

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG060823.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Fri Jun 09 00:44:37 2023
 Response Via : Initial Calibration

Calibration Files

2.5 =BG057665.D 5 =BG057666.D 10 =BG057667.D 20 =BG057668.D 40 =BG057669.D 50 =BG057670.D 60 =BG057671.D 80 =BG057672.D

Compound		2.5	5	10	20	40	50	60	80	Avg	%RSD

1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.356	0.590	1.357	0.566	0.462	0.502	0.429	0.609	55.72	
3)	Pyridine	0.954	0.614	2.382	1.090	1.405	1.456	1.260	1.309	42.32	
4)	n-Nitrosodimet...	0.577	0.588	1.496	0.657	1.115	1.206	1.064	0.958	37.13	
5) S	2-Fluorophenol	0.973	0.563	1.169	0.543	1.147	1.245	1.097	0.962	30.30	
6)	Aniline	1.662	2.336	2.209	2.268	1.909	2.128	1.876	2.056	11.93	
7) S	Phenol-d6	1.263	1.863	1.770	1.318	1.603	1.729	1.534	1.583	14.36	
8)	2-Chlorophenol	1.009	1.330	1.280	1.264	1.161	1.277	1.102	1.203	9.66	
9)	Benzaldehyde	0.771	1.220	1.106	0.791	0.891	0.903	0.750	0.919	19.54	
10) C	Phenol	1.242	1.844	1.677	1.320	1.485	1.643	1.437	1.521	13.94	
11)	bis(2-Chloroet...	1.081	1.391	1.233	1.357	1.099	1.158	1.026	1.192	11.77	
12)	1,3-Dichlorobe...	1.149	1.360	1.339	1.321	1.235	1.356	1.186	1.278	6.80	
13) C	1,4-Dichlorobe...	1.423	1.466	1.401	1.348	1.269	1.354	1.198	1.351	6.83	
14)	1,2-Dichlorobe...	1.482	1.249	1.394	1.294	1.194	1.332	1.143	1.298	8.98	
15)	Benzyl Alcohol	1.172	1.200	1.378	1.295	1.292	1.422	1.259	1.288	6.97	
16)	2,2'-oxybis(1-...	2.836	2.613	2.693	2.651	2.198	2.318	2.027	2.477	12.01	
17)	2-Methylphenol	1.359	1.054	1.281	1.265	1.065	1.213	1.056	1.185	10.62	
18)	Hexachloroethane	0.475	0.532	0.576	0.503	0.499	0.526	0.466	0.511	7.35	
19) P	n-Nitroso-di-n...	1.092	1.185	1.171	1.259	1.166	1.007	1.090	0.985	1.119	8.33
20)	3+4-Methylphenols	1.642	1.708	1.767	1.735	1.471	1.627	1.425	1.625	8.06	
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.557	0.889	0.262	0.558	0.557	0.555	0.530	0.558	32.54	
23) S	Nitrobenzene-d5	0.435	0.692	0.224	0.449	0.466	0.461	0.432	0.451	30.13	
24)	Nitrobenzene	0.424	0.674	0.224	0.446	0.439	0.445	0.434	0.441	29.55	
25)	Isophorone	0.785	0.925	0.622	0.836	0.793	0.793	0.767	0.789	11.47	
26) C	2-Nitrophenol	0.154	0.177	0.180	0.185	0.185	0.189	0.187	0.179	6.64	
27)	2,4-Dimethylph...	0.297	0.325	0.343	0.335	0.337	0.333	0.326	0.328	4.60	
28)	bis(2-Chloroet...	0.465	0.292	0.494	0.473	0.396	0.399	0.386	0.415	16.68	
29) C	2,4-Dichloroph...	0.285	0.199	0.333	0.318	0.322	0.328	0.311	0.299	15.62	
30)	1,2,4-Trichlor...	0.364	0.362	0.369	0.352	0.364	0.361	0.352	0.361	1.78	
31)	Naphthalene	1.102	1.129	1.137	1.082	1.074	1.080	1.008	1.087	3.92	
32)	Benzoic acid		0.325	0.097	0.125	0.127	0.145	0.150	0.162	50.78	
33)	4-Chloroaniline	0.462	0.516	0.491	0.468	0.467	0.469	0.440	0.473	5.09	
34) C	Hexachlorobuta...	0.244	0.211	0.256	0.236	0.264	0.267	0.259	0.248	7.99	
35)	Caprolactam	0.073	0.074	0.105	0.104	0.095	0.102	0.096	0.093	14.67	
36) C	4-Chloro-3-met...	0.343	0.238	0.387	0.382	0.389	0.383	0.384	0.358	15.46	
37)	2-Methylnaphth...	0.774	0.497	0.824	0.772	0.779	0.800	0.758	0.743	14.91	
38)	1-Methylnaphth...	0.741	0.469	0.749	0.725	0.730	0.732	0.706	0.693	14.40	

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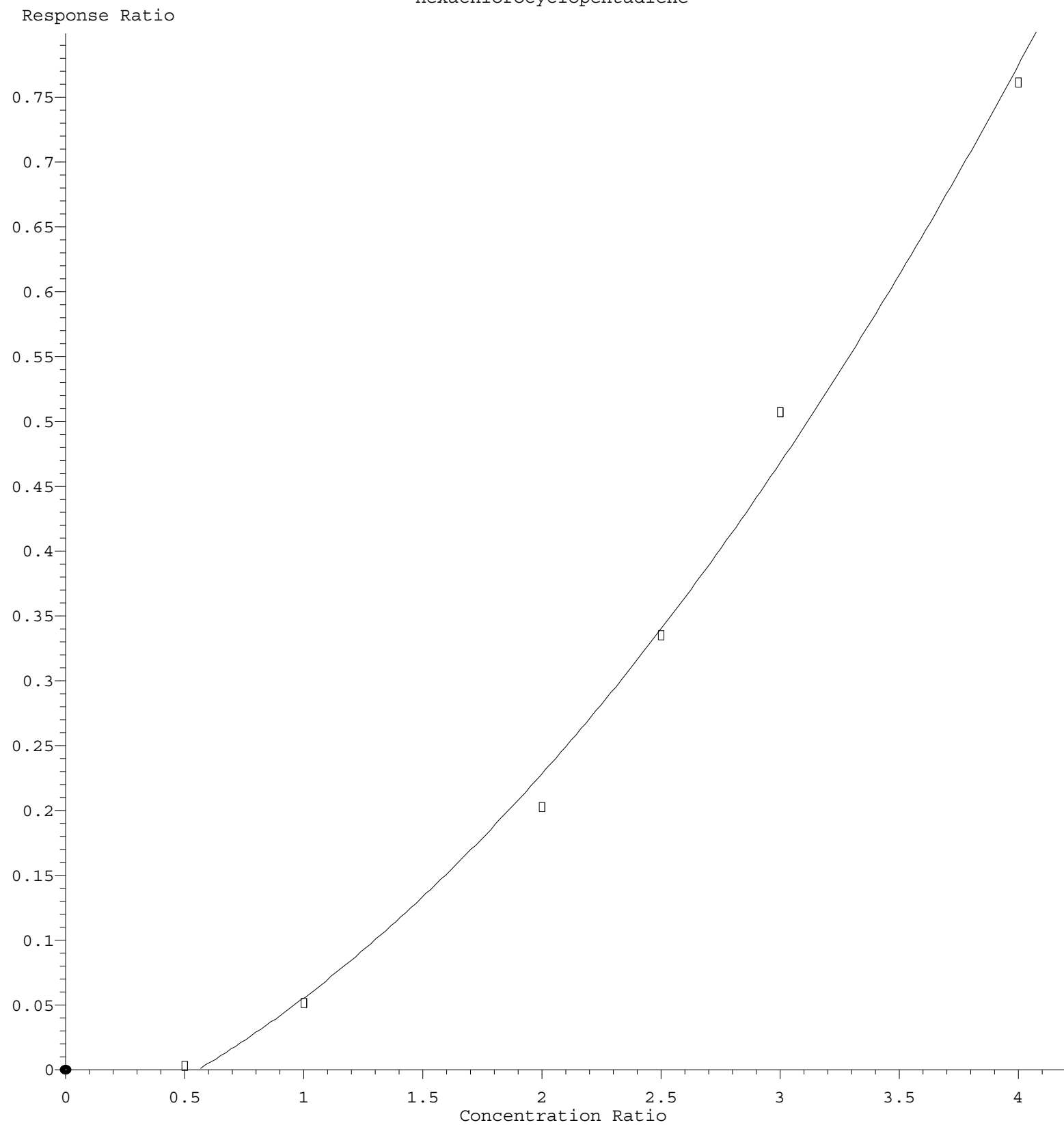
39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.647	0.629	0.643	0.624	0.657	0.679	0.635	0.645	2.93	
41) P	Hexachlorocycl...	0.006	0.051	0.101	0.134	0.169	0.190	0.109		64.88	
42) S	2,4,6-Tribromo...	0.191	0.313	0.283	0.273	0.327	0.346	0.324	0.294	17.65	
43) C	2,4,6-Trichlor...	0.333	0.369	0.400	0.397	0.423	0.432	0.409	0.395	8.61	
44)	2,4,5-Trichlor...	0.408	0.437	0.477	0.446	0.484	0.511	0.466	0.461	7.37	
45) S	2-Fluorobiphenyl	1.616	1.539	1.521	1.388	1.460	1.518	1.390	1.490	5.56	
46)	1,1'-Biphenyl	1.625	1.466	1.522	1.420	1.442	1.504	1.379	1.480	5.44	
47)	2-Chloronaphth...	1.116	1.113	1.158	1.065	1.092	1.131	1.021	1.099	4.12	
48)	2-Nitroaniline	0.420	0.420	0.450	0.437	0.477	0.495	0.447	0.449	6.23	
49)	Acenaphthylene	1.977	1.816	1.927	1.798	1.854	1.893	1.714	1.854	4.73	
50)	Dimethylphthalate	1.606	1.513	1.558	1.470	1.537	1.573	1.461	1.531	3.49	
51)	2,6-Dinitrotol...	0.347	0.307	0.339	0.327	0.333	0.326	0.309	0.327	4.46	
52) C	Acenaphthene	1.091	1.138	1.118	1.064	1.094	1.121	1.021	1.092	3.63	
53)	3-Nitroaniline	0.318	0.312	0.353	0.339	0.345	0.348	0.324	0.334	4.81	
54) P	2,4-Dinitrophenol	0.037	0.100	0.128	0.147	0.171	0.161	0.124		39.98	
55)	Dibenzofuran	1.389	1.862	1.901	1.791	1.862	1.875	1.754	1.776	10.03	
56) P	4-Nitrophenol	0.140	0.174	0.196	0.220	0.207	0.203	0.190		15.17	
57)	2,4-Dinitrotol...	0.325	0.458	0.465	0.456	0.474	0.488	0.448	0.445	12.24	
58)	Fluorene	0.989	1.583	1.561	1.445	1.497	1.529	1.410	1.431	14.28	
59)	2,3,4,6-Tetrac...	0.226	0.398	0.357	0.374	0.408	0.433	0.411	0.373	18.60	
60)	Diethylphthalate	1.208	1.626	1.621	1.490	1.614	1.686	1.501	1.535	10.47	
61)	4-Chlorophenyl...	0.625	0.872	0.828	0.769	0.836	0.875	0.808	0.802	10.72	
62)	4-Nitroaniline	0.218	0.321	0.349	0.351	0.369	0.370	0.329	0.330	15.91	
63)	Azobenzene	1.207	1.613	1.622	1.537	1.560	1.590	1.465	1.513	9.60	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.074	0.079	0.110	0.251	0.117	0.122	0.126	0.126	47.17	
66) c	n-Nitrosodiphe...	0.696	0.560	0.575	1.199	0.521	0.550	0.546	0.664	36.53#	
67)	4-Bromophenyl-...	0.282	0.231	0.221	0.325	0.226	0.235	0.235	0.251	15.39	
68)	Hexachlorobenzene	0.249	0.240	0.236	0.229	0.235	0.245	0.243	0.240	2.83	
69)	Atrazine	0.237	0.212	0.214	0.190	0.203	0.208	0.200	0.209	7.03	
70) C	Pentachlorophenol	0.043	0.075	0.089	0.097	0.105	0.115	0.087		29.53	
71)	Phenanthrene	1.323	1.036	1.048	1.001	0.965	0.992	0.971	1.048	11.93	
72)	Anthracene	1.383	1.080	1.078	1.021	1.003	1.022	1.008	1.085	12.45	
73)	Carbazole	1.396	1.000	0.979	0.969	0.889	0.908	0.887	1.004	17.80	
74)	Di-n-butylphth...	2.196	1.209	1.176	1.567	1.105	1.115	1.089	1.351	30.17	
75) C	Fluoranthene	2.513	2.078	1.279	1.312	1.205	1.227	1.200	1.545	34.28#	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.512	0.348	0.504	0.663	0.521	0.529	0.471	0.507	18.32	
78)	Pyrene	1.379	1.047	1.422	2.190	1.335	1.407	1.349	1.447	24.29	
79) S	Terphenyl-d14	1.131	0.827	1.111	1.651	1.129	1.167	1.121	1.162	21.02	
80)	Butylbenzylpht...	0.516	0.399	0.569	0.568	0.545	0.560	0.536	0.527	11.31	
81)	Benzo(a)anthra...	1.403	1.436	1.458	1.385	1.370	1.432	1.379	1.409	2.36	
82)	3,3'-Dichlorob...	0.483	0.494	0.528	0.501	0.492	0.516	0.481	0.499	3.48	
83)	Chrysene	1.331	1.349	1.401	1.315	1.296	1.351	1.284	1.333	2.95	
84)	Bis(2-ethylhex...	0.763	0.777	0.835	0.805	0.773	0.809	0.768	0.790	3.38	
85) c	Di-n-octyl pht...	1.352	1.354	1.447	1.378	1.340	1.378	1.302	1.364	3.27	

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86) I	Perylene-d12	-----ISTD-----								
87)	Indeno(1,2,3-c...	1.303	1.337	1.382	1.277	1.311	1.425	1.355	1.342	3.78
88)	Benzo(b)fluora...	1.149	1.165	1.214	1.167	1.162	1.209	1.157	1.175	2.20
89)	Benzo(k)fluora...	1.229	1.223	1.254	1.154	1.117	1.221	1.157	1.194	4.23
90) C	Benzo(a)pyrene	1.118	1.142	1.190	1.124	1.101	1.195	1.122	1.142	3.21
91)	Dibenzo(a,h)an...	1.097	1.110	1.154	1.060	1.107	1.192	1.128	1.121	3.77
92)	Benzo(g,h,i)pe...	1.098	1.106	1.139	0.993	1.092	1.153	1.087	1.095	4.71

(#) = Out of Range

Hexachlorocyclopentadiene



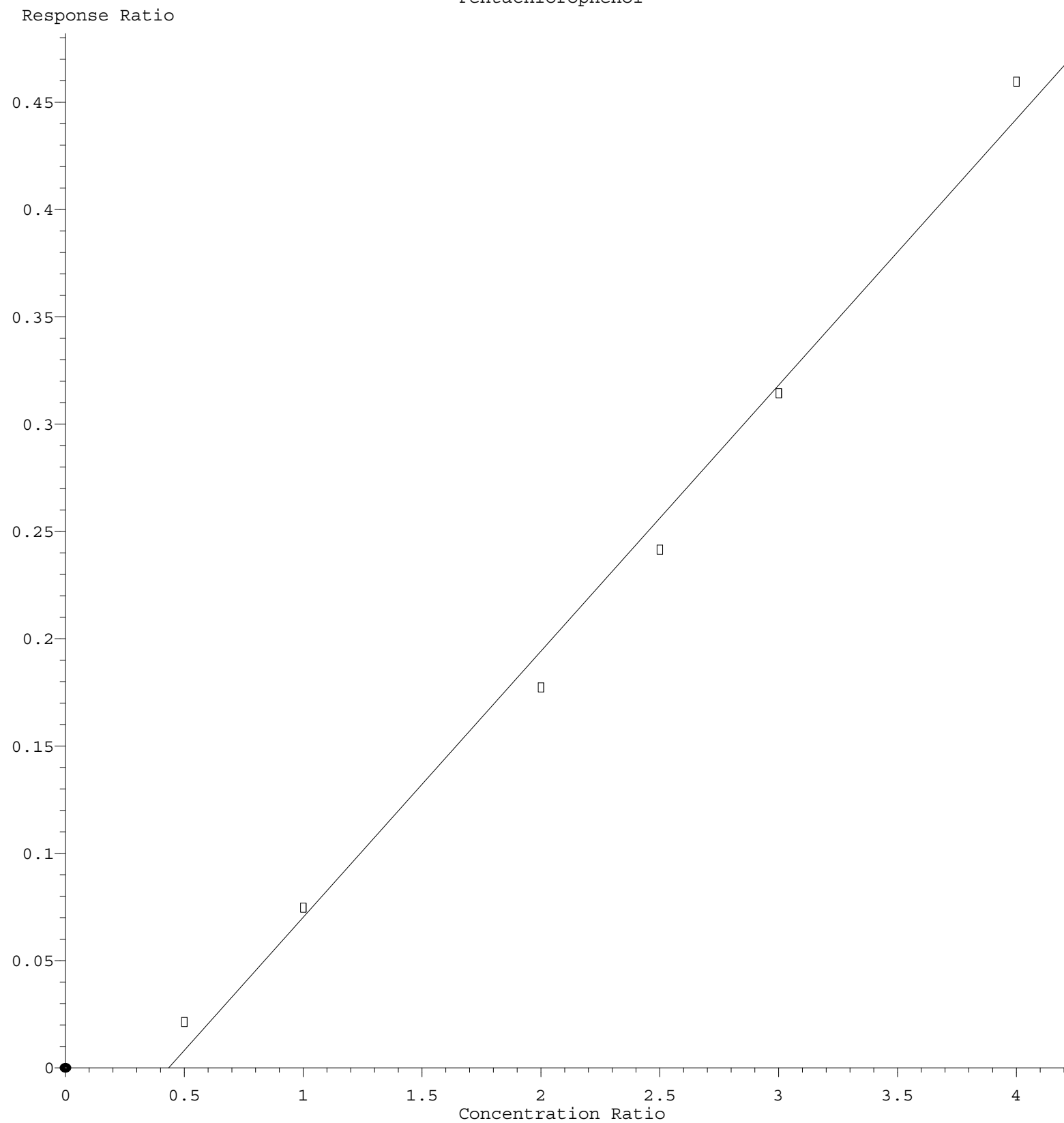
$$R = 3.344e-002 A^2 + 7.260e-002 A - 5.069e-002$$

Coef of Det (r^2) = 0.994035 Curve Fit: Quadratic

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Calibration Table Last Updated: Fri Jun 09 00:44:37 2023

Pentachlorophenol



Response = 1.239e-001 * Amt - 5.375e-002
 Coef of Det (r^2) = 0.992154 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG060823.M
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