

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF060524.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Jun 06 05:19:29 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BF138117.D 5 =BF138118.D 10 =BF138119.D 20 =BF138120.D 40 =BF138121.D 50 =BF138122.D 60 =BF138123.D 80 =BF138124.D

	Compound	2.5	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane		0.579	0.551	0.564	0.527	0.512	0.532	0.500	0.538	5.24
3)	Pyridine		1.358	1.336	1.388	1.306	1.273	1.311	1.254	1.318	3.56
4)	n-Nitrosodimet...		0.673	0.662	0.683	0.646	0.627	0.639	0.601	0.647	4.35
5) S	2-Fluorophenol		1.267	1.240	1.270	1.161	1.108	1.147	1.022	1.174	7.82
6)	Aniline		1.584	1.523	1.567	1.499	1.474	1.498	1.393	1.505	4.20
7) S	Phenol-d6		1.616	1.559	1.595	1.470	1.416	1.439	1.299	1.485	7.61
8)	2-Chlorophenol		1.352	1.298	1.361	1.269	1.233	1.254	1.120	1.270	6.41
9)	Benzaldehyde			1.040	0.956	0.767	0.676	0.680	0.631	0.791	21.20
10) C	Phenol		1.638	1.577	1.626	1.517	1.469	1.497	1.355	1.525	6.48
11)	bis(2-Chloroet...		1.340	1.287	1.289	1.213	1.163	1.180	1.089	1.223	7.14
12)	1,3-Dichlorobe...		1.603	1.525	1.560	1.447	1.377	1.408	1.264	1.455	8.05
13) C	1,4-Dichlorobe...		1.626	1.555	1.569	1.451	1.383	1.415	1.257	1.465	8.70
14)	1,2-Dichlorobe...		1.545	1.473	1.497	1.371	1.293	1.322	1.126	1.375	10.49
15)	Benzyl Alcohol		1.119	1.084	1.128	1.062	1.023	1.040	0.897	1.051	7.41
16)	2,2'-oxybis(1-...		1.983	1.862	1.917	1.774	1.710	1.731	1.591	1.795	7.47
17)	2-Methylphenol		1.083	1.038	1.071	1.015	1.001	1.018	0.968	1.028	3.88
18)	Hexachloroethane		0.508	0.499	0.527	0.502	0.496	0.507	0.461	0.500	3.98
19) P	n-Nitroso-di-n...	0.959	1.041	0.993	1.006	0.925	0.895	0.905	0.849	0.947	6.81
20)	3+4-Methylphenols		1.413	1.355	1.409	1.298	1.214	1.230	1.033	1.279	10.50
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone		0.530	0.500	0.504	0.456	0.426	0.432	0.377	0.460	11.64
23) S	Nitrobenzene-d5		0.219	0.251	0.304	0.315	0.320	0.336	0.313	0.294	14.44
24)	Nitrobenzene		0.263	0.292	0.338	0.335	0.336	0.348	0.329	0.320	9.64
25)	Isophorone		0.703	0.658	0.678	0.637	0.617	0.633	0.608	0.648	5.23
26) C	2-Nitrophenol		0.047	0.058	0.086	0.116	0.140	0.153	0.157	0.108	41.87#
27)	2,4-Dimethylph...		0.223	0.220	0.232	0.226	0.224	0.230	0.220	0.225	2.09
28)	bis(2-Chloroet...		0.426	0.405	0.421	0.391	0.378	0.388	0.352	0.394	6.51
29) C	2,4-Dichloroph...		0.251	0.260	0.285	0.280	0.272	0.283	0.267	0.271	4.61
30)	1,2,4-Trichlor...		0.348	0.340	0.351	0.325	0.311	0.323	0.292	0.327	6.49
31)	Naphthalene		1.127	1.063	1.072	0.957	0.883	0.906	0.763	0.967	13.23
32)	Benzoic acid			0.056	0.097	0.137	0.179	0.202	0.216	0.148	42.49
33)	4-Chloroaniline		0.400	0.375	0.390	0.357	0.346	0.351	0.327	0.364	7.11
34) C	Hexachlorobuta...		0.207	0.200	0.205	0.193	0.183	0.189	0.169	0.192	6.86
35)	Caprolactam		0.086	0.085	0.093	0.091	0.089	0.094	0.096	0.091	4.53
36) C	4-Chloro-3-met...		0.275	0.271	0.293	0.281	0.274	0.286	0.265	0.278	3.53
37)	2-Methylnaphth...		0.743	0.708	0.715	0.647	0.606	0.622	0.524	0.652	11.70
38)	1-Methylnaphth...		0.730	0.692	0.701	0.639	0.589	0.605	0.515	0.639	11.77

Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
 Method File : 8270-BF060524.M

39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.599	0.584	0.610	0.568	0.542	0.550	0.499	0.565	6.74	
41) P	Hexachlorocycl...	0.143	0.160	0.189	0.199	0.206	0.212	0.196	0.186	13.72	
42) S	2,4,6-Tribromo...	0.167	0.185	0.210	0.209	0.196	0.204	0.177	0.192	8.69	
43) C	2,4,6-Trichlor...	0.291	0.308	0.355	0.359	0.363	0.367	0.356	0.343	8.89	
44)	2,4,5-Trichlor...	0.327	0.357	0.393	0.394	0.390	0.406	0.378	0.378	7.26	
45) S	2-Fluorobiphenyl	1.444	1.380	1.124	1.033	1.021	0.861	1.144		19.69	
46)	1,1'-Biphenyl	1.603	1.556	1.566	1.389	1.296	1.310	1.114	1.405	12.80	
47)	2-Chloronaphth...	1.245	1.199	1.216	1.115	1.050	1.065	0.948	1.120	9.56	
48)	2-Nitroaniline	0.114	0.150	0.215	0.258	0.280	0.296	0.288	0.229	31.54	
49)	Acenaphthylene	1.769	1.697	1.742	1.576	1.485	1.513	1.309	1.584	10.39	
50)	Dimethylphthalate	1.389	1.329	1.359	1.247	1.167	1.219	1.087	1.257	8.68	
51)	2,6-Dinitrotol...	0.131	0.176	0.240	0.262	0.272	0.290	0.277	0.235	25.12	
52) C	Acenaphthene	1.175	1.145	1.170	1.057	1.016	1.023	0.907	1.070	9.23	
53)	3-Nitroaniline	0.149	0.185	0.247	0.263	0.272	0.291	0.274	0.240	21.94	
54) P	2,4-Dinitrophenol	0.019	0.025	0.038	0.058	0.080	0.096	0.102	0.060	56.81	
55)	Dibenzofuran	1.726	1.663	1.664	1.508	1.394	1.422	1.221	1.514	12.02	
56) P	4-Nitrophenol	0.125	0.180	0.202	0.206	0.228	0.205	0.191		18.61	
57)	2,4-Dinitrotol...	0.133	0.179	0.268	0.315	0.329	0.360	0.321	0.272	31.21	
58)	Fluorene	1.422	1.371	1.377	1.205	1.068	1.089	0.888	1.203	16.53	
59)	2,3,4,6-Tetrac...	0.252	0.273	0.310	0.320	0.313	0.324	0.295	0.298	8.94	
60)	Diethylphthalate	1.363	1.324	1.348	1.238	1.135	1.173	1.006	1.227	10.67	
61)	4-Chlorophenyl...	0.713	0.686	0.689	0.607	0.556	0.557	0.470	0.611	14.64	
62)	4-Nitroaniline	0.156	0.194	0.248	0.264	0.255	0.279	0.246	0.235	18.55	
63)	Azobenzene	1.332	1.285	1.298	1.183	1.085	1.126	0.963	1.182	11.30	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.024	0.040	0.061	0.077	0.087	0.092	0.063		42.84	
66) c	n-Nitrosodiphe...	0.653	0.626	0.629	0.595	0.575	0.569	0.533	0.597	6.96	
67)	4-Bromophenyl-...	0.237	0.231	0.236	0.228	0.225	0.223	0.214	0.228	3.52	
68)	Hexachlorobenzene	0.285	0.273	0.276	0.261	0.258	0.258	0.243	0.265	5.31	
69)	Atrazine	0.207	0.205	0.202	0.191	0.184	0.183	0.173	0.192	6.80	
70) C	Pentachlorophenol	0.094	0.113	0.142	0.156	0.161	0.167	0.160	0.142	19.69	
71)	Phenanthrene	1.114	1.071	1.059	0.956	0.889	0.867	0.783	0.963	12.77	
72)	Anthracene	1.095	1.058	1.060	0.945	0.887	0.897	0.779	0.960	12.03	
73)	Carbazole	0.987	0.963	0.966	0.869	0.817	0.820	0.725	0.878	11.12	
74)	Di-n-butylphth...	1.165	1.158	1.167	1.035	0.935	0.946	0.819	1.032	13.35	
75) C	Fluoranthene	1.206	1.194	1.190	1.032	0.934	0.933	0.782	1.039	15.87	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.089	0.136	0.265	0.278	0.241	0.353	0.079	0.206	50.99	
78)	Pyrene	1.761	1.691	1.828	1.830	1.859	1.943	1.750	1.809	4.57	
79) S	Terphenyl-d14	1.464	1.413	1.439	1.296	1.266	1.323	1.159	1.337	8.14	
80)	Butylbenzylpht...	0.416	0.482	0.602	0.652	0.666	0.707	0.655	0.597	17.96	
81)	Benzo(a)anthra...	1.369	1.332	1.366	1.240	1.190	1.228	1.107	1.262	7.80	
82)	3,3'-Dichlorob...	0.303	0.312	0.339	0.322	0.315	0.346	0.302	0.320	5.34	
83)	Chrysene	1.252	1.251	1.273	1.223	1.142	1.208	1.128	1.211	4.63	
84)	Bis(2-ethylhex...	0.790	0.843	0.932	0.916	0.889	0.900	0.822	0.870	6.07	
85) c	Di-n-octyl pht...	1.017	1.125	1.257	1.235	1.250	1.274	1.227	1.198	7.80	

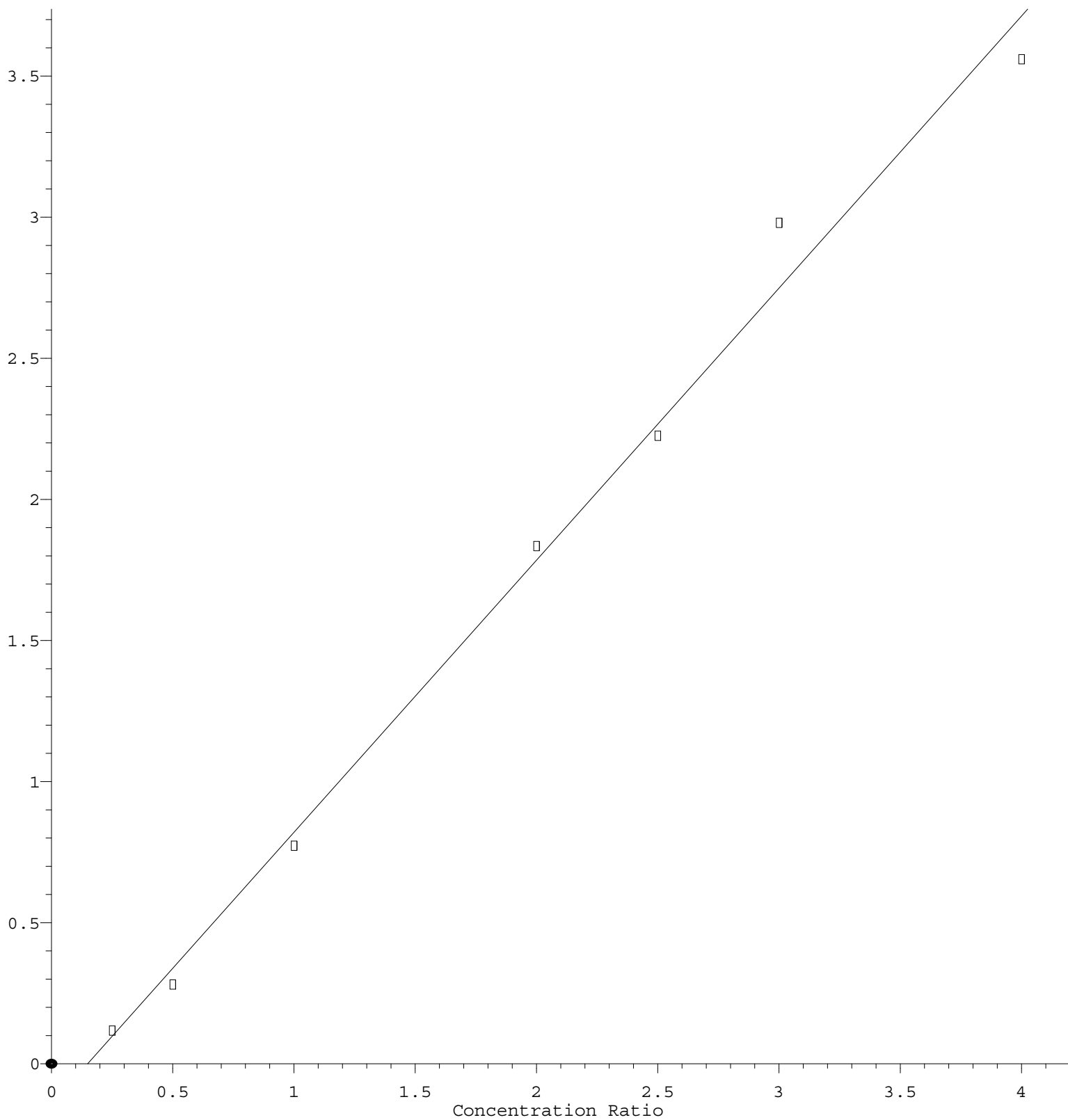
Method Path : Z:\svoasrv\HPCHEM1\BNA_F\Methods\
Method File : 8270-BF060524.M

		-----ISTD-----	
86) I	Perylene-d12		
87)	Indeno(1,2,3-c...	0.579 0.679 0.909 1.076 1.058 1.185 1.095 0.940	24.40
88)	Benzo(b)fluora...	1.331 1.357 1.375 1.214 1.167 1.226 1.131 1.257	7.71
89)	Benzo(k)fluora...	1.255 1.285 1.344 1.190 1.165 1.092 1.063 1.199	8.53
90) C	Benzo(a)pyrene	0.930 0.930 0.976 0.972 0.940 0.982 0.936 0.952	2.44
91)	Dibenzo(a,h)an...	0.455 0.531 0.723 0.860 0.842 0.944 0.881 0.748	25.08
92)	Benzo(g,h,i)pe...	0.471 0.562 0.773 0.917 0.890 0.993 0.890 0.785	25.03

(#) = Out of Range

Benzo(g,h,i)perylene

Response Ratio



$$\text{Response} = 9.643\text{e-}001 * \text{Amt} - 1.439\text{e-}001$$

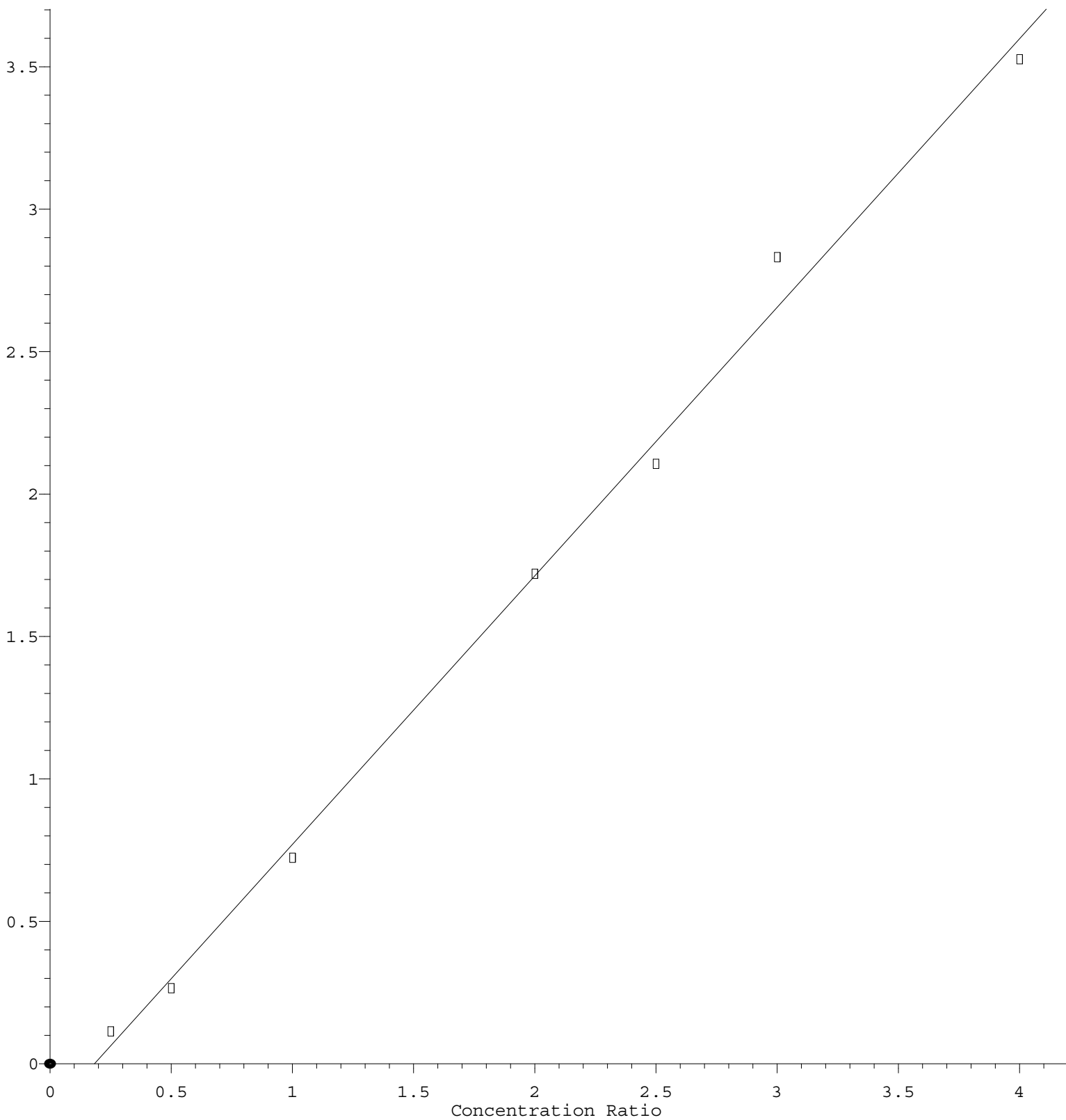
Coef of Det (r^2) = 0.991918 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

Dibenzo (a,h) anthracene

Response Ratio



$$\text{Response} = 9.428\text{e-}001 * \text{Amt} - 1.723\text{e-}001$$

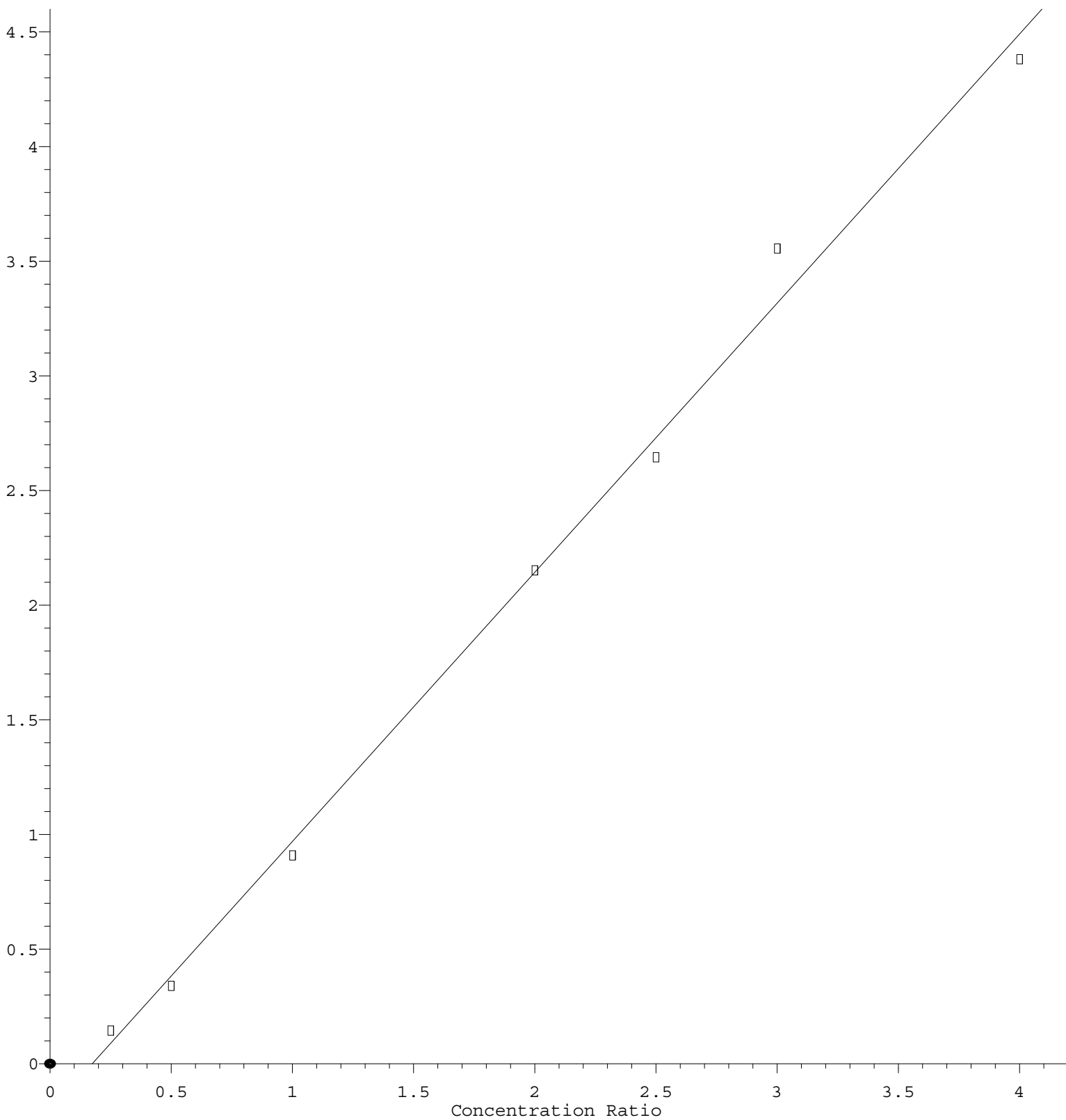
Coef of Det (r^2) = 0.995292 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

Indeno(1,2,3-cd)pyrene

Response Ratio



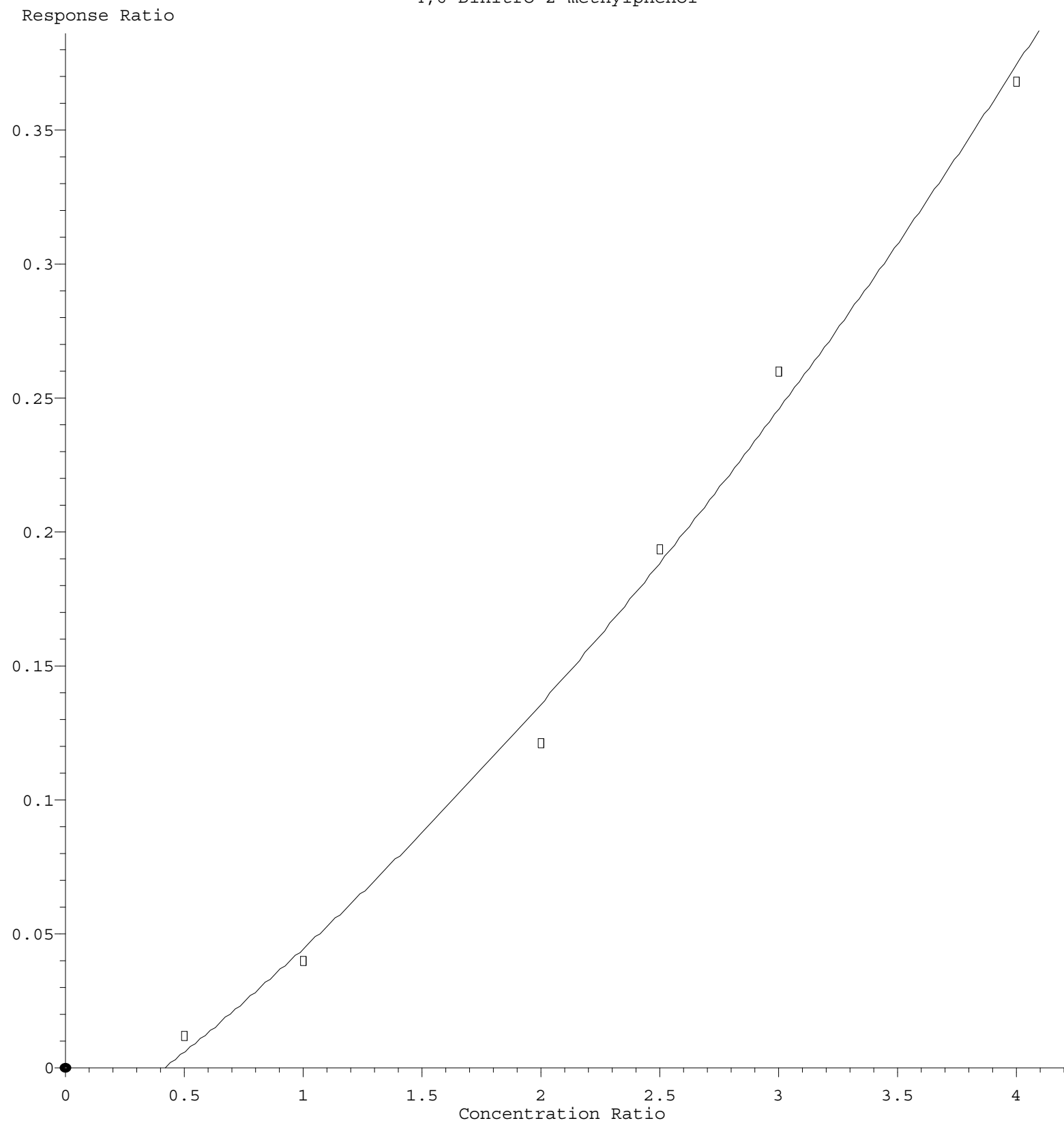
$$\text{Response} = 1.174\text{e}+000 * \text{Amt} - 2.044\text{e}-001$$

Coef of Det (r^2) = 0.994674 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

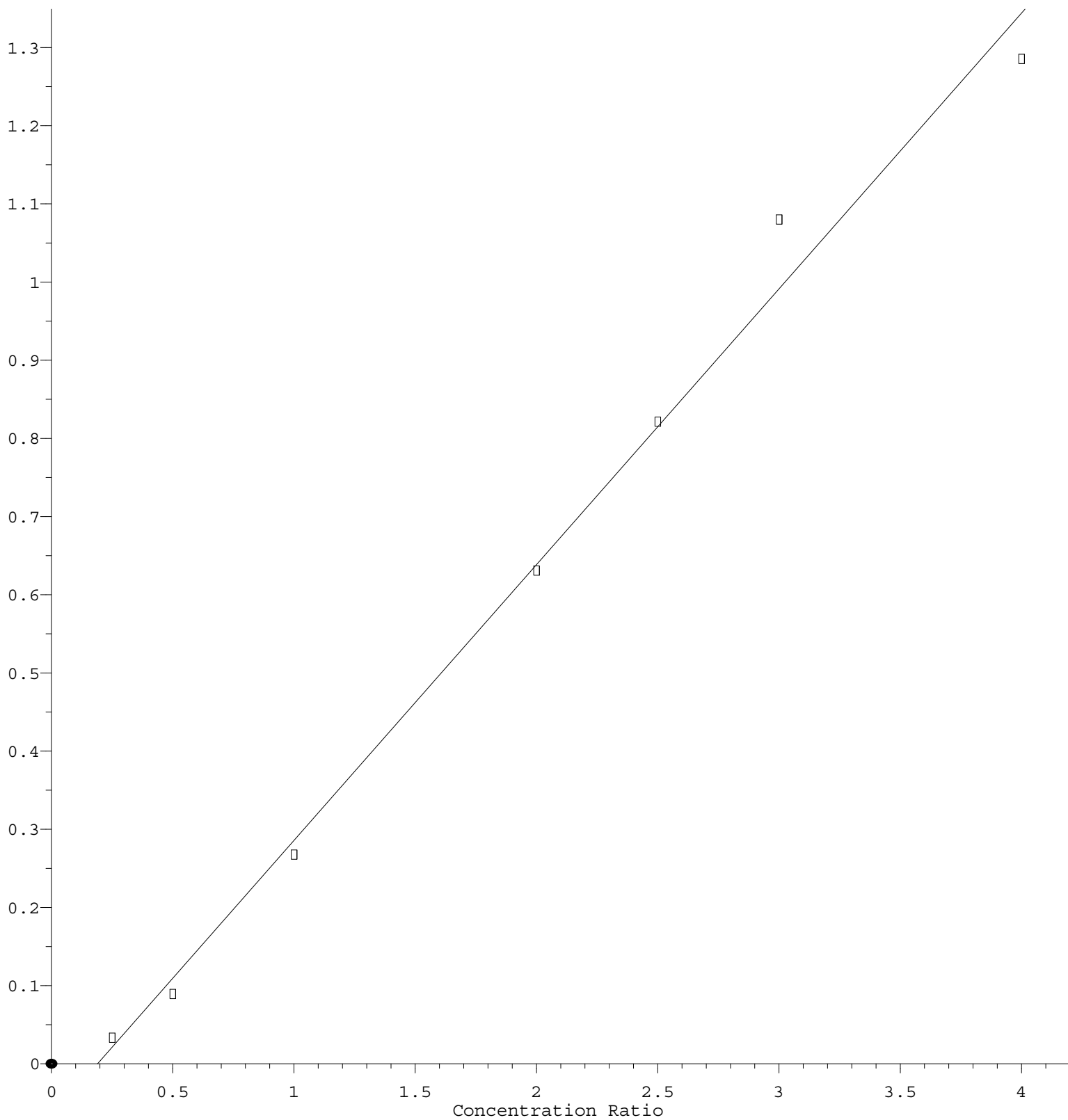
4,6-Dinitro-2-methylphenol



$R = 9.302e-003 A^2 + 6.338e-002 A - 2.820e-002$
 Coef of Det (r^2) = 0.994183 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M
 Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

2,4-Dinitrotoluene

Response Ratio



$$\text{Response} = 3.529\text{e-}001 * \text{Amt} - 6.695\text{e-}002$$

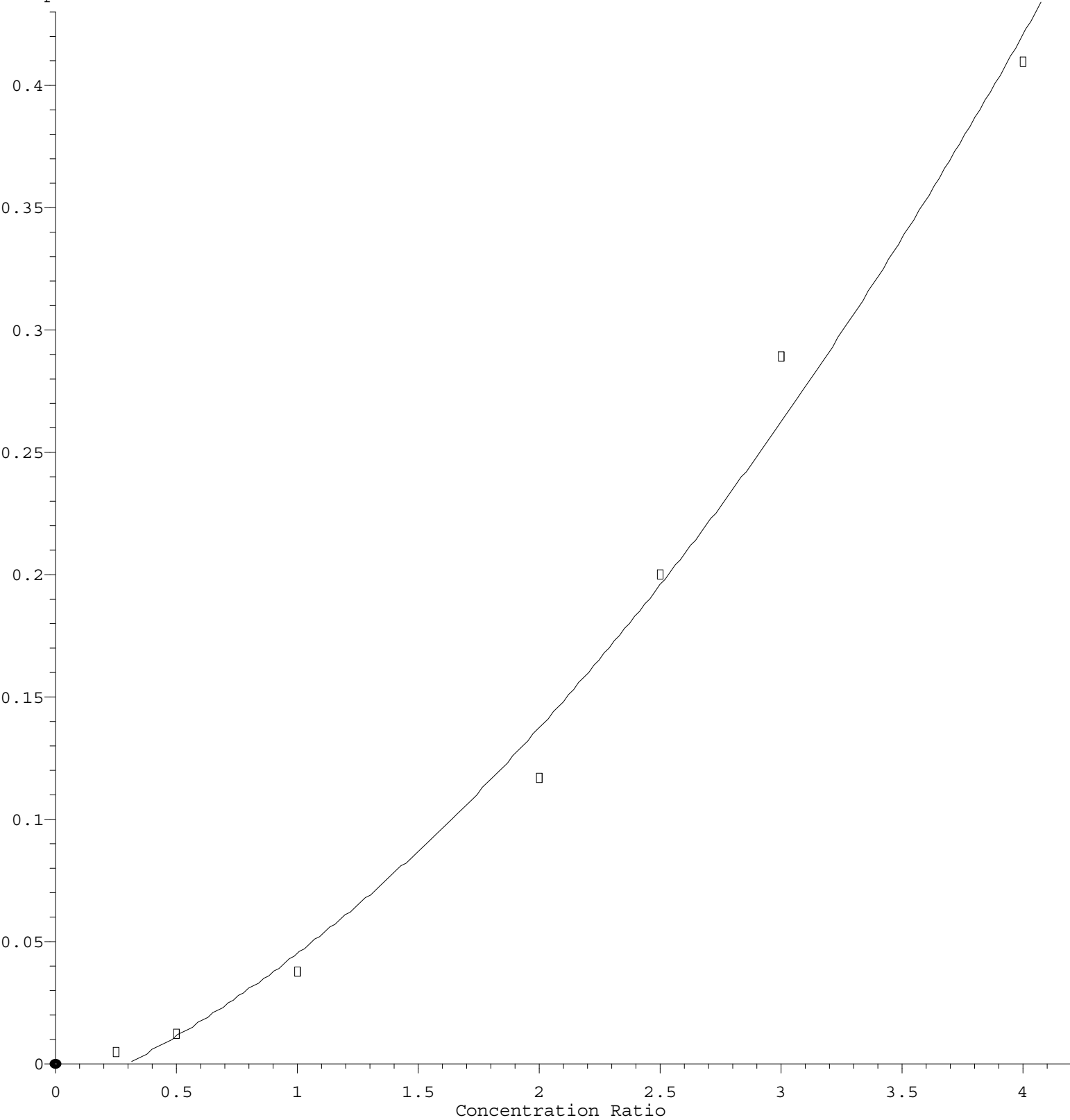
Coef of Det (r^2) = 0.991485 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

2,4-Dinitrophenol

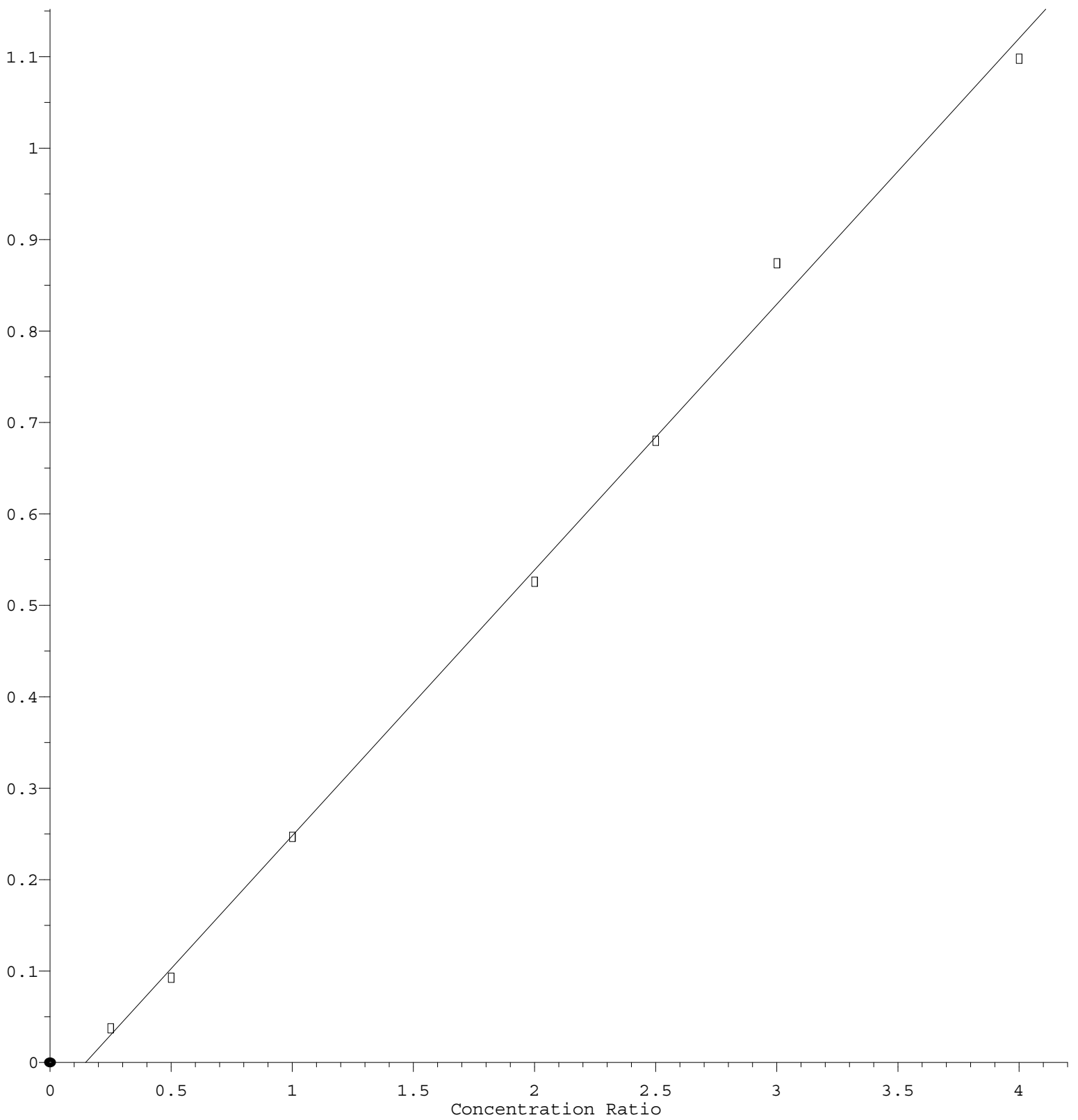
Response Ratio



$R = 1.650e-002 A^2 + 4.272e-002 A - 1.409e-002$
 Coef of Det (r^2) = 0.990387 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M
 Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

3-Nitroaniline

Response Ratio



$$\text{Response} = 2.908\text{e-}001 * \text{Amt} - 4.282\text{e-}002$$

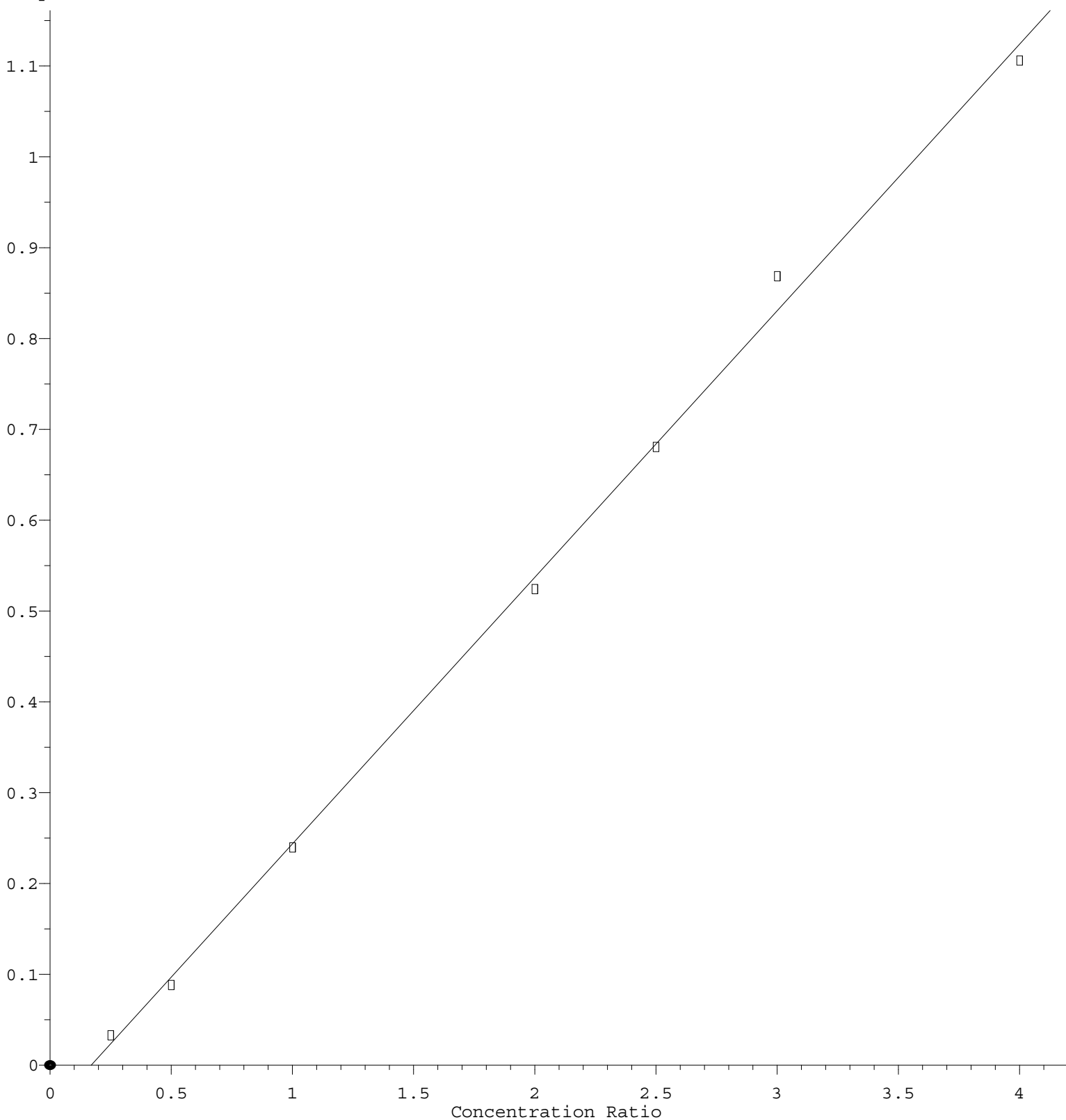
Coef of Det (r^2) = 0.997090 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

2,6-Dinitrotoluene

Response Ratio



$$\text{Response} = 2.936\text{e-}001 * \text{Amt} - 4.994\text{e-}002$$

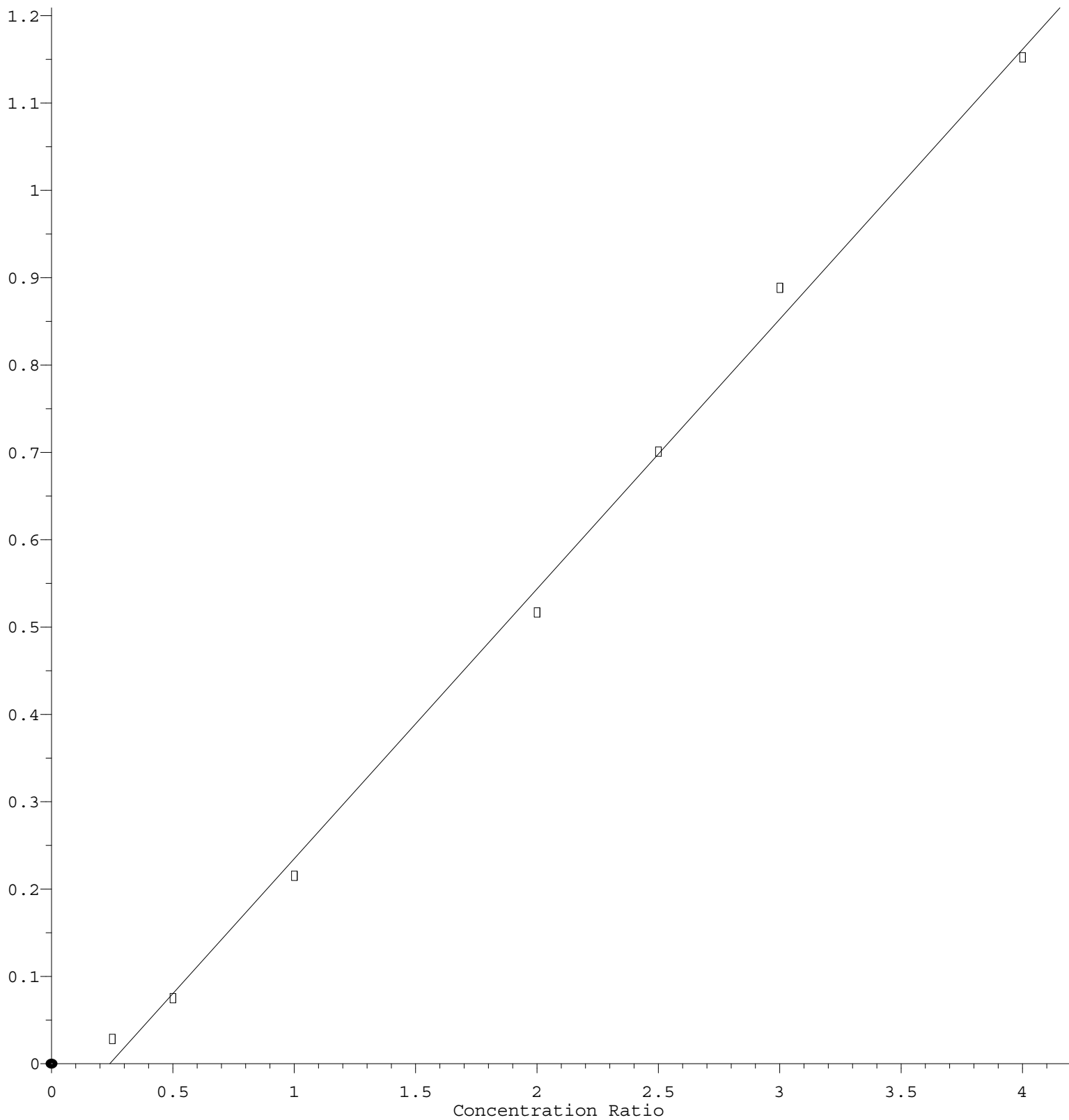
Coef of Det (r^2) = 0.997857 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

2-Nitroaniline

Response Ratio



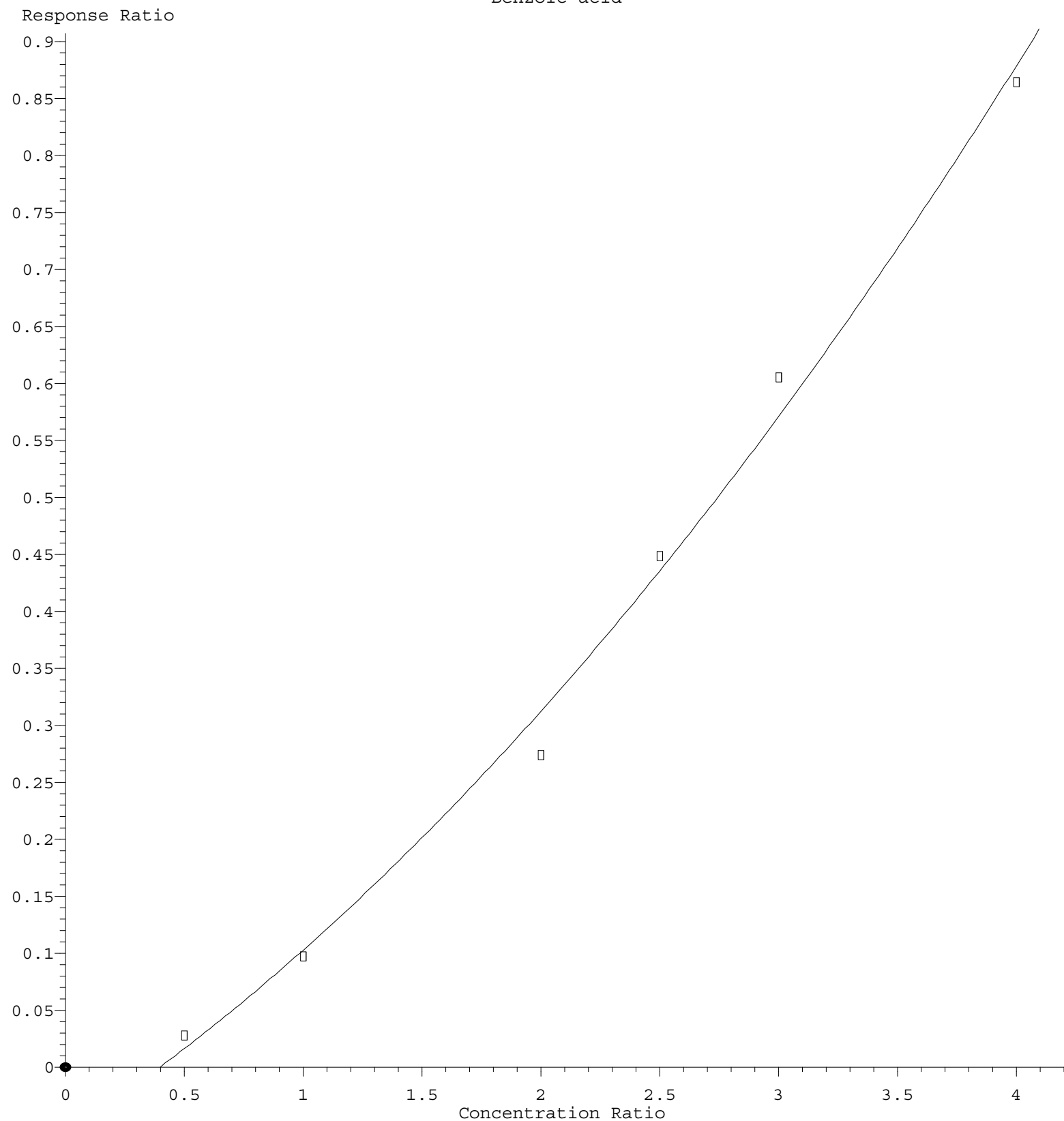
$$\text{Response} = 3.091\text{e-}001 * \text{Amt} - 7.413\text{e-}002$$

Coef of Det (r^2) = 0.997129 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

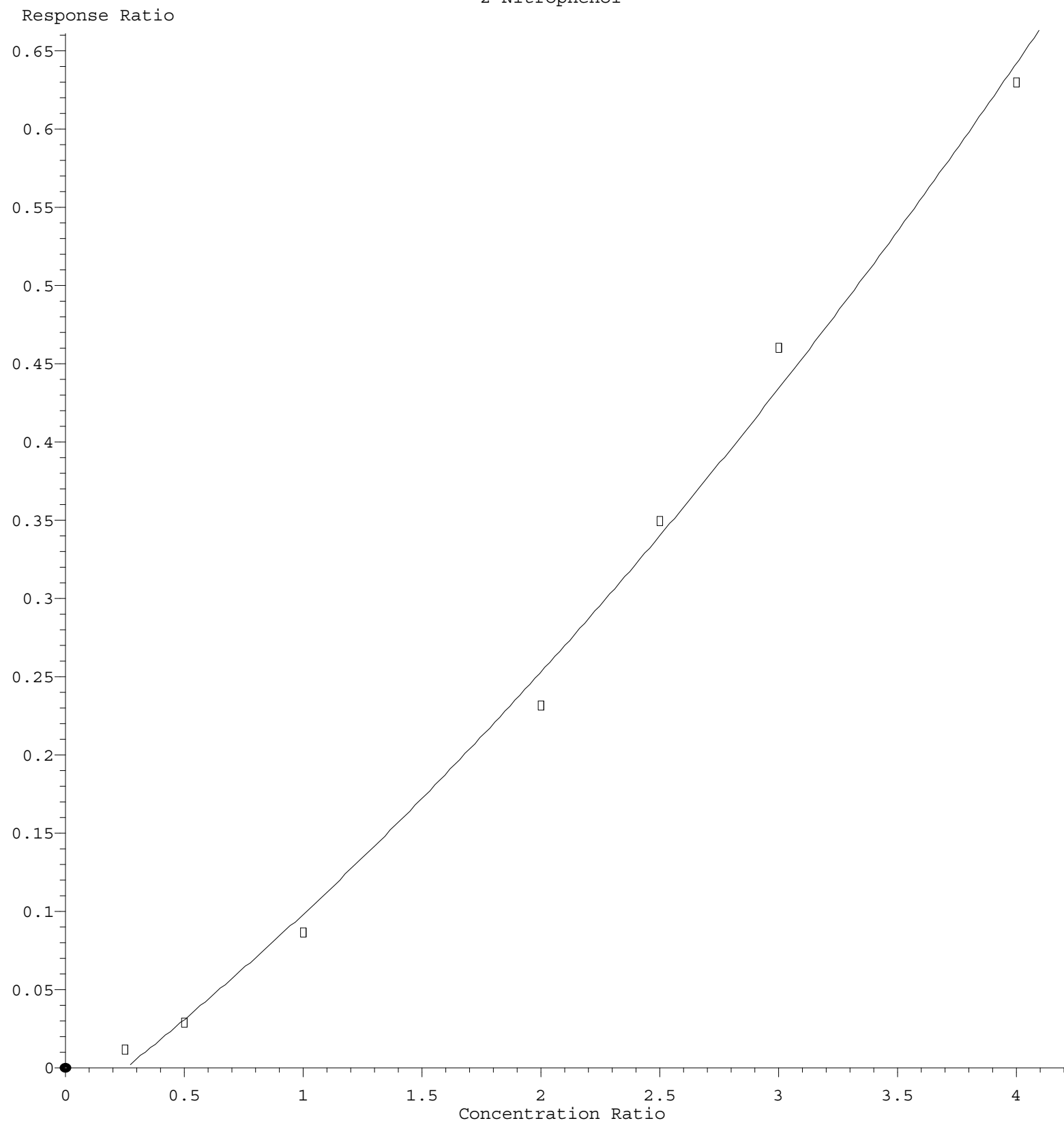
Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

Benzoic acid



$R = 2.450e-002 A^2 + 1.361e-001 A - 5.786e-002$
 Coef of Det (r^2) = 0.993688 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M
 Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

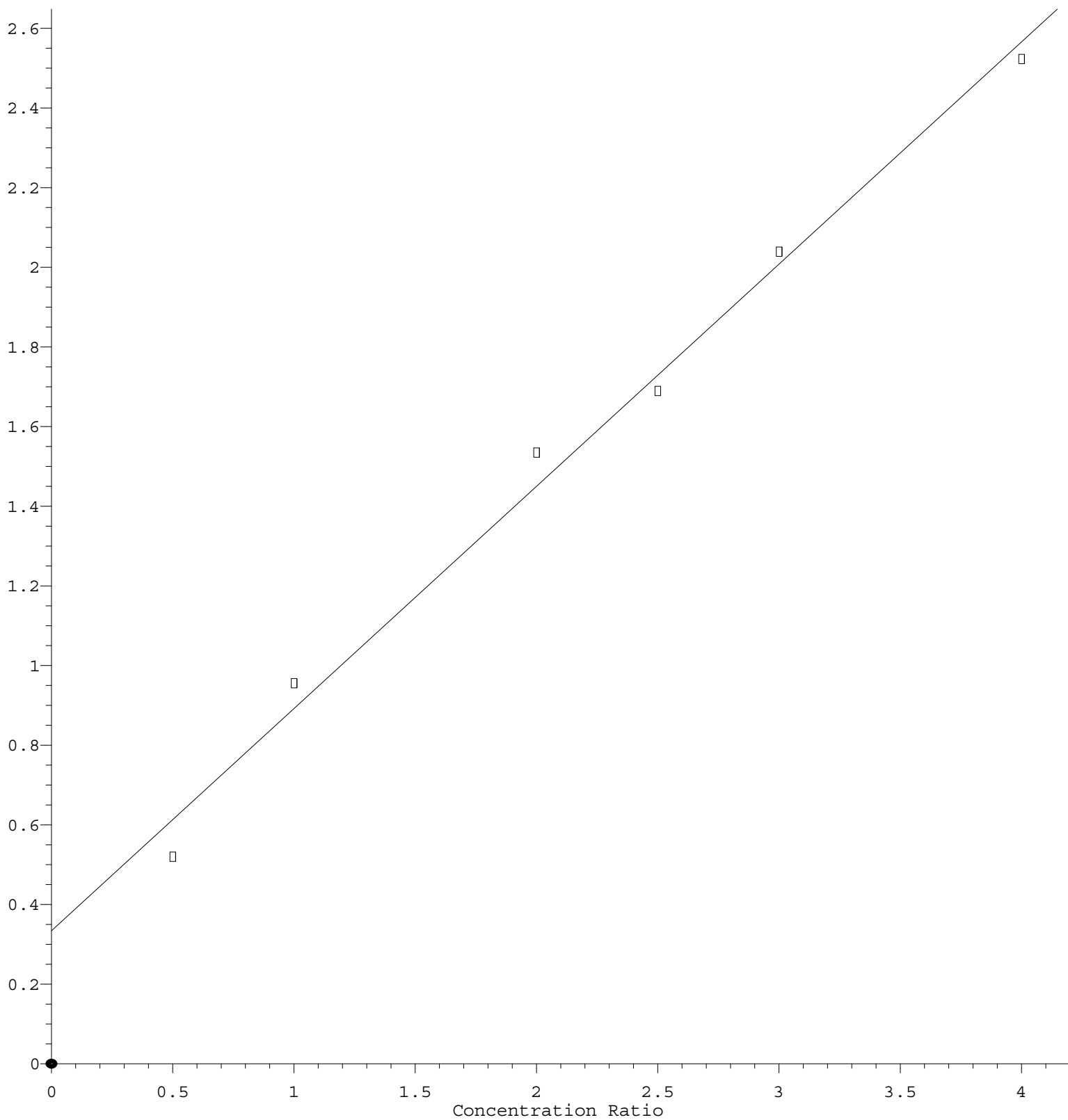
2-Nitrophenol



$R = 1.329e-002 A^2 + 1.148e-001 A - 2.983e-002$
 Coef of Det (r^2) = 0.995000 Curve Fit: Quadratic
 Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M
 Calibration Table Last Updated: Thu Jun 06 05:19:29 2024

Benzaldehyde

Response Ratio



$$\text{Response} = 5.580\text{e-}001 * \text{Amt} + 3.344\text{e-}001$$

Coef of Det (r^2) = 0.990718 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA F\Methods\8270-BF060524.M

Calibration Table Last Updated: Thu Jun 06 05:19:29 2024