

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG091922.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Mon Sep 19 17:41:40 2022
 Response Via : Initial Calibration

Calibration Files

5 =BG054969.D 10 =BG054970.D 20 =BG054971.D 40 =BG054972.D 50 =BG054973.D 60 =BG054974.D 80 =BG054975.D

	Compound	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----								
2)	1,4-Dioxane	0.564	0.516	0.491	0.467	0.485	0.488	0.466	0.497	6.87
3)	Pyridine	1.331	1.328	1.399	1.337	1.347	1.381	1.324	1.349	2.14
4)	n-Nitrosodimet...	0.487	0.480	0.518	0.481	0.500	0.498	0.482	0.492	2.85
5) S	2-Fluorophenol	1.032	1.014	1.064	1.013	1.023	1.038	0.971	1.022	2.79
6)	Aniline	1.683	1.707	1.810	1.750	1.707	1.726	1.613	1.714	3.54
7) S	Phenol-d6	1.306	1.364	1.444	1.408	1.365	1.401	1.302	1.370	3.84
8)	2-Chlorophenol	1.207	1.205	1.283	1.211	1.180	1.208	1.133	1.204	3.70
9)	Benzaldehyde	1.019	1.009	0.968	0.852	0.814	0.771	0.677	0.873	14.92
10) C	Phenol	1.439	1.467	1.575	1.527	1.473	1.509	1.404	1.485	3.85
11)	bis(2-Chloroet...	1.248	1.171	1.218	1.169	1.144	1.169	1.101	1.174	4.07
12)	1,3-Dichlorobe...	1.489	1.413	1.473	1.360	1.377	1.406	1.313	1.404	4.41
13) C	1,4-Dichlorobe...	1.511	1.442	1.506	1.399	1.400	1.426	1.334	1.431	4.38
14)	1,2-Dichlorobe...	1.405	1.379	1.435	1.336	1.309	1.331	1.246	1.349	4.70
15)	Benzyl Alcohol	0.906	0.959	1.057	1.063	1.004	1.041	0.984	1.002	5.70
16)	2,2'-oxybis(1-...	1.439	1.468	1.503	1.454	1.416	1.439	1.354	1.439	3.22
17)	2-Methylphenol	0.971	1.027	1.120	1.095	1.031	1.076	1.006	1.047	5.02
18)	Hexachloroethane	0.502	0.506	0.514	0.480	0.479	0.499	0.464	0.492	3.66
19) P	n-Nitroso-di-n...	0.900	0.927	1.008	0.997	0.928	0.949	0.892	0.943	4.74
20)	3+4-Methylphenols	1.381	1.436	1.528	1.536	1.450	1.505	1.410	1.464	4.11
21) I	Naphthalene-d8	-----ISTD-----								
22)	Acetophenone	0.476	0.467	0.490	0.465	0.466	0.480	0.464	0.473	2.08
23) S	Nitrobenzene-d5	0.336	0.337	0.355	0.339	0.346	0.354	0.348	0.345	2.31
24)	Nitrobenzene	0.333	0.338	0.359	0.343	0.345	0.356	0.347	0.346	2.62
25)	Isophorone	0.654	0.624	0.671	0.661	0.648	0.671	0.658	0.655	2.50
26) C	2-Nitrophenol	0.168	0.169	0.184	0.184	0.186	0.193	0.190	0.182	5.49
27)	2,4-Dimethylph...	0.242	0.243	0.269	0.260	0.257	0.267	0.261	0.257	4.21
28)	bis(2-Chloroet...	0.364	0.355	0.382	0.368	0.367	0.376	0.370	0.369	2.26
29) C	2,4-Dichloroph...	0.271	0.289	0.320	0.312	0.313	0.325	0.317	0.307	6.29
30)	1,2,4-Trichlor...	0.350	0.349	0.367	0.348	0.347	0.361	0.353	0.354	2.11
31)	Naphthalene	1.066	1.051	1.084	1.021	1.026	1.049	1.023	1.046	2.28
32)	Benzoic acid	0.243	0.005	0.034	0.038	0.048	0.052	0.070		123.47
33)	4-Chloroaniline	0.400	0.401	0.443	0.433	0.428	0.443	0.436	0.426	4.37
34) C	Hexachlorobuta...	0.244	0.228	0.243	0.227	0.233	0.237	0.233	0.235	2.84
35)	Caprolactam	0.101	0.102	0.118	0.116	0.113	0.119	0.114	0.112	6.82
36) C	4-Chloro-3-met...	0.296	0.309	0.343	0.340	0.328	0.343	0.334	0.328	5.55
37)	2-Methylnaphth...	0.722	0.727	0.779	0.754	0.731	0.760	0.741	0.745	2.74
38)	1-Methylnaphth...	0.721	0.718	0.764	0.734	0.717	0.735	0.712	0.729	2.47

Method	Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\									
Method	File : 8270-BG091922.M									
39) I	Acenaphthene-d10	-----ISTD-----								
40)	1,2,4,5-Tetrac...	0.589	0.577	0.601	0.568	0.580	0.595	0.584	0.585	1.92
41) P	Hexachlorocycl...			0.014	0.044	0.056	0.078	0.092	0.057	52.76
42) S	2,4,6-Tribromo...	0.207	0.209	0.238	0.243	0.241	0.258	0.245	0.234	8.07
43) C	2,4,6-Trichlor...	0.296	0.303	0.337	0.344	0.348	0.363	0.361	0.336	7.89
44)	2,4,5-Trichlor...	0.341	0.338	0.401	0.394	0.404	0.412	0.410	0.386	8.32
45) S	2-Fluorobiphenyl	1.354	1.287	1.328	1.231	1.237	1.242	1.205	1.269	4.37
46)	1,1'-Biphenyl	1.392	1.339	1.398	1.314	1.326	1.356	1.320	1.349	2.54
47)	2-Chloronaphth...	1.074	1.047	1.094	1.030	1.040	1.061	1.037	1.055	2.17
48)	2-Nitroaniline	0.284	0.281	0.317	0.316	0.316	0.332	0.321	0.310	6.25
49)	Acenaphthylene	1.843	1.783	1.859	1.746	1.741	1.783	1.723	1.783	2.91
50)	Dimethylphthalate	1.590	1.500	1.579	1.495	1.477	1.513	1.448	1.514	3.45
51)	2,6-Dinitrotol...	0.314	0.309	0.331	0.323	0.321	0.333	0.322	0.322	2.62
52) C	Acenaphthene	1.092	1.055	1.138	1.088	1.089	1.127	1.100	1.098	2.49
53)	3-Nitroaniline	0.332	0.321	0.361	0.356	0.354	0.369	0.353	0.349	4.83
54) P	2,4-Dinitrophenol		0.020	0.052	0.087	0.088	0.110	0.113	0.078	45.94
55)	Dibenzofuran	1.811	1.725	1.795	1.705	1.693	1.698	1.651	1.726	3.35
56) P	4-Nitrophenol		0.135	0.178	0.203	0.205	0.228	0.213	0.194	17.02
57)	2,4-Dinitrotol...	0.428	0.429	0.469	0.469	0.464	0.484	0.463	0.458	4.65
58)	Fluorene	1.530	1.468	1.523	1.447	1.428	1.462	1.394	1.465	3.33
59)	2,3,4,6-Tetrac...	0.265	0.246	0.319	0.359	0.346	0.372	0.370	0.325	15.68
60)	Diethylphthalate	1.603	1.508	1.569	1.498	1.467	1.510	1.430	1.512	3.87
61)	4-Chlorophenyl...	0.764	0.732	0.762	0.730	0.727	0.746	0.719	0.740	2.41
62)	4-Nitroaniline	0.335	0.344	0.379	0.375	0.365	0.387	0.360	0.364	5.16
63)	Azobenzene	1.233	1.210	1.280	1.239	1.221	1.258	1.217	1.237	2.02
64) I	Phenanthrene-d10	-----ISTD-----								
65)	4,6-Dinitro-2-...		0.065	0.087	0.097	0.097	0.107	0.107	0.093	16.97
66) c	n-Nitrosodiphe...	0.535	0.530	0.563	0.527	0.523	0.535	0.531	0.535	2.43
67)	4-Bromophenyl-...	0.217	0.214	0.226	0.221	0.218	0.225	0.222	0.221	2.06
68)	Hexachlorobenzene	0.256	0.250	0.267	0.249	0.247	0.255	0.250	0.253	2.66
69)	Atrazine	0.221	0.203	0.211	0.200	0.197	0.200	0.190	0.203	4.91
70) C	Pentachlorophenol		0.021	0.039	0.049	0.056	0.065	0.069	0.050	35.80#
71)	Phenanthrene	1.068	1.018	1.071	0.999	0.988	1.014	0.972	1.019	3.72
72)	Anthracene	1.030	1.000	1.051	0.982	0.970	0.980	0.942	0.994	3.69
73)	Carbazole	0.947	0.902	0.959	0.906	0.882	0.910	0.862	0.910	3.73
74)	Di-n-butylphth...	1.215	1.129	1.190	1.126	1.093	1.105	1.041	1.128	5.22
75) C	Fluoranthene	1.427	1.303	1.356	1.267	1.231	1.263	1.183	1.290	6.30
76) I	Chrysene-d12	-----ISTD-----								
77)	Benzidine	0.622	0.491	0.469	0.484	0.463	0.495	0.443	0.496	11.87
78)	Pyrene	1.333	1.335	1.405	1.304	1.288	1.292	1.245	1.315	3.83
79) S	Terphenyl-d14	1.032	1.013	1.041	0.935	0.918	0.899	0.855	0.956	7.62
80)	Butylbenzylpht...	0.503	0.501	0.530	0.504	0.512	0.519	0.507	0.511	2.03
81)	Benzo(a)anthra...	1.335	1.276	1.334	1.264	1.264	1.282	1.248	1.286	2.71
82)	3,3'-Dichlorob...	0.442	0.446	0.467	0.440	0.443	0.450	0.435	0.446	2.29
83)	Chrysene	1.285	1.233	1.284	1.198	1.205	1.229	1.186	1.231	3.25
84)	Bis(2-ethylhex...	0.723	0.719	0.758	0.722	0.736	0.741	0.731	0.733	1.85
85) c	Di-n-octyl pht...	1.284	1.261	1.302	1.248	1.263	1.272	1.239	1.267	1.70

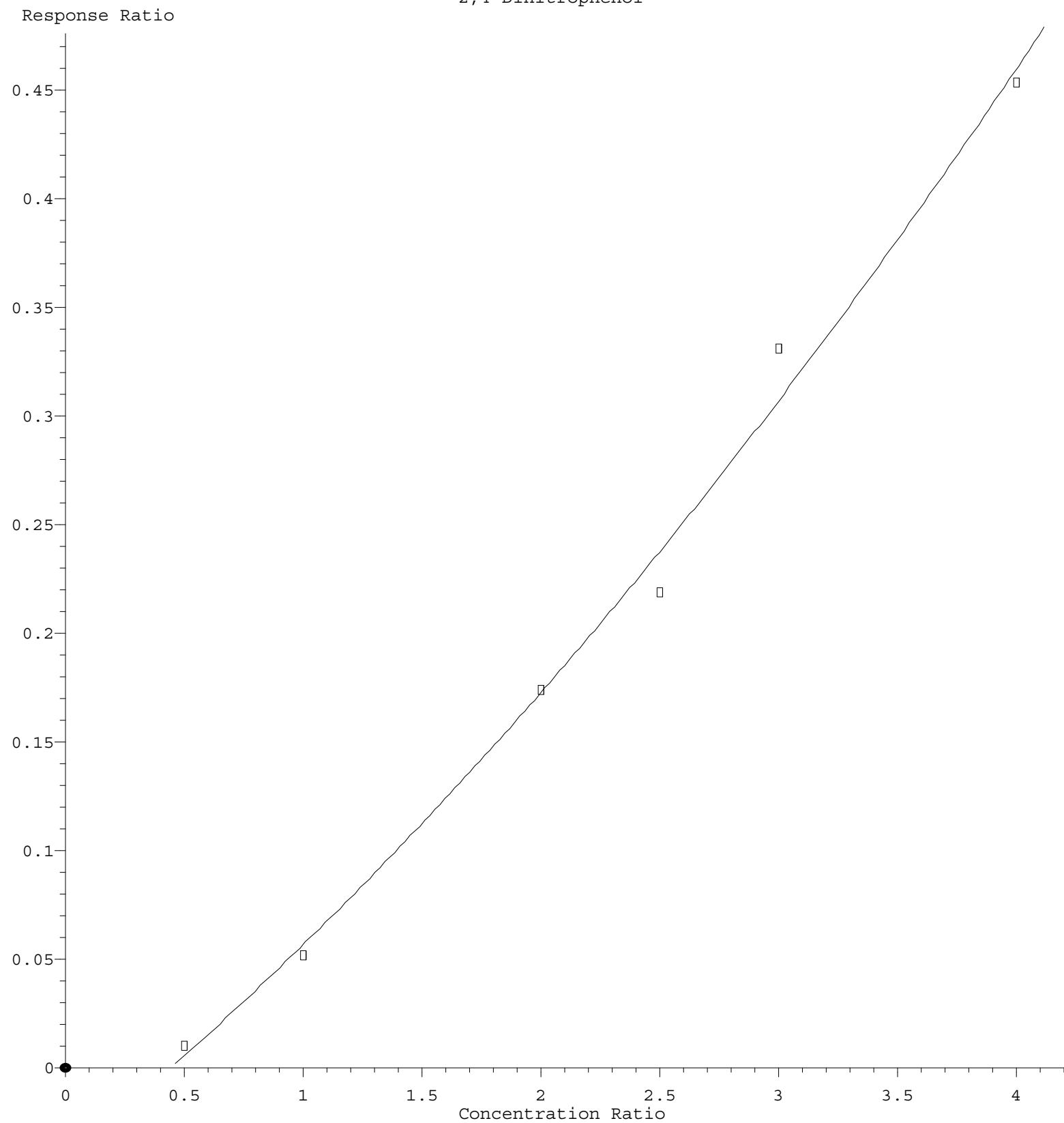
Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\

Method File : 8270-BG091922.M

86) I	Perylene-d12										
87)	Indeno(1,2,3-c...	1.466	1.583	1.495	1.431	1.451	1.504	1.478	1.487		3.30
88)	Benzo(b)fluora...	1.171	1.164	1.211	1.150	1.171	1.208	1.183	1.180		1.92
89)	Benzo(k)fluora...	1.217	1.195	1.240	1.185	1.173	1.178	1.148	1.191		2.54
90) C	Benzo(a)pyrene	1.019	1.013	1.045	0.992	1.004	1.030	1.007	1.016		1.73
91)	Dibenzo(a,h)an...	1.216	1.305	1.234	1.181	1.188	1.233	1.211	1.224		3.34
92)	Benzo(g,h,i)pe...	1.226	1.348	1.235	1.182	1.195	1.238	1.213	1.234		4.41

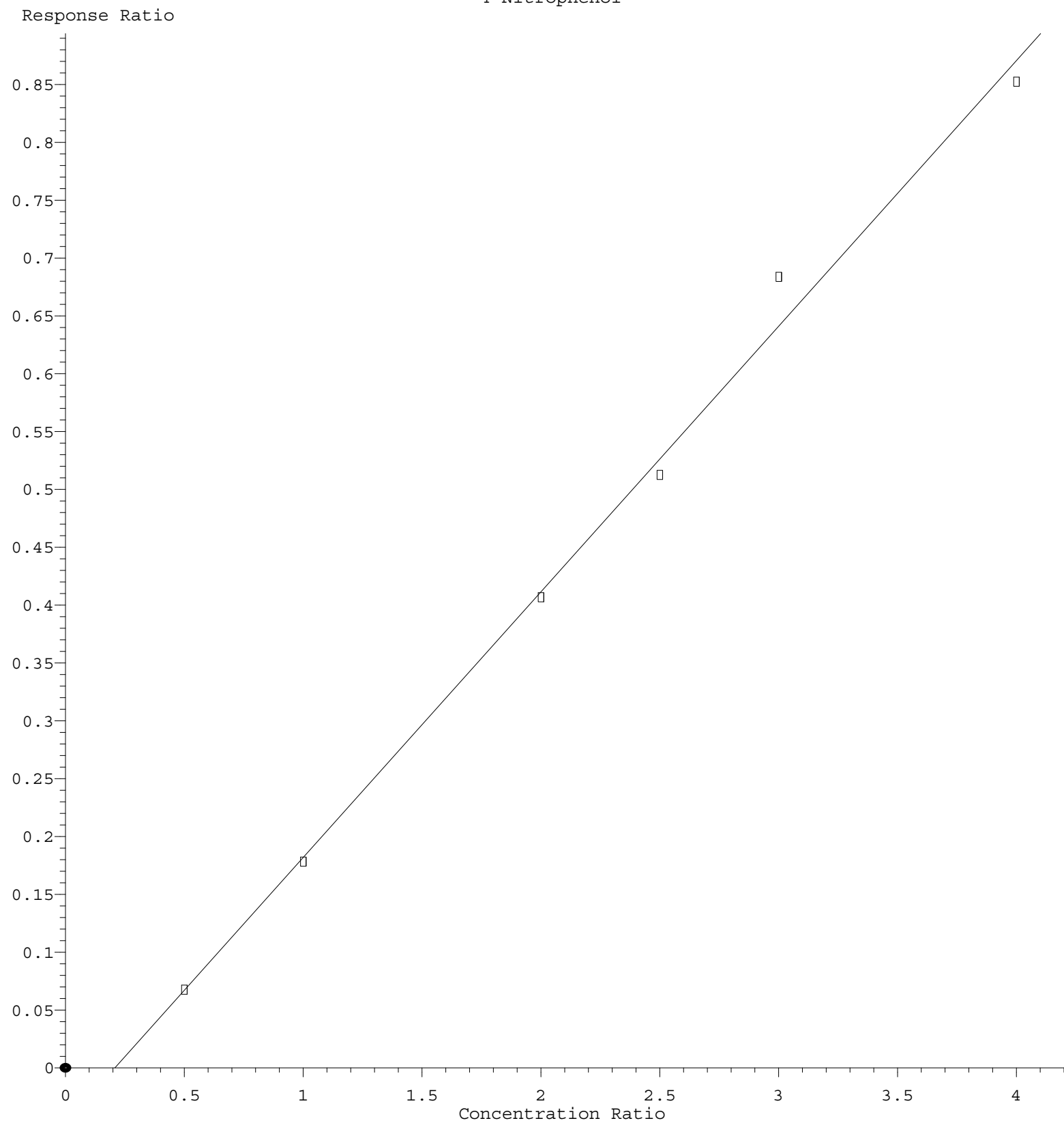
(#) = Out of Range

2,4-Dinitrophenol



$R = 9.139e-003 A^2 + 8.864e-002 A - 4.112e-002$
 Coef of Det (r^2) = 0.992777 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG091922.M
 Calibration Table Last Updated: Mon Sep 19 17:41:40 2022

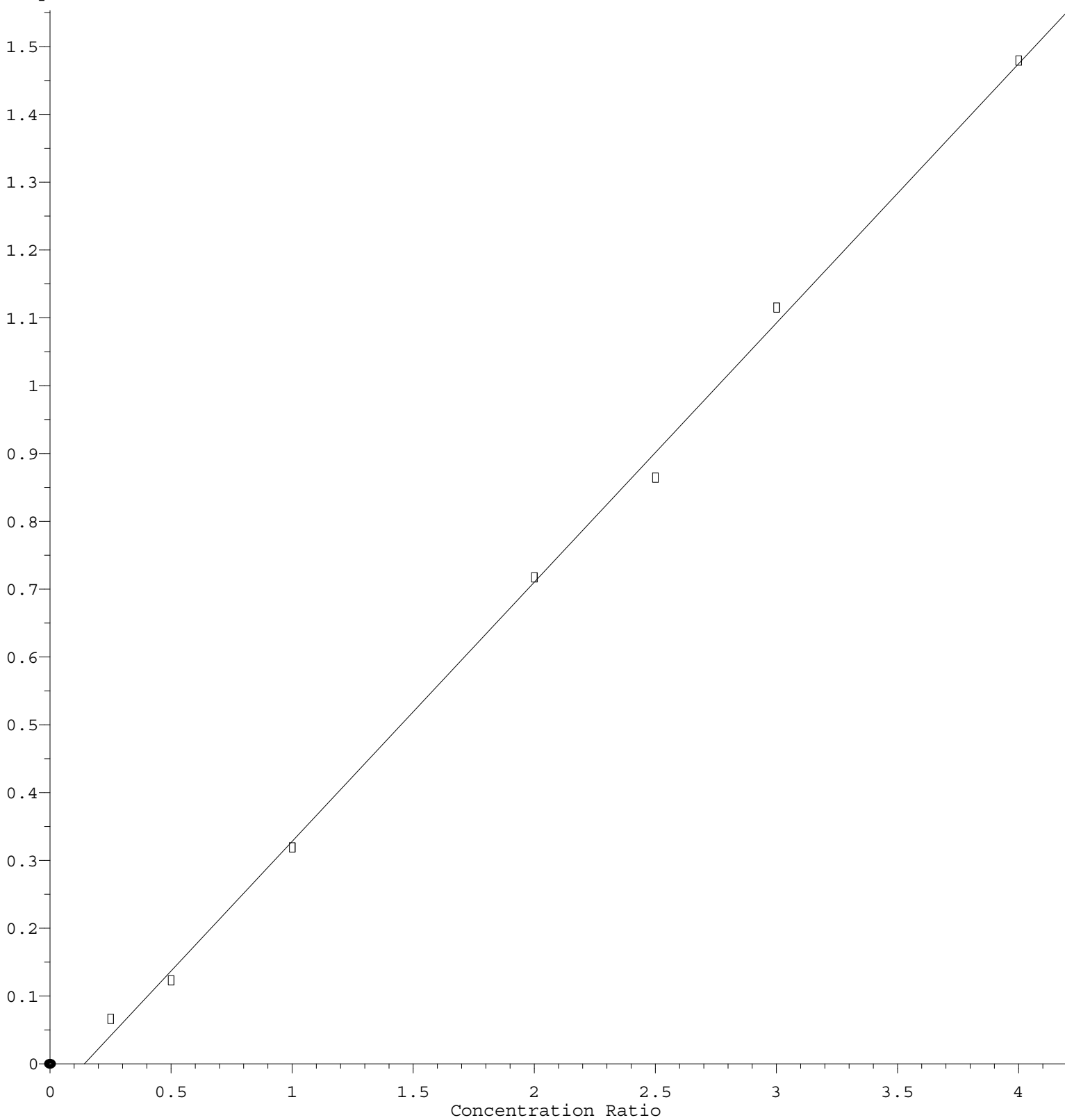
4-Nitrophenol



$\text{Response} = 2.298\text{e-}001 * \text{Amt} - 4.770\text{e-}002$
 Coef of Det (r^2) = 0.994640 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG091922.M
 Calibration Table Last Updated: Mon Sep 19 17:41:40 2022

2,3,4,6-Tetrachlorophenol

Response Ratio



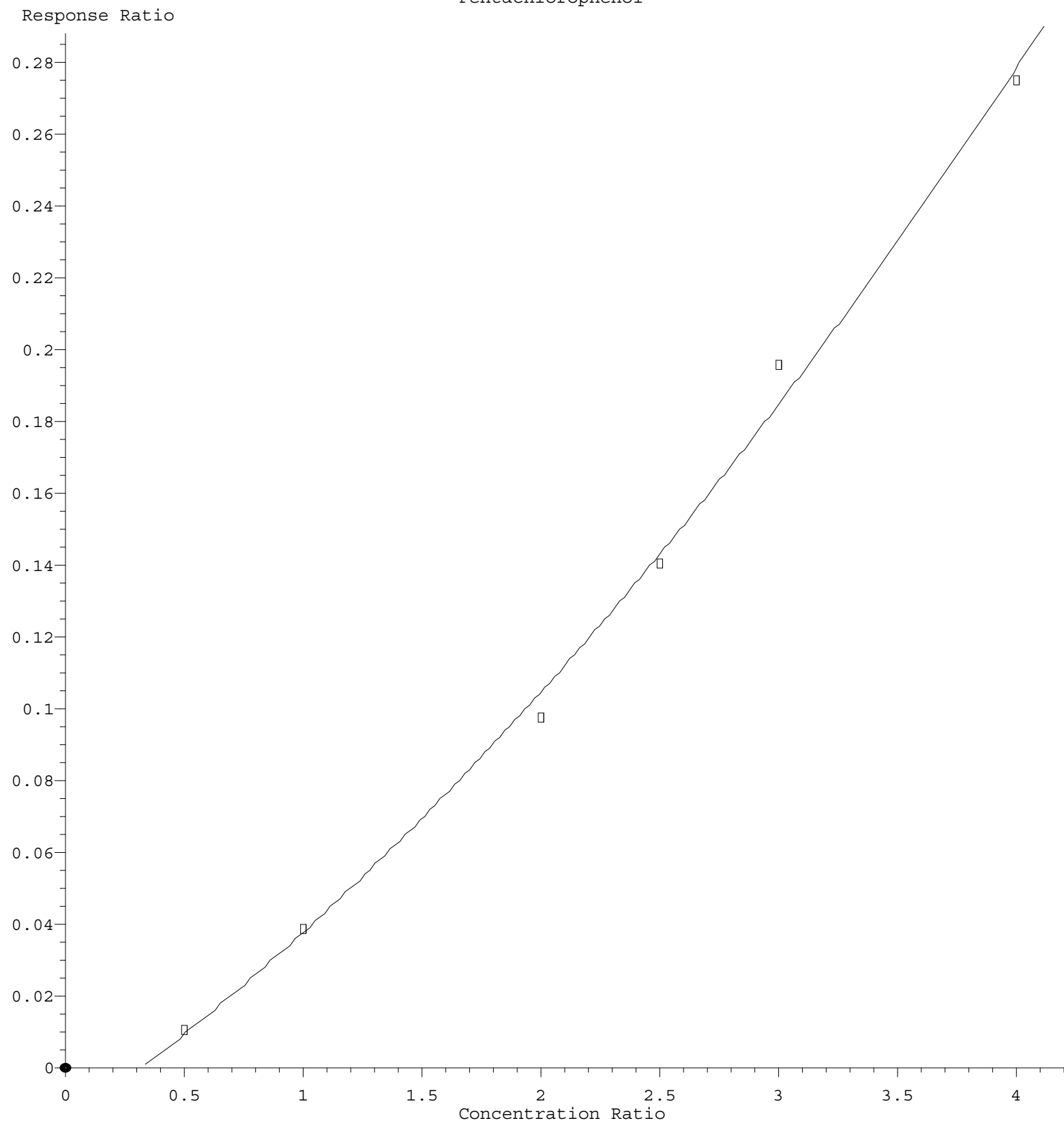
$$\text{Response} = 3.821\text{e-}001 * \text{Amt} - 5.415\text{e-}002$$

Coef of Det (r^2) = 0.998311 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG091922.M

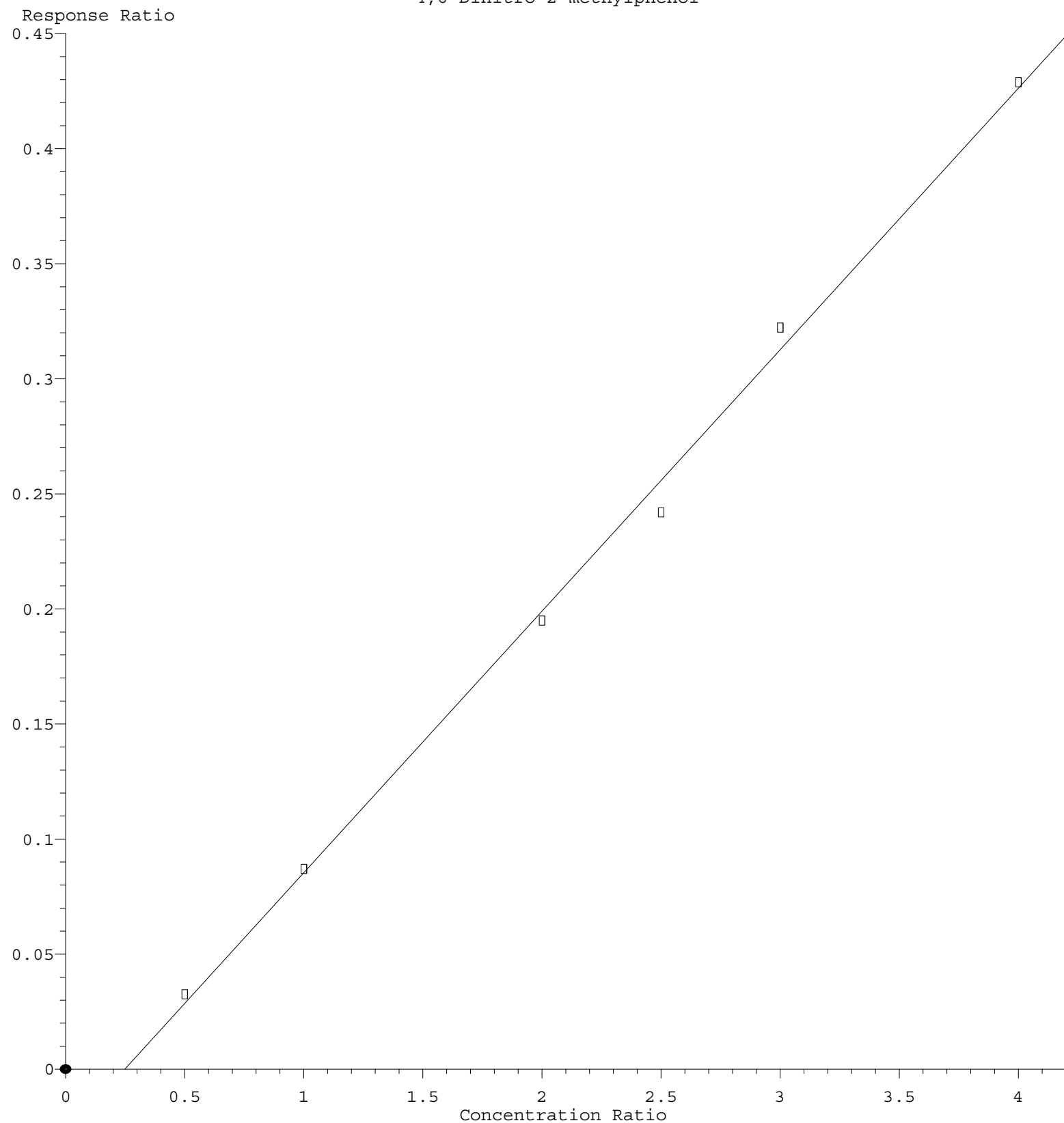
Calibration Table Last Updated: Mon Sep 19 17:41:40 2022

Pentachlorophenol



$R = 6.705e-003 A^2 + 4.671e-002 A - 1.568e-002$
 Coef of Det (r^2) = 0.996110 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG091922.M
 Calibration Table Last Updated: Mon Sep 19 17:41:40 2022

4,6-Dinitro-2-methylphenol



Response = $1.137e-001 * \text{Amt} - 2.835e-002$
 Coef of Det (r^2) = 0.996975 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG091922.M
 Calibration Table Last Updated: Mon Sep 19 17:41:40 2022