

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG040725.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Mon Apr 07 17:00:17 2025
 Response Via : Initial Calibration

Calibration Files

2.5 =BG064196.D 5 =BG064197.D 10 =BG064198.D 20 =BG064199.D 40 =BG064200.D 50 =BG064204.D 60 =BG064202.D 80 =BG064203.D

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

1) I 1,4-Dichlorobenzen...	-----ISTD-----									
2) 1,4-Dioxane	0.556	0.503	0.522	0.473	0.476	0.445	0.474	0.493		7.52
3) Pyridine	1.314	1.222	1.351	1.280	1.359	1.298	1.414	1.320		4.69
4) n-Nitrosodimet...	1.043	0.906	0.994	0.988	1.118	0.978	1.075	1.015		6.91
5) S 2-Fluorophenol	1.152	1.153	1.305	1.227	1.306	1.268	1.293	1.244		5.45
6) Aniline	1.679	1.644	1.847	1.653	1.621	1.544	1.391	1.626		8.49
7) S Phenol-d6	1.596	1.610	1.833	1.771	1.865	1.793	1.857	1.761		6.41
8) 2-Chlorophenol	1.292	1.244	1.456	1.406	1.445	1.401	1.477	1.389		6.32
9) Benzaldehyde	1.192	1.094	1.107	0.916	0.963	0.756		1.004		15.73
10) C Phenol	1.539	1.591	1.785	1.761	1.874	1.784	1.883	1.745		7.59
11) bis(2-Chloroet...	1.290	1.337	1.404	1.355	1.353	1.338	1.381	1.351		2.66
12) 1,3-Dichlorobe...	1.533	1.422	1.584	1.517	1.499	1.471	1.526	1.507		3.39
13) C 1,4-Dichlorobe...	1.558	1.450	1.578	1.491	1.525	1.487	1.541	1.519		2.95
14) 1,2-Dichlorobe...	1.406	1.490	1.649	1.502	1.491	1.466	1.533	1.505		4.94
15) Benzyl Alcohol	1.361	1.410	1.629	1.518	1.603	1.578	1.623	1.532		7.00
16) 2,2'-oxybis(1-...	3.060	2.949	3.113	2.989	3.031	2.922	3.048	3.016		2.22
17) 2-Methylphenol	1.220	1.186	1.265	1.252	1.278	1.273	1.317	1.256		3.39
18) Hexachloroethane	0.567	0.552	0.617	0.597	0.609	0.600	0.610	0.593		4.11
19) P n-Nitroso-di-n...	1.176	1.358	1.367	1.370	1.308	1.323	1.307	1.340	1.319	4.76
20) 3+4-Methylphenols		1.676	1.645	1.879	1.783	1.813	1.774	1.831	1.771	4.72
21) I Naphthalene-d8	-----ISTD-----									
22) Acetophenone	0.545	0.536	0.569	0.543	0.536	0.552	0.547	0.547		2.07
23) S Nitrobenzene-d5	0.377	0.377	0.403	0.408	0.404	0.416	0.421	0.401		4.37
24) Nitrobenzene	0.378	0.374	0.419	0.407	0.398	0.411	0.408	0.399		4.34
25) Isophorone	0.750	0.717	0.758	0.740	0.721	0.728	0.728	0.735		2.07
26) C 2-Nitrophenol	0.142	0.153	0.169	0.174	0.171	0.177	0.184	0.167		8.79
27) 2,4-Dimethylph...	0.210	0.220	0.236	0.228	0.228	0.235	0.235	0.227		4.28
28) bis(2-Chloroet...	0.422	0.412	0.440	0.428	0.418	0.423	0.426	0.424		2.10
29) C 2,4-Dichloroph...	0.284	0.270	0.307	0.300	0.301	0.317	0.315	0.299		5.69
30) 1,2,4-Trichlor...	0.337	0.312	0.323	0.335	0.325	0.338	0.338	0.330		3.07
31) Naphthalene	1.123	1.066	1.104	1.083	1.076	1.094	1.087	1.090		1.74
32) Benzoic acid		0.171	0.204	0.245	0.259	0.266	0.272	0.236		17.00
33) 4-Chloroaniline	0.387	0.383	0.414	0.404	0.391	0.393	0.390	0.395		2.72
34) C Hexachlorobuta...	0.214	0.231	0.242	0.231	0.226	0.235	0.233	0.230		3.73
35) Caprolactam	0.119	0.106	0.114	0.118	0.114	0.112	0.111	0.113		3.80
36) C 4-Chloro-3-met...	0.408	0.388	0.402	0.405	0.398	0.404	0.400	0.401		1.59
37) 2-Methylnaphth...	0.816	0.774	0.816	0.789	0.773	0.793	0.792	0.793		2.21
38) 1-Methylnaphth...	0.809	0.774	0.813	0.784	0.754	0.784	0.783	0.786		2.57

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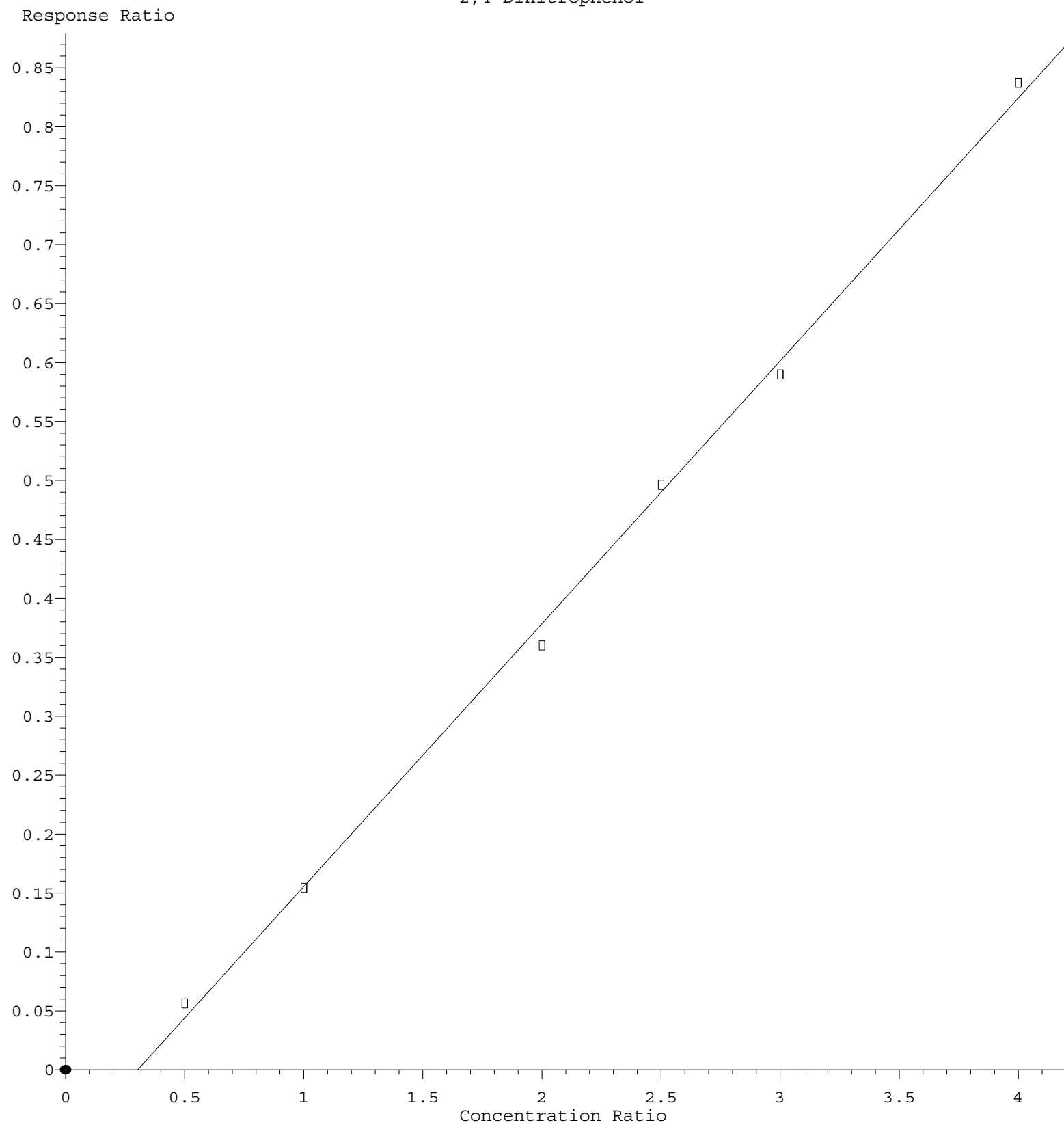
39) I	Acenaphthene-d10	-----ISTD-----									
40)	1,2,4,5-Tetrac...	0.577	0.570	0.606	0.602	0.572	0.612	0.608	0.593	3.15	
41) P	Hexachlorocycl...	0.162	0.176	0.220	0.232	0.221	0.242	0.238	0.213	14.76	
42) S	2,4,6-Tribromo...	0.256	0.267	0.269	0.274	0.268	0.272	0.273	0.269	2.25	
43) C	2,4,6-Trichlor...	0.334	0.365	0.373	0.390	0.389	0.397	0.399	0.378	6.16	
44)	2,4,5-Trichlor...	0.380	0.413	0.425	0.438	0.426	0.437	0.450	0.424	5.38	
45) S	2-Fluorobiphenyl	1.371	1.364	1.426	1.400	1.329	1.371	1.366	1.375	2.22	
46)	1,1'-Biphenyl	1.529	1.534	1.576	1.570	1.517	1.569	1.586	1.554	1.75	
47)	2-Chloronaphth...	1.104	1.108	1.135	1.131	1.083	1.139	1.140	1.120	1.96	
48)	2-Nitroaniline	0.344	0.390	0.407	0.422	0.423	0.444	0.442	0.410	8.52	
49)	Acenaphthylene	1.733	1.725	1.778	1.795	1.718	1.777	1.788	1.759	1.85	
50)	Dimethylphthalate	1.608	1.547	1.518	1.562	1.487	1.511	1.504	1.534	2.71	
51)	2,6-Dinitrotol...	0.271	0.288	0.318	0.329	0.325	0.334	0.329	0.313	7.72	
52) C	Acenaphthene	1.201	1.144	1.187	1.176	1.135	1.178	1.179	1.172	2.00	
53)	3-Nitroaniline	0.262	0.281	0.310	0.334	0.307	0.327	0.326	0.307	8.57	
54) P	2,4-Dinitrophenol		0.113	0.154	0.180	0.198	0.197	0.209	0.175	20.65	
55)	Dibenzofuran	1.891	1.889	1.904	1.917	1.856	1.897	1.886	1.891	0.99	
56) P	4-Nitrophenol	0.191	0.211	0.248	0.274	0.262	0.276	0.277	0.249	13.83	
57)	2,4-Dinitrotol...	0.399	0.415	0.442	0.471	0.459	0.467	0.464	0.445	6.31	
58)	Fluorene	1.592	1.510	1.488	1.503	1.451	1.497	1.460	1.500	3.07	
59)	2,3,4,6-Tetrac...	0.383	0.397	0.415	0.424	0.411	0.416	0.425	0.410	3.70	
60)	Diethylphthalate	1.641	1.645	1.622	1.680	1.597	1.649	1.630	1.638	1.57	
61)	4-Chlorophenyl...	0.787	0.754	0.750	0.756	0.729	0.746	0.741	0.752	2.37	
62)	4-Nitroaniline	0.255	0.280	0.306	0.337	0.327	0.342	0.339	0.312	10.82	
63)	Azobenzene	1.613	1.610	1.598	1.654	1.561	1.607	1.598	1.606	1.71	
64) I	Phenanthrene-d10	-----ISTD-----									
65)	4,6-Dinitro-2-...	0.087	0.092	0.116	0.124	0.131	0.136	0.137	0.117	17.43	
66) c	n-Nitrosodiphe...	0.573	0.569	0.598	0.597	0.580	0.588	0.583	0.584	1.93	
67)	4-Bromophenyl-...	0.223	0.213	0.224	0.228	0.226	0.229	0.226	0.224	2.40	
68)	Hexachlorobenzene	0.248	0.233	0.247	0.242	0.243	0.245	0.240	0.243	2.03	
69)	Atrazine	0.205	0.194	0.168	0.128	0.186			0.176	17.17	
70) C	Pentachlorophenol	0.098	0.132	0.158	0.168	0.165	0.171	0.172	0.152	18.03	
71)	Phenanthrene	1.090	1.049	1.114	1.089	1.067	1.093	1.052	1.079	2.20	
72)	Anthracene	1.034	1.057	1.107	1.101	1.070	1.089	1.062	1.074	2.44	
73)	Carbazole	0.930	0.953	1.002	1.007	0.992	1.010	0.975	0.981	3.09	
74)	Di-n-butylphth...	1.166	1.208	1.258	1.260	1.234	1.277	1.211	1.231	3.12	
75) C	Fluoranthene	1.289	1.278	1.340	1.310	1.271	1.312	1.246	1.292	2.41	
76) I	Chrysene-d12	-----ISTD-----									
77)	Benzidine	0.242	0.223	0.143	0.361	0.305	0.212	0.360	0.264	30.91	
78)	Pyrene	1.315	1.236	1.309	1.315	1.269	1.279	1.275	1.285	2.28	
79) S	Terphenyl-d14	1.023	0.990	1.052	1.023	0.989	0.978	0.963	1.003	3.09	
80)	Butylbenzylpht...	0.477	0.478	0.531	0.549	0.536	0.547	0.540	0.523	5.99	
81)	Benzo(a)anthra...	1.270	1.242	1.321	1.323	1.298	1.297	1.304	1.294	2.22	
82)	3,3'-Dichlorob...	0.413	0.418	0.445	0.459	0.439	0.436	0.442	0.436	3.61	
83)	Chrysene	1.303	1.264	1.266	1.277	1.253	1.261	1.259	1.269	1.31	
84)	Bis(2-ethylhex...	0.733	0.758	0.811	0.806	0.780	0.796	0.781	0.781	3.55	
85) c	Di-n-octyl pht...	1.306	1.272	1.369	1.387	1.369	1.389	1.385	1.354	3.40	

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86) I	Perylene-d12	-----ISTD-----								
87)	Indeno(1,2,3-c...	1.354	1.322	1.407	1.398	1.369	1.432	1.410	1.384	2.75
88)	Benzo(b)fluora...	1.189	1.154	1.274	1.256	1.243	1.271	1.233	1.232	3.63
89)	Benzo(k)fluora...	1.228	1.162	1.228	1.190	1.161	1.233	1.226	1.204	2.68
90) C	Benzo(a)pyrene	0.995	1.041	1.109	1.080	1.069	1.132	1.103	1.076	4.32
91)	Dibenzo(a,h)an...	1.072	1.079	1.163	1.160	1.140	1.206	1.167	1.141	4.30
92)	Benzo(g,h,i)pe...	1.087	1.108	1.198	1.164	1.151	1.199	1.178	1.155	3.75

(#) = Out of Range

2,4-Dinitrophenol



Response = 2.230e-001 * Amt - 6.747e-002
Coef of Det (r^2) = 0.998000 Curve Fit: Linear
Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG040725.M
Calibration Table Last Updated: Mon Apr 07 17:00:17 2025