

Method Path : Z:\svoasrv\HPCHEM1\BNA_M\Methods\
 Method File : 8270-BM062724.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Fri Jun 28 03:23:08 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BM046334.D 5 =BM046335.D 10 =BM046336.D 20 =BM046337.D 40 =BM046338.D 50 =BM046339.D 60 =BM046340.D 80 =BM046341.D

	Compound	2.5	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.543	0.525	0.462	0.481	0.461	0.450	0.437	0.480		8.29
3)	Pyridine	1.190	1.322	1.420	1.433	1.360	1.329	1.330	1.341		5.97
4)	n-Nitrosodimet...	1.009	1.019	1.095	1.098	1.027	1.029	0.990	1.038		4.04
5) S	2-Fluorophenol	1.202	1.182	1.258	1.314	1.243	1.283	1.227	1.244		3.66
6)	Aniline	1.506	1.573	1.710	1.761	1.601	1.553	1.532	1.605		5.92
7) S	Phenol-d6	1.599	1.679	1.850	1.905	1.776	1.815	1.750	1.768		5.85
8)	2-Chlorophenol	1.322	1.344	1.417	1.470	1.372	1.413	1.367	1.386		3.62
9)	Benzaldehyde	1.327	1.287	1.196	1.052	0.898	0.813		1.096		19.19
10) C	Phenol	1.708	1.639	1.800	1.864	1.729	1.781	1.723	1.749		4.16
11)	bis(2-Chloroet...	1.415	1.350	1.427	1.518	1.409	1.421	1.403	1.420		3.52
12)	1,3-Dichlorobe...	1.485	1.518	1.573	1.606	1.508	1.563	1.506	1.537		2.86
13) C	1,4-Dichlorobe...	1.622	1.552	1.597	1.636	1.550	1.593	1.521	1.581		2.65
14)	1,2-Dichlorobe...	1.544	1.554	1.589	1.641	1.540	1.554	1.492	1.559		2.96
15)	Benzyl Alcohol	1.256	1.320	1.513	1.615	1.458	1.539	1.471	1.453		8.62
16)	2,2'-oxybis(1-...	2.426	2.509	2.529	2.557	2.324	2.354	2.221	2.417		5.11
17)	2-Methylphenol	1.153	1.138	1.262	1.331	1.212	1.245	1.198	1.220		5.44
18)	Hexachloroethane	0.716	0.710	0.723	0.737	0.693	0.720	0.687	0.712		2.42
19) P	n-Nitroso-di-n...	1.029	1.206	1.305	1.476	1.516	1.361	1.459	1.383	1.342	12.02
20)	3+4-Methylphenols	1.543	1.614	1.848	1.946	1.760	1.874	1.797	1.769		8.14
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.492	0.525	0.565	0.590	0.560	0.604	0.575	0.559		6.91
23) S	Nitrobenzene-d5	0.454	0.467	0.489	0.508	0.484	0.494	0.481	0.482		3.66
24)	Nitrobenzene	0.465	0.454	0.482	0.498	0.474	0.479	0.466	0.474		3.00
25)	Isophorone	0.792	0.806	0.846	0.868	0.801	0.822	0.791	0.818		3.57
26) C	2-Nitrophenol	0.128	0.143	0.172	0.188	0.179	0.193	0.185	0.170		14.65
27)	2,4-Dimethylph...	0.186	0.205	0.219	0.228	0.213	0.223	0.214	0.213		6.49
28)	bis(2-Chloroet...	0.416	0.429	0.464	0.479	0.448	0.474	0.454	0.452		5.10
29) C	2,4-Dichloroph...	0.234	0.262	0.296	0.310	0.294	0.312	0.301	0.287		9.93
30)	1,2,4-Trichlor...	0.312	0.310	0.326	0.337	0.320	0.342	0.329	0.325		3.69
31)	Naphthalene	1.076	1.058	1.105	1.136	1.070	1.132	1.097	1.096		2.75
32)	Benzoic acid		0.158	0.244	0.268	0.252	0.263	0.260	0.241		17.27
33)	4-Chloroaniline	0.314	0.358	0.401	0.418	0.396	0.402	0.399	0.384		9.36
34) C	Hexachlorobuta...	0.195	0.190	0.202	0.205	0.196	0.214	0.202	0.200		3.85
35)	Caprolactam	0.088	0.101	0.114	0.115	0.109	0.102	0.108	0.105		8.92
36) C	4-Chloro-3-met...	0.316	0.338	0.370	0.388	0.358	0.367	0.365	0.357		6.60
37)	2-Methylnaphth...	0.699	0.721	0.747	0.777	0.729	0.764	0.738	0.739		3.56
38)	1-Methylnaphth...	0.710	0.700	0.740	0.764	0.710	0.759	0.733	0.731		3.44

Method Path : Z:\svoasrv\HPCHEM1\BNA_M\Methods\
 Method File : 8270-BM062724.M

39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.514	0.495	0.532	0.536	0.534	0.581	0.542	0.534	4.92	
41) P	Hexachlorocycl...	0.170	0.205	0.209	0.209	0.250	0.211	0.209	12.28		
42) S	2,4,6-Tribromo...	0.206	0.230	0.259	0.264	0.252	0.242	0.249	0.243	8.07	
43) C	2,4,6-Trichlor...	0.281	0.317	0.364	0.371	0.366	0.401	0.380	0.354	11.61	
44)	2,4,5-Trichlor...	0.395	0.382	0.402	0.424	0.398	0.438	0.419	0.408	4.75	
45) S	2-Fluorobiphenyl	1.365	1.355	1.461	1.518	1.460	1.636	1.535	1.476	6.68	
46)	1,1'-Biphenyl	1.426	1.415	1.487	1.543	1.478	1.611	1.521	1.497	4.55	
47)	2-Chloronaphth...	1.086	1.082	1.174	1.205	1.159	1.243	1.187	1.162	5.11	
48)	2-Nitroaniline	0.382	0.417	0.481	0.502	0.480	0.484	0.484	0.461	9.54	
49)	Acenaphthylene	1.650	1.681	1.797	1.863	1.801	1.902	1.848	1.792	5.23	
50)	Dimethylphthalate	1.365	1.400	1.484	1.532	1.464	1.497	1.492	1.462	4.02	
51)	2,6-Dinitrotol...	0.288	0.309	0.334	0.345	0.332	0.336	0.335	0.326	6.06	
52) C	Acenaphthene	1.053	1.044	1.120	1.153	1.110	1.174	1.143	1.114	4.45	
53)	3-Nitroaniline	0.266	0.296	0.342	0.359	0.346	0.334	0.359	0.329	10.66	
54) P	2,4-Dinitrophenol	0.106	0.164	0.184	0.178	0.185	0.196	0.169	19.36		
55)	Dibenzofuran	1.681	1.681	1.797	1.857	1.778	1.866	1.833	1.785	4.32	
56) P	4-Nitrophenol	0.151	0.251	0.275	0.265	0.247	0.280	0.245	19.55		
57)	2,4-Dinitrotol...	0.395	0.419	0.489	0.514	0.491	0.481	0.510	0.471	9.73	
58)	Fluorene	1.391	1.407	1.533	1.608	1.535	1.609	1.602	1.527	6.09	
59)	2,3,4,6-Tetrac...	0.332	0.334	0.353	0.361	0.343	0.344	0.343	0.344	2.95	
60)	Diethylphthalate	1.511	1.564	1.636	1.693	1.613	1.601	1.640	1.608	3.64	
61)	4-Chlorophenyl...	0.645	0.649	0.702	0.744	0.709	0.773	0.753	0.711	7.03	
62)	4-Nitroaniline	0.247	0.286	0.344	0.368	0.350	0.307	0.357	0.323	13.71	
63)	Azobenzene	1.837	1.899	2.006	2.018	1.903	1.878	1.875	1.917	3.58	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.085	0.116	0.127	0.120	0.132	0.133	0.119	15.05		
66) c	n-Nitrosodiphe...	0.523	0.541	0.575	0.600	0.570	0.663	0.594	0.581	7.86	
67)	4-Bromophenyl-...	0.188	0.189	0.205	0.210	0.199	0.232	0.207	0.204	7.29	
68)	Hexachlorobenzene	0.230	0.226	0.243	0.246	0.236	0.265	0.242	0.241	5.33	
69)	Atrazine	0.173	0.176	0.188	0.163	0.152	0.141	0.132	0.161	12.53	
70) C	Pentachlorophenol	0.116	0.131	0.159	0.168	0.157	0.166	0.164	0.152	13.07	
71)	Phenanthrene	1.014	1.004	1.066	1.106	1.055	1.113	1.092	1.064	4.04	
72)	Anthracene	0.952	0.975	1.045	1.085	1.032	1.085	1.073	1.035	5.14	
73)	Carbazole	0.942	0.957	1.045	1.085	1.040	0.995	1.072	1.019	5.48	
74)	Di-n-butylphth...	1.161	1.246	1.410	1.488	1.393	1.405	1.483	1.370	8.92	
75) C	Fluoranthene	1.177	1.217	1.296	1.353	1.285	1.140	1.335	1.258	6.45	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.193	0.119	0.157	0.270	0.222	0.284	0.208	31.01		
78)	Pyrene	1.238	1.238	1.379	1.406	1.333	1.707	1.423	1.389	11.44	
79) S	Terphenyl-d14	1.006	1.043	1.198	1.262	1.184	1.510	1.259	1.209	13.73	
80)	Butylbenzylpht...	0.522	0.555	0.663	0.701	0.658	0.756	0.720	0.654	13.15	
81)	Benzo(a)anthra...	1.244	1.251	1.320	1.335	1.306	1.364	1.363	1.312	3.72	
82)	3,3'-Dichlorob...	0.386	0.388	0.446	0.465	0.437	0.412	0.412	0.421	7.09	
83)	Chrysene	1.201	1.202	1.266	1.297	1.247	1.279	1.291	1.255	3.18	
84)	Bis(2-ethylhex...	0.767	0.847	1.039	1.130	1.046	1.154	1.148	1.019	15.06	
85) c	Di-n-octyl pht...	1.355	1.478	1.692	1.818	1.723	1.675	1.826	1.652	10.59	

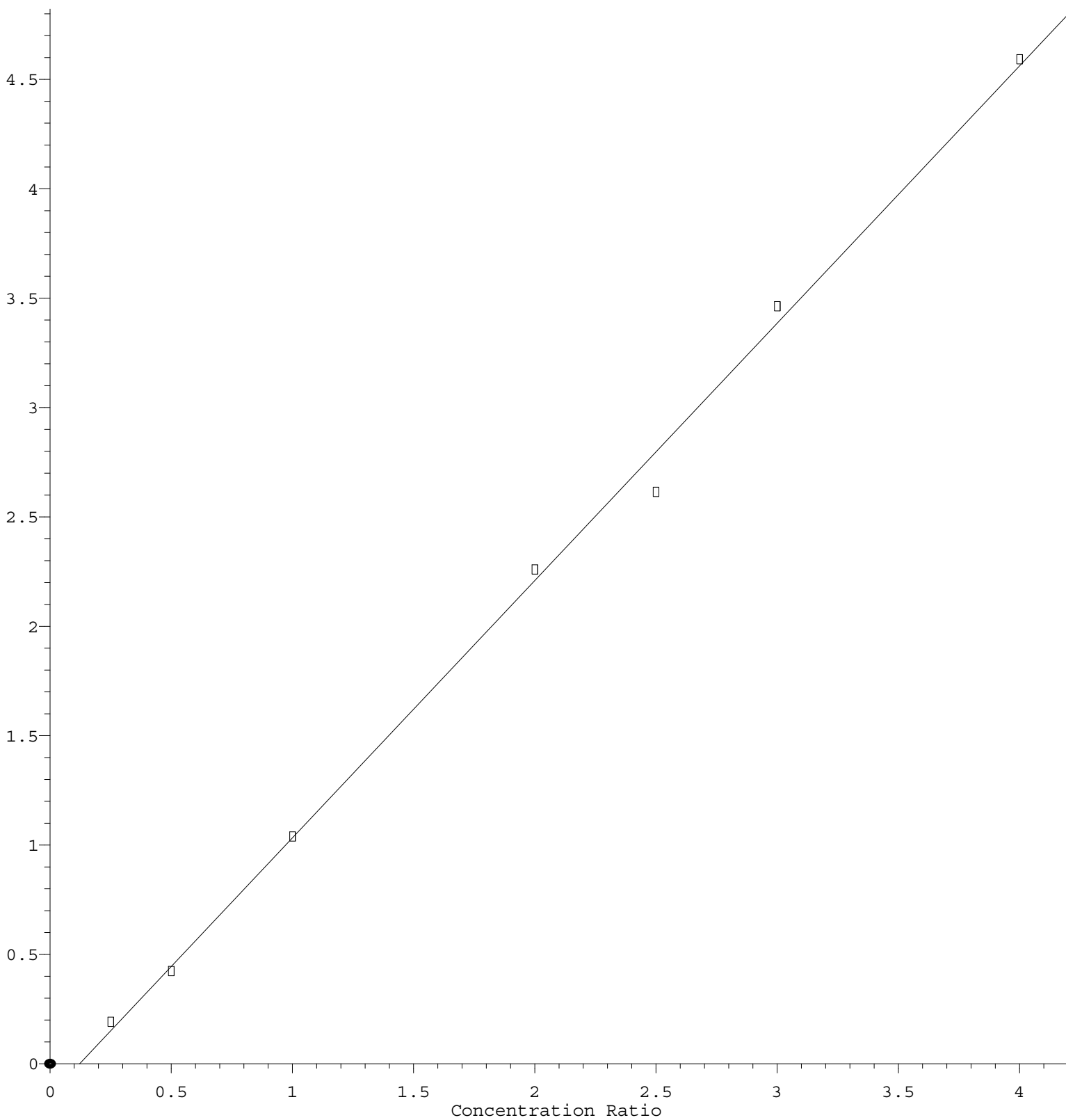
Method Path : Z:\svoasrv\HPCHEM1\BNA_M\Methods\
Method File : 8270-BM062724.M

86) I	Perylene-d12	-----ISTD-----	
87)	Indeno(1,2,3-c...	1.267 1.279 1.257 1.219 1.332 1.390 1.235 1.283	4.62
88)	Benzo(b)fluora...	1.020 1.061 1.230 1.277 1.167 1.280 1.257 1.185	8.97
89)	Benzo(k)fluora...	1.105 1.107 1.192 1.248 1.145 1.224 1.228 1.178	5.04
90) C	Benzo(a)pyrene	0.941 0.952 1.049 1.101 1.041 1.104 1.095 1.040	6.60
91)	Dibenzo(a,h)an...	1.048 1.043 1.040 1.022 1.123 1.160 1.063 1.071	4.72
92)	Benzo(g,h,i)pe...	1.078 1.060 1.024 0.979 1.118 1.171 0.996 1.061	6.44

(#) = Out of Range

Bis(2-ethylhexyl)phthalate

Response Ratio



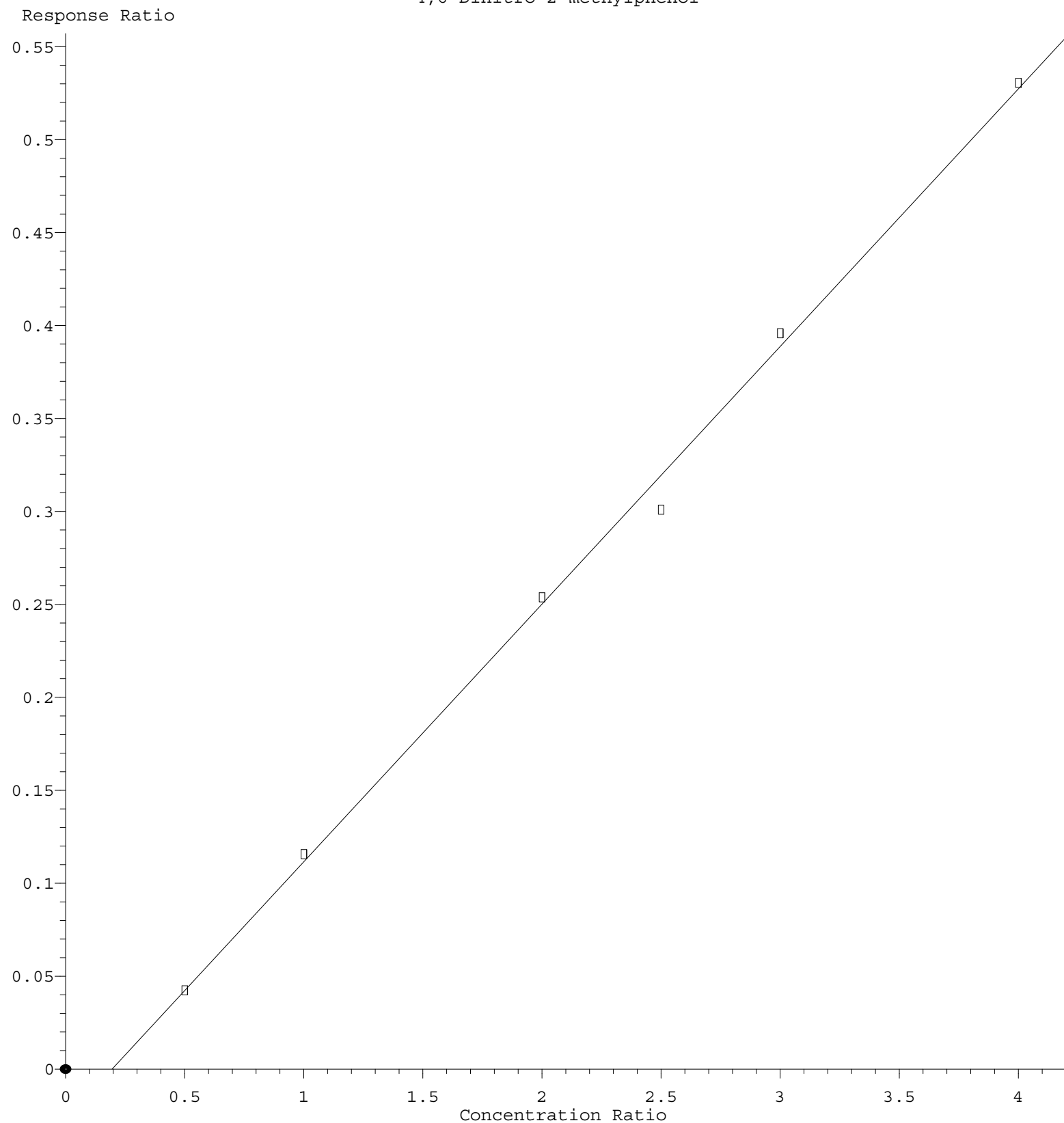
$$\text{Response} = 1.176 \times 10^0 \times \text{Amt} - 1.434 \times 10^{-1}$$

Coef of Det (r^2) = 0.997170 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M

Calibration Table Last Updated: Fri Jun 28 03:23:08 2024

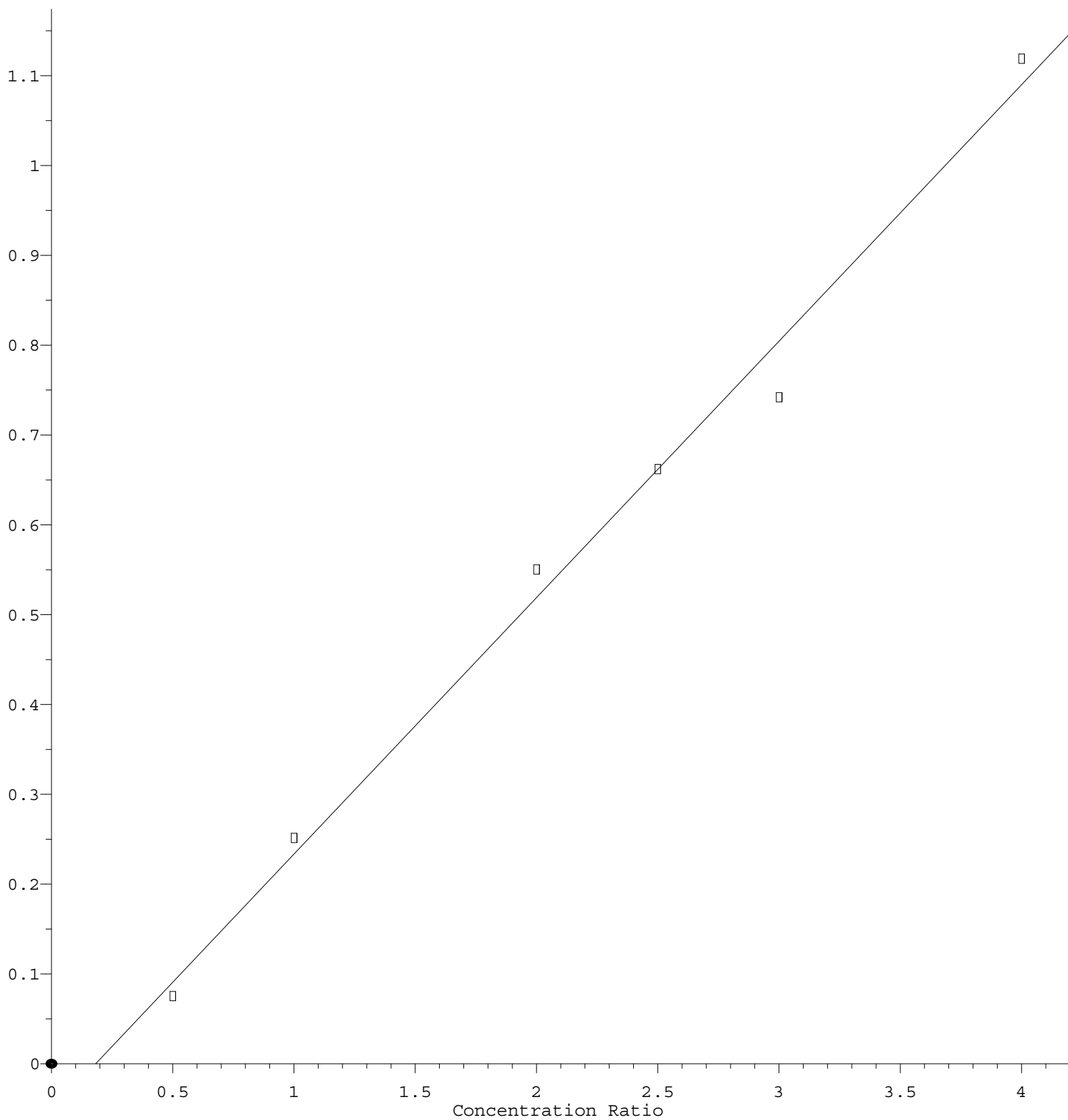
4,6-Dinitro-2-methylphenol



Response = $1.386 \times 10^{-1} \times \text{Amt} - 2.707 \times 10^{-2}$
Coef of Det (r^2) = 0.997300 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M
Calibration Table Last Updated: Fri Jun 28 03:23:08 2024

4-Nitrophenol

Response Ratio



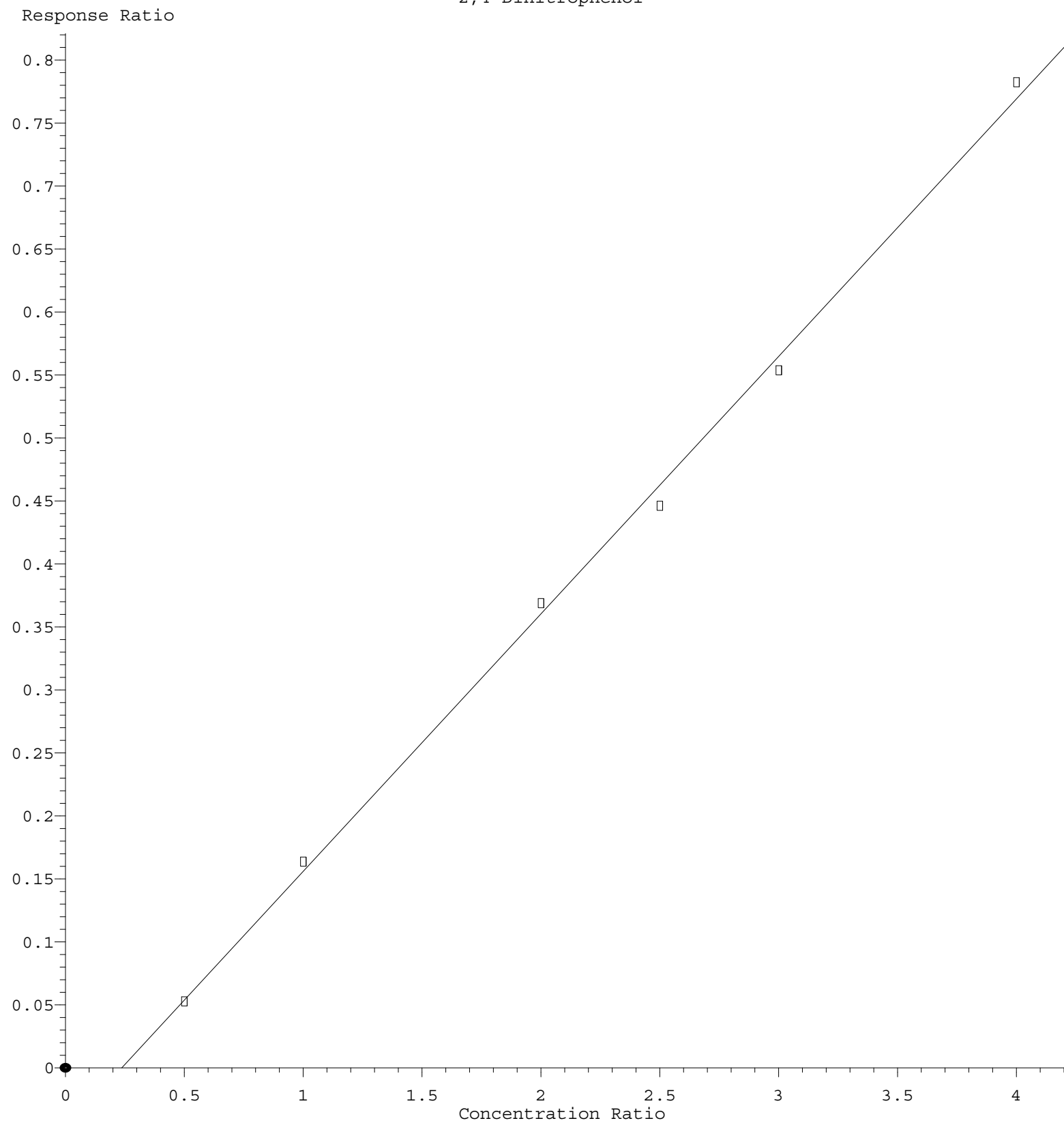
$$\text{Response} = 2.856\text{e-}001 * \text{Amt} - 5.206\text{e-}002$$

Coef of Det (r^2) = 0.990838 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M

Calibration Table Last Updated: Fri Jun 28 03:23:08 2024

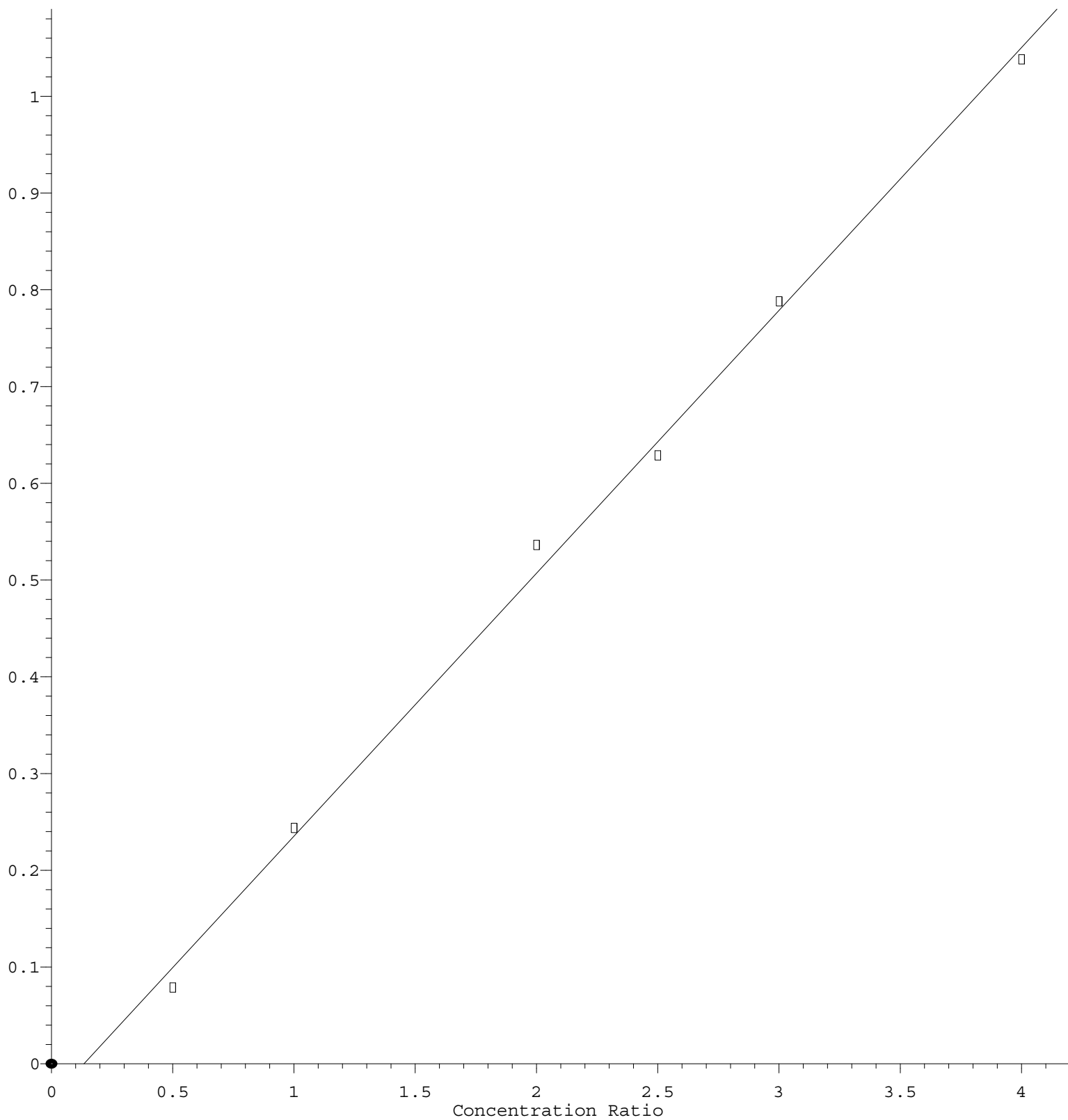
2,4-Dinitrophenol



Response = 2.045e-001 * Amt - 4.846e-002
 Coef of Det (r^2) = 0.997999 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M
 Calibration Table Last Updated: Fri Jun 28 03:23:08 2024

Benzoic acid

Response Ratio



$$\text{Response} = 2.717\text{e-}001 * \text{Amt} - 3.648\text{e-}002$$

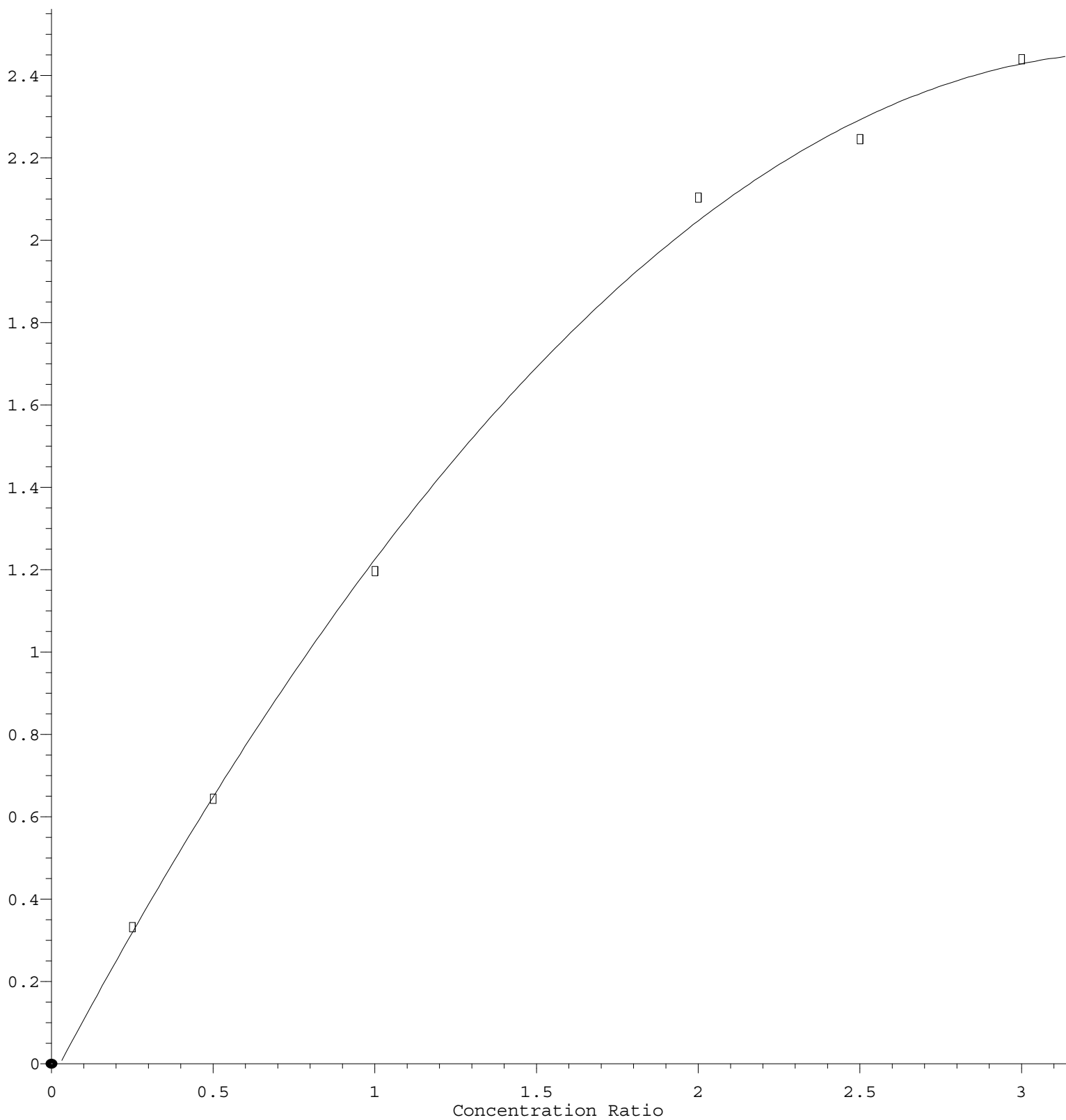
Coef of Det (r^2) = 0.997107 Curve Fit: Linear

Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M

Calibration Table Last Updated: Fri Jun 28 03:23:08 2024

Benzaldehyde

Response Ratio



$R = -2.205e-001 A^2 + 1.484e+000 A - 3.843e-002$
Coef of Det (r^2) = 0.998357 Curve Fit: Quadratic
Method Name: Z:\svoasrv\HPCHEM1\BNA M\Methods\8270-BM062724.M
Calibration Table Last Updated: Fri Jun 28 03:23:08 2024