

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF122023.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Dec 21 01:54:25 2023
 Response Via : Initial Calibration

Calibration Files

2.5 =BF136641.D 5 =BF136642.D 10 =BF136643.D 20 =BF136644.D 40 =BF136645.D 50 =BF136646.D 60 =BF136647.D 80 =BF136648.D

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

1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.574	0.559	0.558	0.520	0.506	0.523	0.499	0.534	5.47	
3)	Pyridine	1.291	1.358	1.339	1.317	1.279	1.310	1.226	1.303	3.32	
4)	n-Nitrosodimet...	0.699	0.672	0.713	0.715	0.687	0.709	0.677	0.696	2.51	
5) S	2-Fluorophenol	1.352	1.364	1.358	1.291	1.216	1.235	1.157	1.282	6.34	
6)	Aniline	1.820	1.749	1.774	1.679	1.552	1.571	1.478	1.661	7.76	
7) S	Phenol-d6	1.780	1.725	1.741	1.669	1.589	1.617	1.505	1.661	5.82	
8)	2-Chlorophenol	1.495	1.451	1.498	1.418	1.326	1.360	1.271	1.403	6.20	
9)	Benzaldehyde	1.123	1.062	1.041	0.933	0.850	0.854	0.751	0.945	14.32	
10) C	Phenol	1.929	1.859	1.870	1.770	1.654	1.721	1.607	1.773	6.73	
11)	bis(2-Chloroet...	1.393	1.403	1.435	1.366	1.285	1.313	1.243	1.348	5.16	
12)	1,3-Dichlorobe...	1.617	1.583	1.608	1.544	1.456	1.490	1.414	1.530	5.13	
13) C	1,4-Dichlorobe...	1.694	1.631	1.646	1.571	1.473	1.517	1.413	1.564	6.46	
14)	1,2-Dichlorobe...	1.599	1.533	1.551	1.460	1.367	1.390	1.284	1.455	7.81	
15)	Benzyl Alcohol	1.137	1.156	1.217	1.145	1.077	1.093	1.020	1.121	5.66	
16)	2,2'-oxybis(1-...	2.360	2.272	2.324	2.237	2.103	2.134	2.002	2.205	5.87	
17)	2-Methylphenol	1.291	1.198	1.200	1.158	1.103	1.119	1.065	1.162	6.49	
18)	Hexachloroethane	0.608	0.581	0.598	0.568	0.547	0.556	0.516	0.568	5.52	
19) P	n-Nitroso-di-n...	1.012	1.090	1.051	1.056	1.001	0.939	0.983	0.923	1.007	5.74
20)	3+4-Methylphenols	1.620	1.556	1.548	1.450	1.342	1.370	1.277	1.452	8.79	
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.548	0.505	0.528	0.510	0.466	0.487	0.466	0.501	6.14	
23) S	Nitrobenzene-d5	0.406	0.388	0.406	0.404	0.370	0.388	0.374	0.391	3.86	
24)	Nitrobenzene	0.402	0.389	0.402	0.409	0.375	0.391	0.372	0.392	3.58	
25)	Isophorone	0.735	0.712	0.732	0.728	0.679	0.706	0.688	0.711	3.08	
26) C	2-Nitrophenol	0.181	0.178	0.196	0.198	0.180	0.192	0.185	0.187	4.35	
27)	2,4-Dimethylph...	0.224	0.225	0.240	0.235	0.220	0.233	0.219	0.228	3.62	
28)	bis(2-Chloroet...	0.451	0.433	0.441	0.450	0.411	0.425	0.416	0.432	3.61	
29) C	2,4-Dichloroph...	0.301	0.294	0.309	0.304	0.282	0.290	0.276	0.294	4.03	
30)	1,2,4-Trichlor...	0.341	0.328	0.346	0.337	0.310	0.321	0.307	0.327	4.60	
31)	Naphthalene	1.191	1.111	1.138	1.126	1.033	1.077	1.017	1.099	5.57	
32)	Benzoic acid	0.171	0.180	0.227	0.246	0.231	0.248	0.242	0.221	14.37	
33)	4-Chloroaniline	0.430	0.410	0.427	0.417	0.382	0.392	0.381	0.406	5.08	
34) C	Hexachlorobuta...	0.212	0.200	0.207	0.203	0.188	0.195	0.186	0.199	4.90	
35)	Caprolactam	0.102	0.099	0.095	0.101	0.092	0.096	0.092	0.097	4.07	
36) C	4-Chloro-3-met...	0.345	0.332	0.347	0.340	0.315	0.324	0.311	0.331	4.31	
37)	2-Methylnaphth...	0.776	0.735	0.740	0.720	0.657	0.676	0.646	0.707	6.83	
38)	1-Methylnaphth...	0.764	0.716	0.739	0.701	0.641	0.661	0.635	0.694	7.15	

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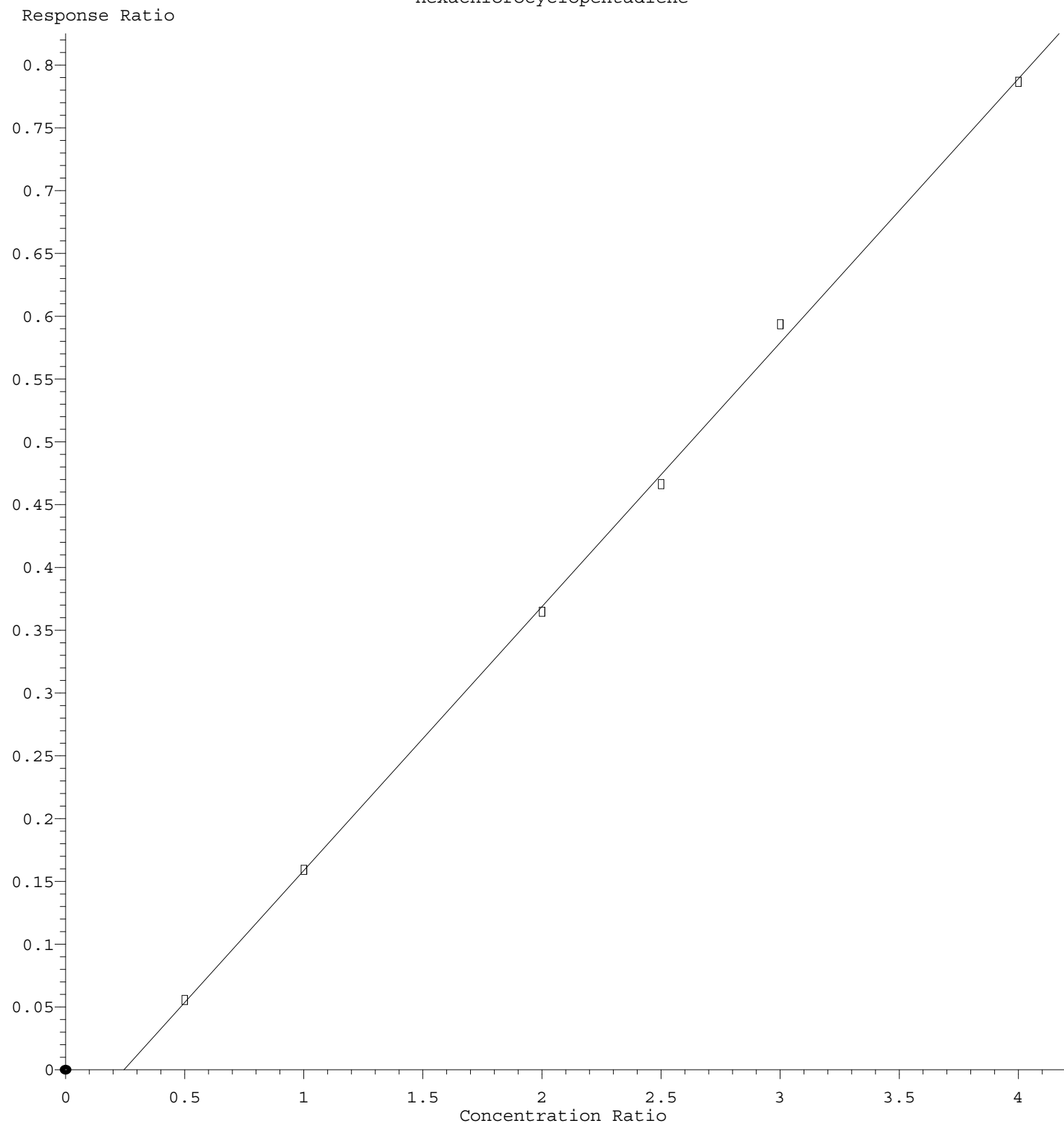
39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.657	0.634	0.653	0.634	0.571	0.614	0.574	0.619	5.64	
41) P	Hexachlorocycl...	0.111	0.159	0.182	0.186	0.198	0.197	0.172	19.18		
42) S	2,4,6-Tribromo...	0.253	0.249	0.253	0.240	0.216	0.226	0.211	0.236	7.47	
43) C	2,4,6-Trichlor...	0.404	0.397	0.414	0.409	0.378	0.400	0.379	0.397	3.46	
44)	2,4,5-Trichlor...	0.456	0.457	0.454	0.455	0.413	0.444	0.420	0.443	4.24	
45) S	2-Fluorobiphenyl	1.709	1.599	1.587	1.474	1.338	1.420	1.282	1.487	10.31	
46)	1,1'-Biphenyl	1.823	1.763	1.777	1.681	1.555	1.636	1.512	1.678	6.98	
47)	2-Chloronaphth...	1.405	1.312	1.331	1.289	1.182	1.248	1.164	1.276	6.65	
48)	2-Nitroaniline	0.399	0.410	0.413	0.409	0.382	0.401	0.384	0.400	3.08	
49)	Acenaphthylene	2.105	2.052	2.068	1.984	1.793	1.899	1.750	1.950	7.15	
50)	Dimethylphthalate	1.563	1.517	1.522	1.463	1.373	1.432	1.355	1.461	5.38	
51)	2,6-Dinitrotol...	0.344	0.334	0.352	0.339	0.314	0.331	0.307	0.332	4.88	
52) C	Acenaphthene	1.324	1.287	1.259	1.215	1.115	1.167	1.093	1.209	7.23	
53)	3-Nitroaniline	0.369	0.364	0.371	0.364	0.325	0.344	0.321	0.351	6.05	
54) P	2,4-Dinitrophenol	0.119	0.153	0.184	0.169	0.185	0.183	0.165	15.68		
55)	Dibenzofuran	2.080	1.962	1.961	1.835	1.663	1.739	1.587	1.832	9.77	
56) P	4-Nitrophenol	0.214	0.234	0.244	0.252	0.237	0.245	0.234	0.237	5.20	
57)	2,4-Dinitrotol...	0.451	0.459	0.466	0.446	0.394	0.416	0.381	0.430	7.83	
58)	Fluorene	1.644	1.599	1.548	1.451	1.309	1.364	1.263	1.454	10.19	
59)	2,3,4,6-Tetrac...	0.338	0.344	0.375	0.356	0.327	0.335	0.311	0.341	6.05	
60)	Diethylphthalate	1.660	1.616	1.590	1.539	1.395	1.456	1.353	1.516	7.69	
61)	4-Chlorophenyl...	0.818	0.766	0.752	0.711	0.638	0.660	0.603	0.707	10.89	
62)	4-Nitroaniline	0.385	0.376	0.346	0.333	0.307	0.320	0.303	0.339	9.60	
63)	Azobenzene	1.513	1.493	1.511	1.421	1.306	1.378	1.288	1.416	6.72	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.084	0.099	0.120	0.130	0.117	0.128	0.125	0.115	14.61	
66) c	n-Nitrosodiphe...	0.689	0.666	0.676	0.675	0.602	0.638	0.601	0.650	5.59	
67)	4-Bromophenyl-...	0.237	0.226	0.232	0.229	0.209	0.216	0.206	0.222	5.33	
68)	Hexachlorobenzene	0.257	0.255	0.263	0.255	0.229	0.242	0.231	0.248	5.45	
69)	Atrazine	0.196	0.188	0.165	0.144	0.140	0.164	0.166	13.77		
70) C	Pentachlorophenol	0.089	0.111	0.121	0.130	0.118	0.122	0.118	0.115	11.33	
71)	Phenanthrene	1.121	1.090	1.075	1.062	0.941	0.974	0.920	1.026	7.75	
72)	Anthracene	1.155	1.088	1.094	1.059	0.953	0.997	0.928	1.039	7.92	
73)	Carbazole	1.081	1.047	1.037	1.004	0.899	0.919	0.871	0.980	8.40	
74)	Di-n-butylphth...	1.284	1.281	1.263	1.225	1.108	1.146	1.062	1.196	7.53	
75) C	Fluoranthene	1.272	1.217	1.161	1.119	0.983	1.005	0.929	1.098	11.73	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.709	0.595	0.664	0.560	0.556	0.585	0.460	0.590	13.58	
78)	Pyrene	1.907	1.886	2.058	2.047	1.908	1.994	1.943	1.963	3.58	
79) S	Terphenyl-d14	1.491	1.447	1.509	1.486	1.348	1.399	1.338	1.431	4.91	
80)	Butylbenzylpht...	0.679	0.678	0.748	0.755	0.698	0.730	0.712	0.714	4.39	
81)	Benzo(a)anthra...	1.501	1.419	1.435	1.428	1.296	1.344	1.298	1.389	5.58	
82)	3,3'-Dichlorob...	0.405	0.397	0.419	0.395	0.374	0.396	0.375	0.394	4.01	
83)	Chrysene	1.441	1.384	1.424	1.373	1.262	1.343	1.281	1.358	4.99	
84)	Bis(2-ethylhex...	0.849	0.869	0.929	0.946	0.891	0.917	0.892	0.899	3.79	
85) c	Di-n-octyl pht...	1.242	1.271	1.379	1.452	1.424	1.504	1.460	1.390	7.14	

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86) I	Perylene-d12	-----ISTD-----								
87)	Indeno(1,2,3-c...	1.323	1.312	1.428	1.517	1.483	1.540	1.505	1.444	6.43
88)	Benzo(b)fluora...	1.483	1.341	1.421	1.321	1.238	1.330	1.218	1.336	7.02
89)	Benzo(k)fluora...	1.416	1.371	1.288	1.304	1.193	1.120	1.144	1.262	8.96
90) C	Benzo(a)pyrene	1.182	1.132	1.153	1.149	1.094	1.101	1.081	1.128	3.25
91)	Dibenzo(a,h)an...	1.091	1.092	1.185	1.264	1.205	1.250	1.206	1.185	5.84
92)	Benzo(g,h,i)pe...	1.073	1.087	1.211	1.285	1.266	1.302	1.268	1.213	7.86

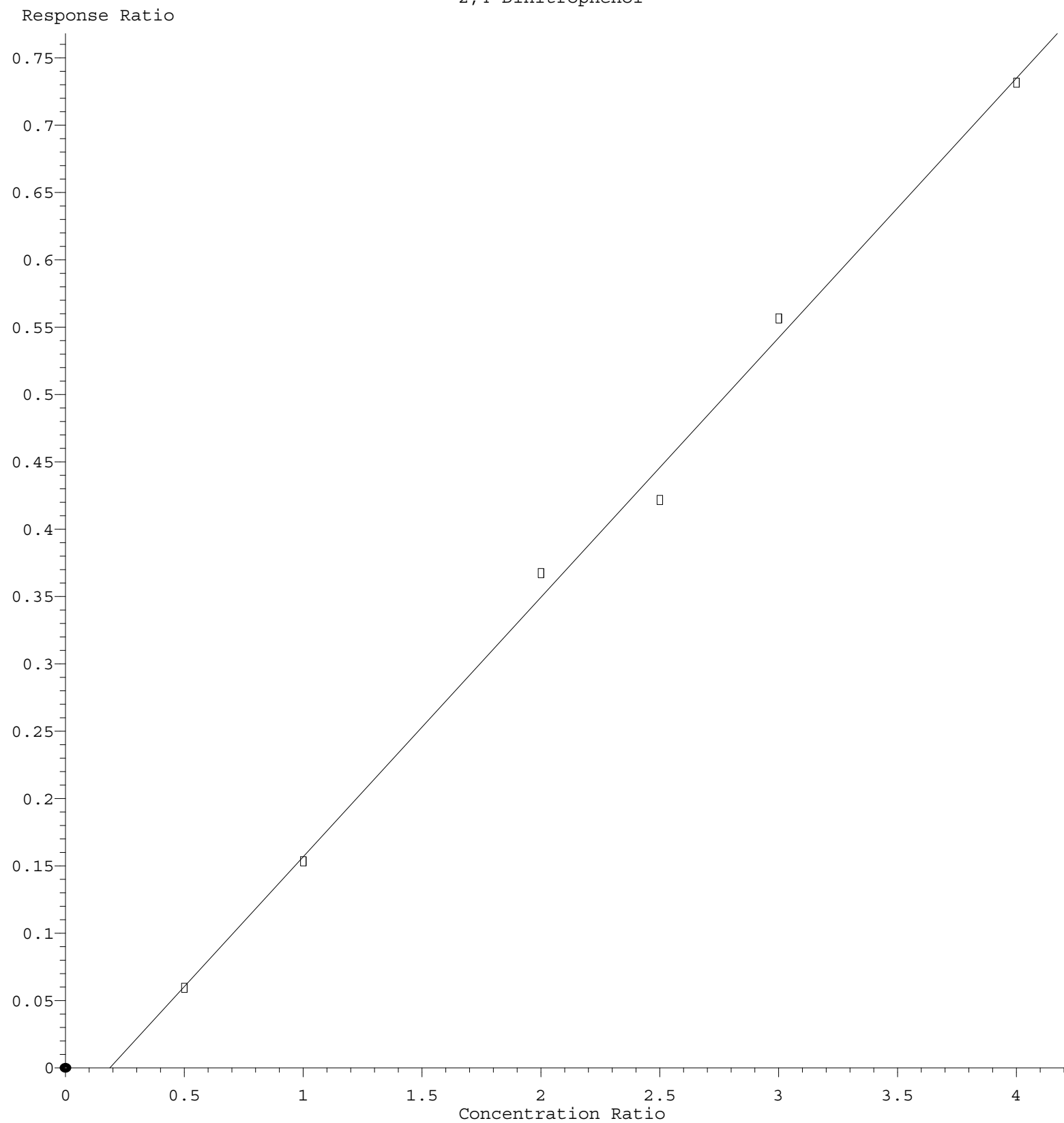
(#) = Out of Range

Hexachlorocyclopentadiene



Response = $2.104 \times 10^{-1} \times \text{Amt} - 5.152 \times 10^{-2}$
Coef of Det (r^2) = 0.999195 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF122023.M
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2,4-Dinitrophenol



Response = 1.927e-001 * Amt - 3.600e-002
 Coef of Det (r^2) = 0.996353 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF122023.M
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