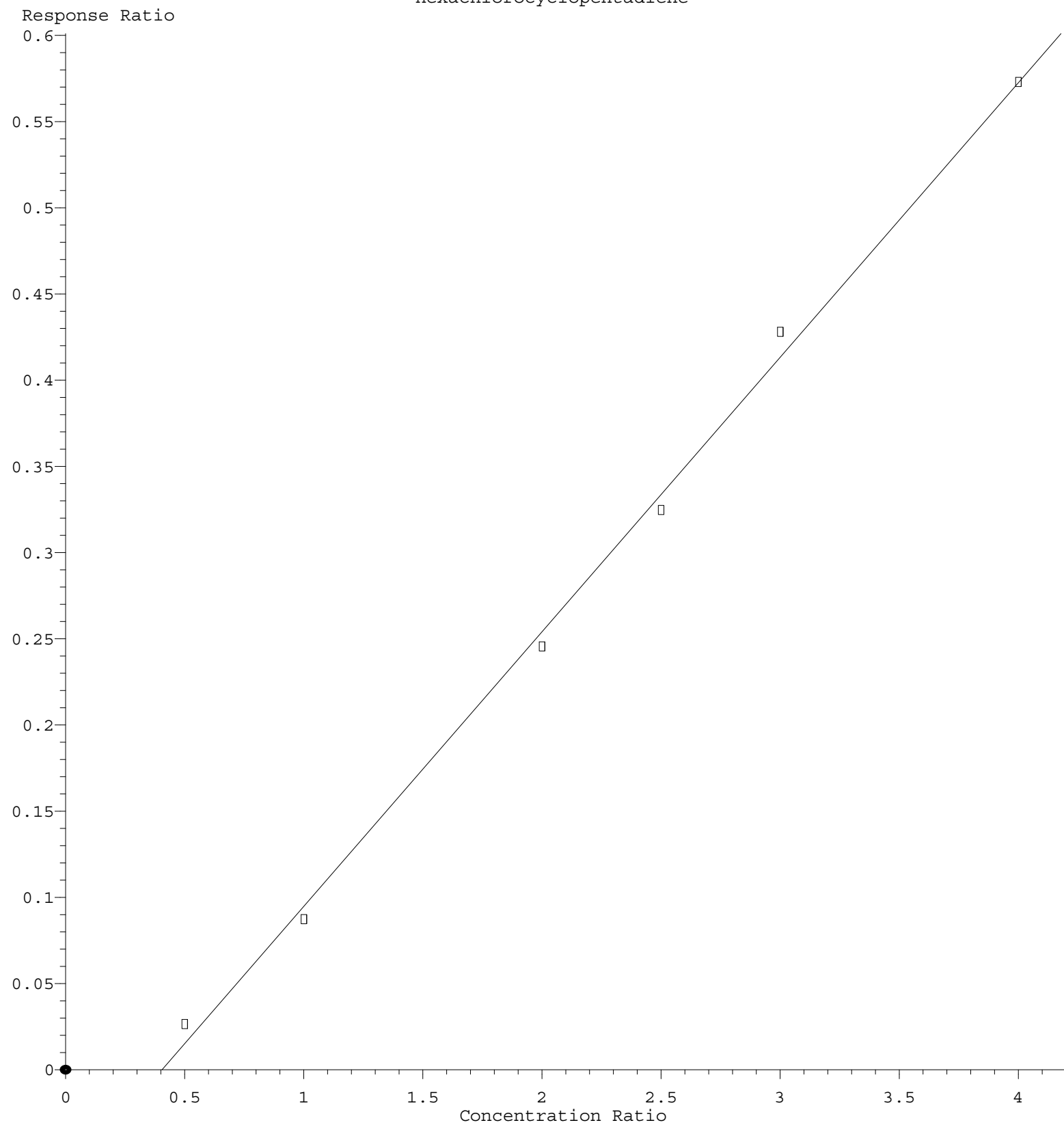
Terphenyl-d14

Response Ratio



$R = -4.761e-002 A^2 + 1.156e+000 A + 3.715e-001$
 Coef of Det (r^2) = 0.997031 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF101524.M
 Calibration Table Last Updated: Tue Oct 15 16:41:54 2024

Hexachlorocyclopentadiene



Response = $1.594 \times 10^{-1} \times \text{Amt} - 6.449 \times 10^{-2}$
Coef of Det (r^2) = 0.997388 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF101524.M
Calibration Table Last Updated: Tue Oct 15 16:41:54 2024