

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG061923.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Tue Jun 20 03:40:38 2023
 Response Via : Initial Calibration

Calibration Files

2.5 =BG057753.D 5 =BG057754.D 10 =BG057755.D 20 =BG057756.D 40 =BG057757.D 50 =BG057758.D 60 =BG057759.D 80 =BG057760.D

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

1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.651	0.622	0.664	0.581	0.539	0.550	0.526	0.590		9.43
3)	Pyridine	1.784	1.818	1.925	1.867	1.748	1.747	1.699	1.798		4.32
4)	n-Nitrosodimet...	0.902	0.849	0.856	0.808	0.781	0.779	0.777	0.822		5.88
5) S	2-Fluorophenol	1.371	1.340	1.382	1.307	1.240	1.251	1.190	1.297		5.58
6)	Aniline	2.577	2.593	2.705	2.592	2.494	2.417	2.362	2.534		4.63
7) S	Phenol-d6	2.006	2.020	2.120	2.030	1.944	1.891	1.843	1.979		4.72
8)	2-Chlorophenol	1.433	1.344	1.431	1.339	1.287	1.269	1.237	1.334		5.72
9)	Benzaldehyde	1.433	1.370	1.293	1.115	1.012			1.244		14.17
10) C	Phenol	1.997	2.009	2.067	1.989	1.914	1.862	1.802	1.949		4.78
11)	bis(2-Chloroet...	1.620	1.522	1.601	1.508	1.445	1.413	1.379	1.498		6.11
12)	1,3-Dichlorobe...	1.430	1.426	1.479	1.374	1.284	1.289	1.224	1.358		6.91
13) C	1,4-Dichlorobe...	1.495	1.410	1.491	1.366	1.317	1.291	1.262	1.376		6.78
14)	1,2-Dichlorobe...	1.412	1.413	1.434	1.344	1.263	1.252	1.203	1.331		6.94
15)	Benzyl Alcohol	1.517	1.486	1.568	1.586	1.504	1.464	1.468	1.513		3.16
16)	2,2'-oxybis(1-...	3.056	2.953	3.080	2.918	2.775	2.726	2.638	2.878		5.86
17)	2-Methylphenol	1.442	1.451	1.541	1.458	1.412	1.370	1.360	1.433		4.28
18)	Hexachloroethane	0.496	0.506	0.515	0.492	0.462	0.474	0.450	0.485		4.91
19) P	n-Nitroso-di-n...	1.328	1.426	1.392	1.500	1.453	1.352	1.302	1.286	1.380	5.50
20)	3+4-Methylphenols	1.904	1.951	2.127	2.087	1.974	1.919	1.896	1.980		4.64
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.572	0.592	0.599	0.559	0.530	0.554	0.520	0.561		5.25
23) S	Nitrobenzene-d5	0.349	0.372	0.407	0.401	0.390	0.415	0.392	0.390		5.76
24)	Nitrobenzene	0.378	0.401	0.423	0.411	0.400	0.426	0.402	0.406		3.99
25)	Isophorone	0.879	0.883	0.940	0.909	0.871	0.896	0.853	0.890		3.17
26) C	2-Nitrophenol	0.106	0.114	0.133	0.144	0.148			0.129		14.18
27)	2,4-Dimethylph...	0.324	0.317	0.334	0.326	0.309	0.319	0.308	0.319		2.96
28)	bis(2-Chloroet...	0.479	0.494	0.505	0.480	0.462	0.481	0.455	0.479		3.60
29) C	2,4-Dichloroph...	0.279	0.289	0.317	0.314	0.304	0.317	0.304	0.303		4.85
30)	1,2,4-Trichlor...	0.337	0.357	0.352	0.330	0.323	0.347	0.322	0.338		4.16
31)	Naphthalene	1.104	1.117	1.120	1.056	1.009	1.053	0.992	1.064		4.84
32)	Benzoic acid		0.182	0.217	0.235	0.247	0.262	0.268	0.235		13.62
33)	4-Chloroaniline	0.474	0.500	0.526	0.498	0.479	0.496	0.480	0.493		3.63
34) C	Hexachlorobuta...	0.213	0.207	0.217	0.200	0.198	0.211	0.198	0.206		3.75
35)	Caprolactam	0.114	0.118	0.138	0.129	0.120	0.123	0.120	0.123		6.41
36) C	4-Chloro-3-met...	0.359	0.389	0.424	0.407	0.396	0.406	0.392	0.396		5.09
37)	2-Methylnaphth...	0.788	0.799	0.823	0.778	0.747	0.769	0.727	0.776		4.14
38)	1-Methylnaphth...	0.729	0.761	0.788	0.738	0.708	0.719	0.685	0.732		4.65

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG061923.M

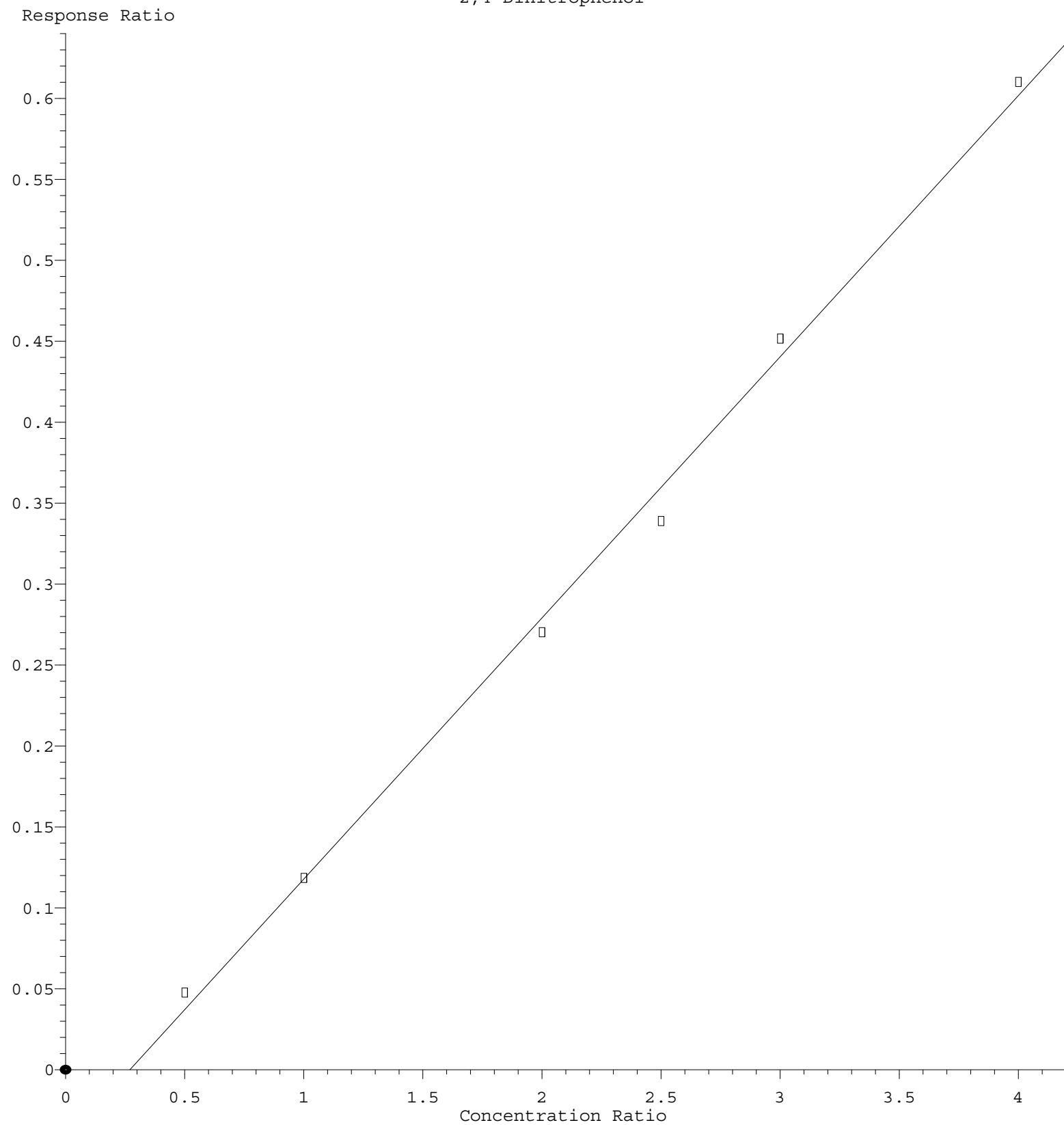
39) I	Acenaphthene-d10	-----ISTD-----								
40)	1,2,4,5-Tetrac...	0.534	0.574	0.586	0.558	0.532	0.569	0.529	0.554	4.13
41) P	Hexachlorocycl...	0.258	0.278	0.297	0.303	0.299	0.317	0.308	0.294	6.88
42) S	2,4,6-Tribromo...	0.257	0.272	0.300	0.289	0.273	0.285	0.275	0.279	4.96
43) C	2,4,6-Trichlor...	0.358	0.384	0.415	0.412	0.392	0.416	0.404	0.397	5.31
44)	2,4,5-Trichlor...	0.396	0.456	0.493	0.473	0.465	0.490	0.466	0.463	6.95
45) S	2-Fluorobiphenyl	1.418	1.430	1.413	1.310	1.222	1.265	1.172	1.319	7.89
46)	1,1'-Biphenyl	1.430	1.452	1.452	1.382	1.283	1.337	1.277	1.373	5.51
47)	2-Chloronaphth...	1.076	1.103	1.114	1.054	1.002	1.047	0.989	1.055	4.49
48)	2-Nitroaniline	0.287	0.312	0.378	0.401	0.402	0.426	0.419	0.375	14.51
49)	Acenaphthylene	1.833	1.908	1.938	1.833	1.717	1.773	1.675	1.811	5.30
50)	Dimethylphthalate	1.580	1.640	1.666	1.541	1.442	1.471	1.414	1.536	6.38
51)	2,6-Dinitrotol...	0.210	0.252	0.299	0.309	0.296	0.314	0.300	0.283	13.38
52) C	Acenaphthene	1.121	1.159	1.181	1.130	1.074	1.111	1.070	1.121	3.65
53)	3-Nitroaniline	0.258	0.301	0.338	0.343	0.326	0.335	0.331	0.319	9.42
54) P	2,4-Dinitrophenol		0.095	0.118	0.135	0.136	0.151	0.153	0.131	16.38
55)	Dibenzofuran	1.856	1.936	1.942	1.823	1.690	1.735	1.637	1.803	6.62
56) P	4-Nitrophenol	0.197	0.233	0.270	0.274	0.259	0.274	0.266	0.253	11.24
57)	2,4-Dinitrotol...	0.284	0.324	0.401	0.419	0.406	0.429	0.422	0.384	14.66
58)	Fluorene	1.545	1.537	1.565	1.436	1.316	1.332	1.257	1.427	8.82
59)	2,3,4,6-Tetrac...	0.365	0.379	0.418	0.416	0.381	0.406	0.395	0.394	5.14
60)	Diethylphthalate	1.631	1.669	1.712	1.607	1.474	1.503	1.435	1.576	6.70
61)	4-Chlorophenyl...	0.796	0.798	0.816	0.761	0.712	0.730	0.698	0.759	6.11
62)	4-Nitroaniline	0.269	0.321	0.353	0.349	0.323	0.333	0.327	0.325	8.51
63)	Azobenzene	1.720	1.704	1.736	1.614	1.498	1.543	1.469	1.612	6.88
64) I	Phenanthrene-d10	-----ISTD-----								
65)	4,6-Dinitro-2-...		0.062	0.076	0.085	0.088	0.096	0.098	0.084	15.88
66) c	n-Nitrosodiphe...	0.525	0.547	0.552	0.532	0.509	0.528	0.503	0.528	3.43
67)	4-Bromophenyl-...	0.206	0.219	0.216	0.213	0.205	0.218	0.207	0.212	2.85
68)	Hexachlorobenzene	0.213	0.221	0.220	0.217	0.210	0.223	0.210	0.216	2.48
69)	Atrazine	0.184	0.192	0.199	0.182	0.171	0.174	0.166	0.181	6.59
70) C	Pentachlorophenol	0.133	0.145	0.165	0.167	0.168	0.173	0.167	0.160	9.39
71)	Phenanthrene	1.037	1.028	1.031	0.961	0.907	0.934	0.880	0.968	6.65
72)	Anthracene	1.043	1.035	1.054	0.993	0.933	0.958	0.906	0.989	5.88
73)	Carbazole	0.966	0.973	0.988	0.906	0.849	0.872	0.828	0.912	7.09
74)	Di-n-butylphth...	1.079	1.108	1.172	1.101	1.031	1.061	1.004	1.080	5.12
75) C	Fluoranthene	1.245	1.229	1.259	1.146	1.062	1.093	1.041	1.154	7.91
76) I	Chrysene-d12	-----ISTD-----								
77)	Benzidine	0.543	0.494	0.442	0.505	0.456	0.422	0.420	0.469	9.86
78)	Pyrene	1.481	1.506	1.503	1.451	1.359	1.392	1.295	1.427	5.64
79) S	Terphenyl-d14	1.183	1.171	1.152	1.065	1.000	1.013	0.924	1.073	9.23
80)	Butylbenzylpht...	0.484	0.510	0.557	0.569	0.552	0.580	0.547	0.543	6.26
81)	Benzo(a)anthra...	1.399	1.420	1.433	1.387	1.319	1.369	1.278	1.372	4.05
82)	3,3'-Dichlorob...	0.453	0.475	0.507	0.486	0.455	0.464	0.428	0.467	5.47
83)	Chrysene	1.386	1.341	1.375	1.321	1.262	1.312	1.223	1.317	4.45
84)	Bis(2-ethylhex...	0.750	0.806	0.830	0.815	0.767	0.808	0.740	0.788	4.46
85) c	Di-n-octyl pht...	1.230	1.341	1.419	1.456	1.389	1.467	1.373	1.382	5.83

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
Method File : 8270-BG061923.M

86) I	Perylene-d12	-----ISTD-----									
87)	Indeno(1,2,3-c...	1.267	1.301	1.355	1.340	1.281	1.329	1.280	1.308	2.61	
88)	Benzo(b)fluora...	1.098	1.118	1.140	1.116	1.053	1.103	1.043	1.096	3.23	
89)	Benzo(k)fluora...	1.145	1.154	1.171	1.130	1.081	1.098	1.035	1.116	4.27	
90) C	Benzo(a)pyrene	1.047	1.086	1.131	1.108	1.047	1.088	1.027	1.076	3.47	
91)	Dibenzo(a,h)an...	1.082	1.084	1.125	1.109	1.055	1.092	1.048	1.085	2.53	
92)	Benzo(g,h,i)pe...	1.066	1.082	1.120	1.107	1.065	1.105	1.059	1.086	2.25	

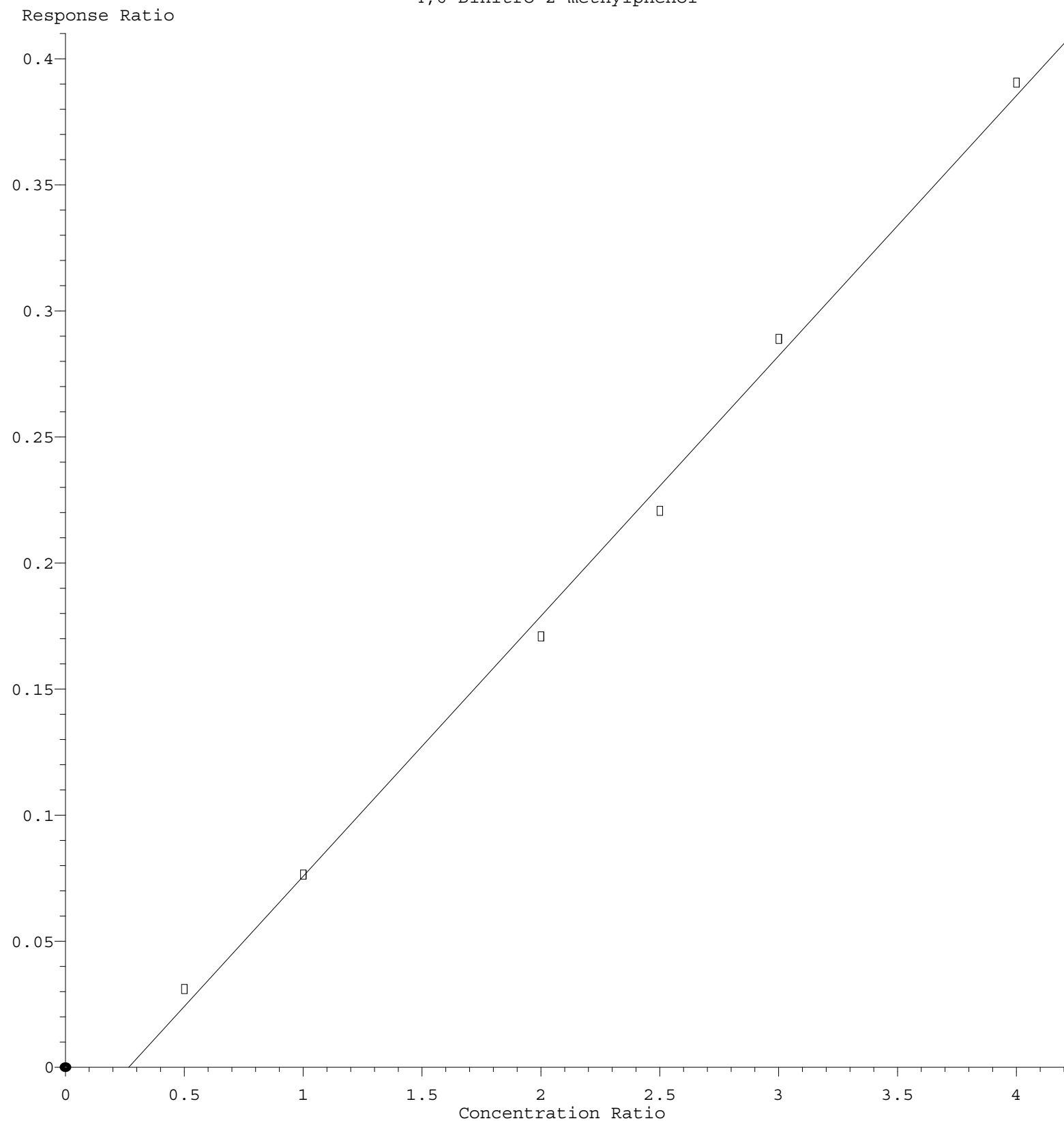
(#) = Out of Range

2,4-Dinitrophenol



Response = $1.614 \times 10^{-1} \times \text{Amt} - 4.358 \times 10^{-2}$
Coef of Det (r^2) = 0.996241 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG061923.M
Calibration Table Last Updated: Tue Jun 20 03:40:38 2023

4,6-Dinitro-2-methylphenol



$\text{Response} = 1.033\text{e-}001 * \text{Amt} - 2.744\text{e-}002$
 Coef of Det (r^2) = 0.996850 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG061923.M
 Calibration Table Last Updated: Tue Jun 20 03:40:38 2023