

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG043024.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Wed May 01 05:11:25 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BG061058.D 5 =BG061059.D 10 =BG061060.D 20 =BG061061.D 40 =BG061062.D 50 =BG061063.D 60 =BG061064.D 80 =BG061065.D

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

1) I 1,4-Dichlorobenzen...	-----ISTD-----									
2) 1,4-Dioxane	0.457	0.464	0.337	0.401	0.367	0.381	0.377	0.398	11.77	
3) Pyridine	1.143	1.136	1.111	1.242	1.182	1.206	1.170	1.170	3.81	
4) n-Nitrosodimet...	1.074	1.202	1.072	1.152	1.114	1.124	1.132	1.124	4.00	
5) S 2-Fluorophenol	0.992	0.989	0.925	1.023	0.978	0.994	0.983	0.983	3.01	
6) Aniline	1.346	1.419	1.530	1.452	1.450	1.509	1.389	1.442	4.46	
7) S Phenol-d6	1.375	1.382	1.540	1.466	1.433	1.510	1.477	1.455	4.26	
8) 2-Chlorophenol	1.121	1.211	1.271	1.228	1.178	1.207	1.229	1.206	3.89	
9) Benzaldehyde	0.983	0.996	0.908	0.803	0.714	0.709	0.627	0.820	17.69	
10) C Phenol	1.388	1.415	1.569	1.465	1.481	1.534	1.504	1.479	4.30	
11) bis(2-Chloroet...	0.972	1.012	1.106	1.007	0.991	1.032	1.034	1.022	4.21	
12) 1,3-Dichlorobe...	1.396	1.425	1.559	1.419	1.379	1.381	1.397	1.422	4.41	
13) C 1,4-Dichlorobe...	1.491	1.467	1.553	1.496	1.411	1.429	1.442	1.470	3.30	
14) 1,2-Dichlorobe...	1.419	1.393	1.467	1.415	1.386	1.414	1.433	1.418	1.88	
15) Benzyl Alcohol	1.295	1.158	1.446	1.302	1.282	1.353	1.351	1.312	6.69	
16) 2,2'-oxybis(1-...	2.139	2.208	2.110	2.241	2.221	2.310	2.296	2.218	3.34	
17) 2-Methylphenol	0.958	1.041	1.012	1.095	1.065	1.096	1.114	1.054	5.23	
18) Hexachloroethane	0.507	0.569	0.469	0.564	0.516	0.521	0.526	0.525	6.50	
19) P n-Nitroso-di-n...	1.009	1.200	1.057	1.035	1.069	1.104	1.140	1.097	1.089	5.59
20) 3+4-Methylphenols	1.355	1.528	1.476	1.560	1.535	1.602	1.591	1.521	5.56	
21) I Naphthalene-d8	-----ISTD-----									
22) Acetophenone	0.499	0.485	0.525	0.491	0.485	0.497	0.504	0.498	2.73	
23) S Nitrobenzene-d5	0.410	0.390	0.417	0.388	0.401	0.402	0.417	0.403	2.92	
24) Nitrobenzene	0.384	0.377	0.415	0.392	0.392	0.397	0.406	0.395	3.27	
25) Isophorone	0.791	0.764	0.772	0.705	0.711	0.738	0.739	0.746	4.28	
26) C 2-Nitrophenol	0.180	0.189	0.190	0.181	0.184	0.192	0.192	0.187	2.70	
27) 2,4-Dimethylph...	0.205	0.213	0.226	0.208	0.220	0.217	0.220	0.216	3.50	
28) bis(2-Chloroet...	0.383	0.366	0.381	0.348	0.354	0.364	0.366	0.366	3.48	
29) C 2,4-Dichloroph...	0.333	0.323	0.357	0.328	0.334	0.337	0.351	0.338	3.60	
30) 1,2,4-Trichlor...	0.415	0.390	0.429	0.396	0.396	0.408	0.415	0.407	3.43	
31) Naphthalene	1.080	1.019	1.090	1.017	0.996	1.022	1.052	1.039	3.39	
32) Benzoic acid	0.195	0.224	0.275	0.248	0.263	0.265	0.274	0.249	11.92	
33) 4-Chloroaniline	0.400	0.398	0.439	0.404	0.412	0.421	0.427	0.415	3.69	
34) C Hexachlorobuta...	0.355	0.358	0.360	0.332	0.341	0.346	0.353	0.349	2.87	
35) Caprolactam	0.134	0.128	0.128	0.112	0.114	0.119	0.117	0.122	6.80	
36) C 4-Chloro-3-met...	0.392	0.418	0.426	0.401	0.390	0.409	0.415	0.407	3.35	
37) 2-Methylnaphth...	0.768	0.797	0.823	0.755	0.749	0.770	0.790	0.779	3.34	
38) 1-Methylnaphth...	0.834	0.773	0.824	0.740	0.750	0.774	0.794	0.784	4.51	

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39) I	Acenaphthene-d10	-----ISTD-----								
40)	1,2,4,5-Tetrac...	0.680	0.612	0.682	0.681	0.656	0.666	0.695	0.667	4.14
41) P	Hexachlorocycl...	0.203	0.212	0.248	0.285	0.272	0.279	0.289	0.256	13.80
42) S	2,4,6-Tribromo...	0.409	0.395	0.409	0.409	0.391	0.391	0.395	0.400	2.15
43) C	2,4,6-Trichlor...	0.414	0.403	0.431	0.442	0.428	0.436	0.445	0.429	3.49
44)	2,4,5-Trichlor...	0.452	0.470	0.505	0.500	0.486	0.484	0.497	0.485	3.86
45) S	2-Fluorobiphenyl	1.340	1.312	1.362	1.384	1.332	1.362	1.377	1.353	1.91
46)	1,1'-Biphenyl	1.306	1.244	1.285	1.306	1.258	1.278	1.298	1.282	1.88
47)	2-Chloronaphth...	1.018	0.983	1.048	1.061	1.017	1.015	1.046	1.027	2.58
48)	2-Nitroaniline	0.380	0.370	0.374	0.389	0.379	0.386	0.389	0.381	1.86
49)	Acenaphthylene	1.571	1.507	1.619	1.620	1.558	1.577	1.588	1.577	2.46
50)	Dimethylphthalate	1.647	1.534	1.593	1.552	1.492	1.496	1.499	1.545	3.75
51)	2,6-Dinitrotol...	0.354	0.326	0.323	0.338	0.326	0.323	0.323	0.330	3.55
52) C	Acenaphthene	0.984	0.937	1.017	1.009	0.971	0.982	0.987	0.984	2.65
53)	3-Nitroaniline	0.317	0.305	0.319	0.299	0.289	0.296	0.292	0.302	3.92
54) P	2,4-Dinitrophenol	0.185	0.180	0.203	0.227	0.234	0.228	0.236	0.213	11.12
55)	Dibenzofuran	1.695	1.651	1.684	1.695	1.640	1.636	1.631	1.662	1.72
56) P	4-Nitrophenol	0.261	0.238	0.239	0.255	0.242	0.239	0.237	0.244	3.88
57)	2,4-Dinitrotol...	0.477	0.491	0.513	0.506	0.487	0.479	0.469	0.489	3.26
58)	Fluorene	1.449	1.426	1.482	1.485	1.397	1.410	1.415	1.438	2.45
59)	2,3,4,6-Tetrac...	0.499	0.484	0.518	0.516	0.488	0.494	0.495	0.499	2.60
60)	Diethylphthalate	1.711	1.561	1.602	1.575	1.487	1.484	1.485	1.558	5.33
61)	4-Chlorophenyl...	0.873	0.881	0.886	0.871	0.844	0.840	0.849	0.863	2.17
62)	4-Nitroaniline	0.321	0.327	0.347	0.338	0.325	0.319	0.320	0.328	3.19
63)	Azobenzene	1.260	1.217	1.235	1.235	1.167	1.186	1.188	1.213	2.74
64) I	Phenanthrene-d10	-----ISTD-----								
65)	4,6-Dinitro-2-...	0.113	0.112	0.128	0.124	0.129	0.130	0.134	0.124	6.81
66) c	n-Nitrosodiphe...	0.489	0.480	0.517	0.484	0.490	0.502	0.509	0.496	2.81
67)	4-Bromophenyl-...	0.219	0.223	0.243	0.234	0.237	0.245	0.247	0.236	4.69
68)	Hexachlorobenzene	0.297	0.287	0.296	0.284	0.287	0.298	0.304	0.293	2.49
69)	Atrazine	0.232	0.209	0.211	0.200	0.197	0.194	0.190	0.205	6.98
70) C	Pentachlorophenol	0.155	0.163	0.185	0.185	0.183	0.194	0.196	0.180	8.59
71)	Phenanthrene	0.937	0.896	0.932	0.913	0.898	0.908	0.916	0.914	1.69
72)	Anthracene	0.904	0.890	0.952	0.909	0.891	0.918	0.923	0.912	2.35
73)	Carbazole	0.855	0.832	0.805	0.801	0.772	0.773	0.778	0.802	3.95
74)	Di-n-butylphth...	1.022	0.960	0.965	0.965	0.916	0.921	0.929	0.954	3.87
75) C	Fluoranthene	1.360	1.283	1.278	1.238	1.209	1.198	1.203	1.253	4.68
76) I	Chrysene-d12	-----ISTD-----								
77)	Benzidine	0.179	0.208	0.349	0.413	0.386	0.377	0.385	0.328	28.75
78)	Pyrene	1.267	1.237	1.307	1.264	1.264	1.272	1.291	1.272	1.76
79) S	Terphenyl-d14	1.219	1.172	1.264	1.185	1.181	1.194	1.202	1.203	2.60
80)	Butylbenzylphth...	0.405	0.379	0.417	0.404	0.397	0.404	0.411	0.402	3.03
81)	Benzo(a)anthra...	1.383	1.264	1.336	1.282	1.282	1.294	1.320	1.309	3.11
82)	3,3'-Dichlorob...	0.482	0.451	0.486	0.472	0.463	0.468	0.474	0.471	2.52
83)	Chrysene	1.301	1.204	1.233	1.200	1.196	1.212	1.237	1.226	3.00
84)	Bis(2-ethylhex...	0.524	0.497	0.529	0.521	0.516	0.540	0.553	0.526	3.43
85) c	Di-n-octyl pht...	0.929	0.858	0.938	0.933	0.922	0.936	0.975	0.927	3.76

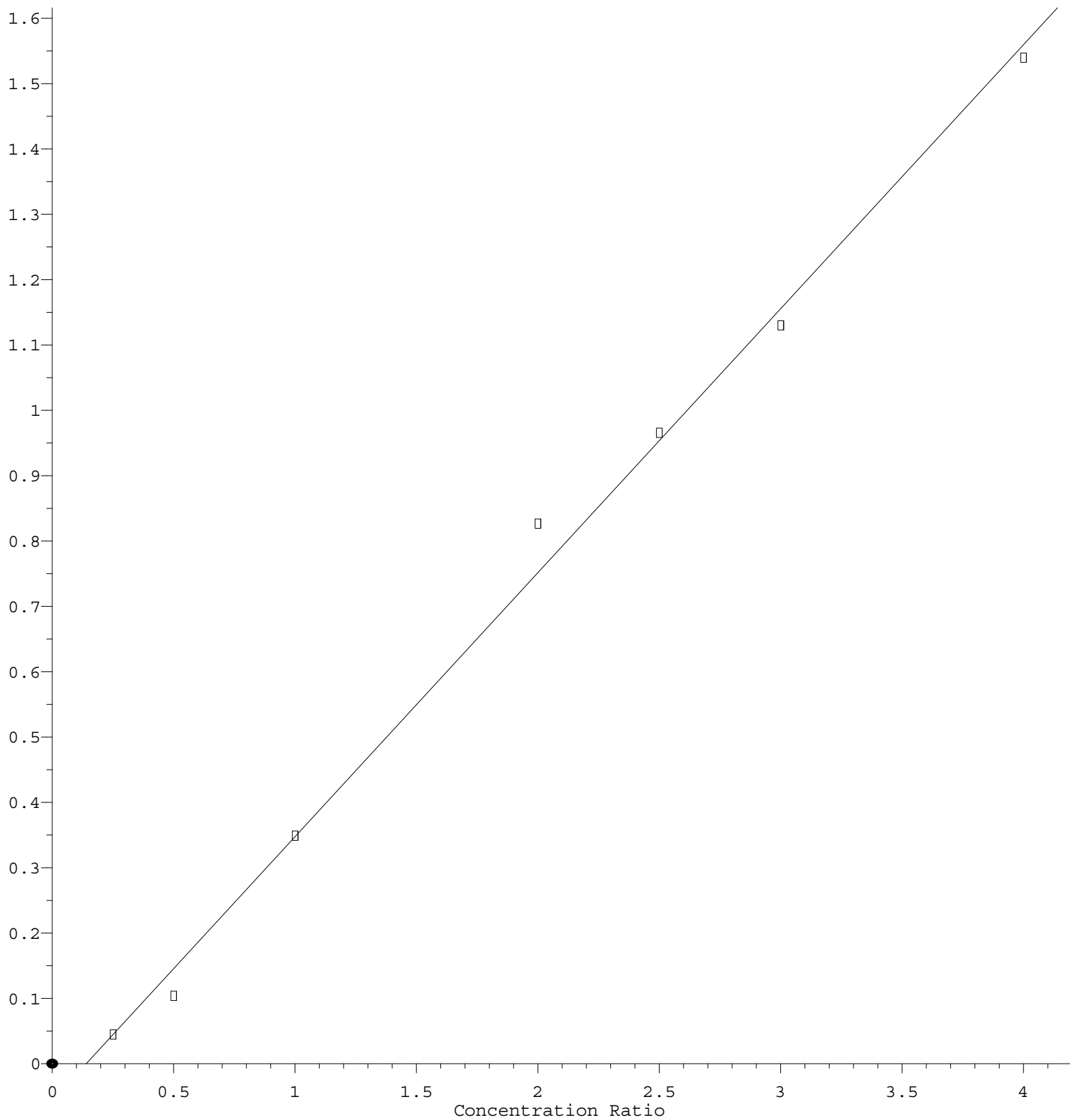
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86) I	Perylene-d12	-----ISTD-----									
87)	Indeno(1,2,3-c...	1.380	1.256	1.345	1.298	1.297	1.347	1.355	1.325	3.25	
88)	Benzo(b)fluora...	1.164	1.062	1.145	1.081	1.092	1.141	1.139	1.118	3.45	
89)	Benzo(k)fluora...	1.160	1.053	1.086	1.062	1.036	1.058	1.072	1.075	3.75	
90) C	Benzo(a)pyrene	1.043	0.922	0.988	0.951	0.945	0.985	0.982	0.974	4.02	
91)	Dibenzo(a,h)an...	1.136	1.051	1.120	1.070	1.070	1.120	1.116	1.098	3.00	
92)	Benzo(g,h,i)pe...	1.110	1.057	1.107	1.076	1.066	1.103	1.116	1.091	2.18	

(#) = Out of Range

Benzidine

Response Ratio



$$\text{Response} = 4.041\text{e-}001 * \text{Amt} - 5.648\text{e-}002$$

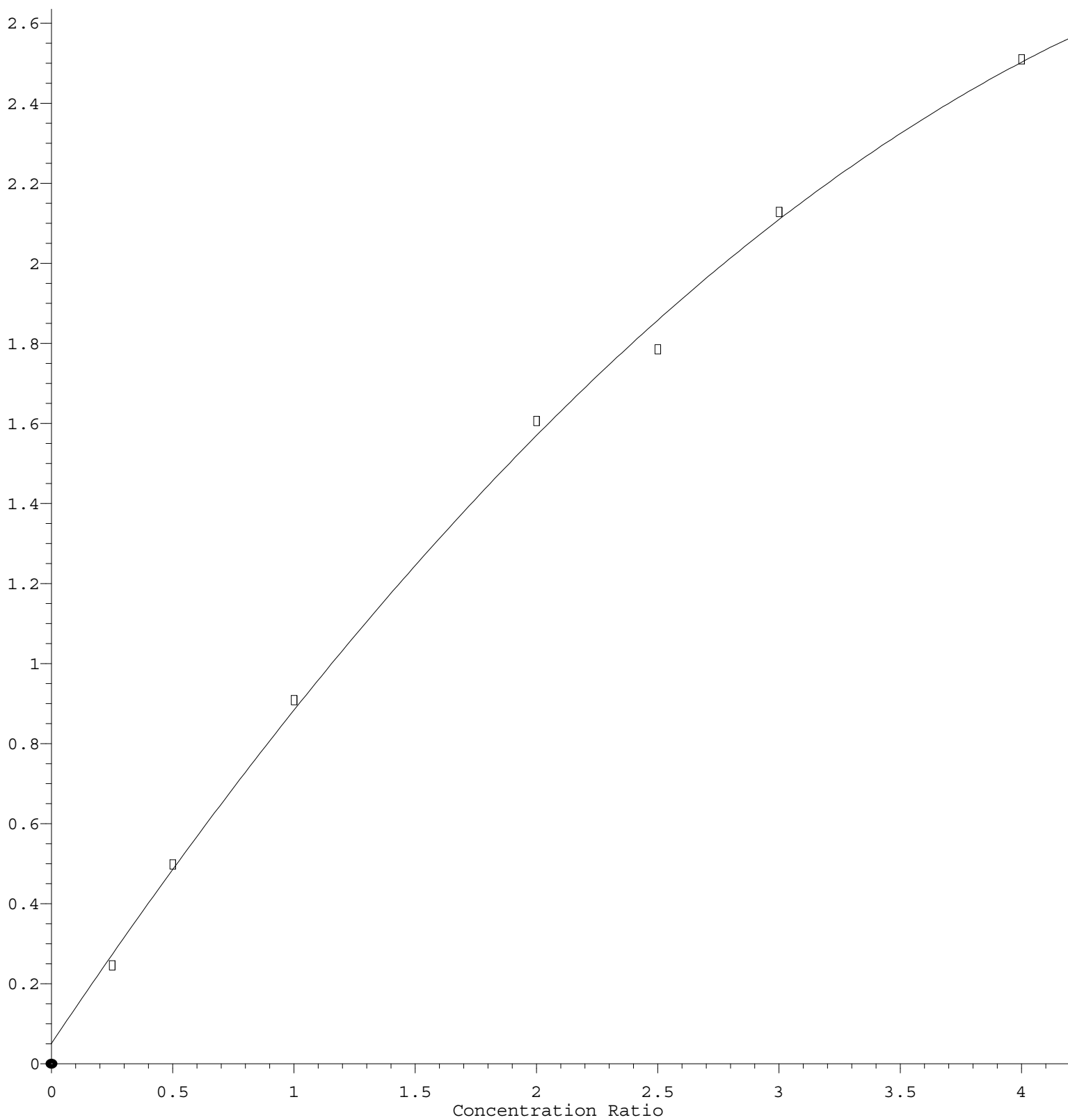
Coef of Det (r^2) = 0.995464 Curve Fit: Linear

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Benzaldehyde

Response Ratio



$R = -7.330e-002 A^2 + 9.061e-001 A + 5.074e-002$
Coef of Det (r^2) = 0.998031 Curve Fit: Quadratic
Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG043024.M
Calibration Table Last Updated: Wed May 01 05:11:25 2024