

Method Path : Z:\svoasrv\HPCHEM1\BNA_G\Methods\
 Method File : 8270-BG041724.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Apr 18 02:02:19 2024
 Response Via : Initial Calibration

Calibration Files

2.5 =BG060916.D 5 =BG060917.D 10 =BG060918.D 20 =BG060919.D 40 =BG060920.D 50 =BG060921.D 60 =BG060922.D 80 =BG060923.D

	Compound	2.5	5	10	20	40	50	60	80	Avg	%RSD
1) I	1,4-Dichlorobenzen...	-----ISTD-----									
2)	1,4-Dioxane	0.491	0.533	0.518	0.416	0.438	0.433	0.428	0.465	10.21	
3)	Pyridine	1.530	1.485	1.538	1.381	1.491	1.492	1.449	1.481	3.57	
4)	n-Nitrosodimet...	0.983	0.952	0.977	0.878	0.936	0.969	0.920	0.945	3.94	
5) S	2-Fluorophenol	1.207	1.214	1.197	1.052	1.134	1.164	1.092	1.151	5.37	
6)	Aniline	1.953	2.042	2.193	1.899	1.990	2.041	2.027	2.021	4.57	
7) S	Phenol-d6	1.611	1.633	1.743	1.598	1.678	1.699	1.649	1.659	3.08	
8)	2-Chlorophenol	1.296	1.330	1.385	1.227	1.297	1.342	1.293	1.310	3.77	
9)	Benzaldehyde	0.989	0.965	1.049	0.816	0.844	0.826	0.747	0.891	12.36	
10) C	Phenol	1.609	1.644	1.710	1.604	1.687	1.727	1.656	1.662	2.87	
11)	bis(2-Chloroet...	1.296	1.294	1.356	1.233	1.280	1.308	1.259	1.289	3.02	
12)	1,3-Dichlorobe...	1.406	1.434	1.477	1.277	1.405	1.402	1.342	1.392	4.65	
13) C	1,4-Dichlorobe...	1.464	1.464	1.493	1.325	1.410	1.432	1.386	1.425	3.99	
14)	1,2-Dichlorobe...	1.401	1.370	1.467	1.300	1.383	1.408	1.369	1.386	3.63	
15)	Benzyl Alcohol	1.330	1.362	1.492	1.373	1.470	1.496	1.465	1.427	4.87	
16)	2,2'-oxybis(1-...	3.108	3.088	3.223	2.905	3.151	3.195	3.144	3.116	3.34	
17)	2-Methylphenol	1.235	1.270	1.371	1.246	1.347	1.378	1.348	1.313	4.67	
18)	Hexachloroethane	0.518	0.518	0.523	0.470	0.510	0.510	0.492	0.506	3.71	
19) P	n-Nitroso-di-n...	1.294	1.173	1.212	1.320	1.205	1.307	1.328	1.269	1.263	4.67
20)	3+4-Methylphenols	1.636	1.751	1.937	1.771	1.888	1.921	1.853	1.822	5.95	
21) I	Naphthalene-d8	-----ISTD-----									
22)	Acetophenone	0.574	0.570	0.601	0.538	0.573	0.576	0.554	0.570	3.42	
23) S	Nitrobenzene-d5	0.426	0.432	0.465	0.423	0.444	0.456	0.437	0.441	3.56	
24)	Nitrobenzene	0.412	0.429	0.463	0.415	0.429	0.448	0.430	0.432	4.13	
25)	Isophorone	0.783	0.782	0.872	0.806	0.822	0.847	0.780	0.813	4.43	
26) C	2-Nitrophenol	0.160	0.171	0.189	0.180	0.191	0.194	0.186	0.182	6.72	
27)	2,4-Dimethylph...	0.335	0.317	0.325	0.307	0.315	0.328	0.314	0.320	3.00	
28)	bis(2-Chloroet...	0.471	0.464	0.477	0.433	0.450	0.465	0.440	0.457	3.57	
29) C	2,4-Dichloroph...	0.340	0.359	0.369	0.332	0.342	0.351	0.338	0.347	3.72	
30)	1,2,4-Trichlor...	0.369	0.381	0.390	0.353	0.368	0.383	0.360	0.372	3.56	
31)	Naphthalene	1.088	1.089	1.138	1.032	1.080	1.097	1.048	1.082	3.18	
32)	Benzoic acid	0.210	0.227	0.281	0.251	0.287	0.282	0.277	0.259	11.80	
33)	4-Chloroaniline	0.480	0.503	0.534	0.503	0.504	0.527	0.499	0.507	3.59	
34) C	Hexachlorobuta...	0.280	0.269	0.274	0.249	0.261	0.267	0.253	0.265	4.22	
35)	Caprolactam	0.127	0.120	0.138	0.128	0.129	0.132	0.121	0.128	4.78	
36) C	4-Chloro-3-met...	0.413	0.405	0.460	0.407	0.417	0.433	0.402	0.420	4.92	
37)	2-Methylnaphth...	0.807	0.810	0.873	0.796	0.815	0.833	0.794	0.818	3.35	
38)	1-Methylnaphth...	0.758	0.770	0.834	0.752	0.772	0.780	0.742	0.772	3.86	

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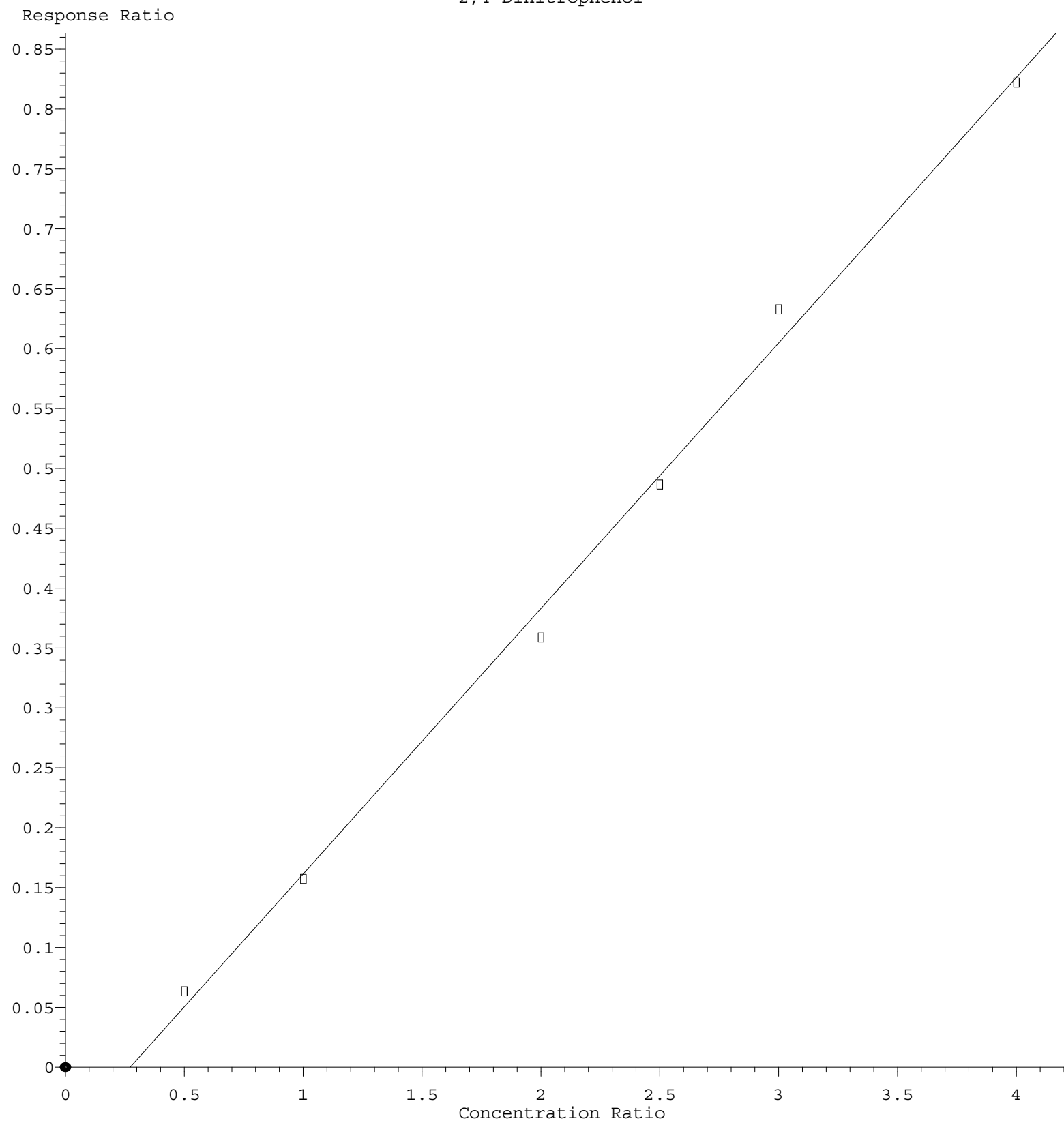
39) I	Acenaphthene-d10	-----ISTD-----								
40)	1,2,4,5-Tetrac...	0.599	0.637	0.628	0.592	0.627	0.628	0.609	0.617	2.78
41) P	Hexachlorocycl...	0.260	0.294	0.308	0.315	0.340	0.349	0.343	0.316	10.10
42) S	2,4,6-Tribromo...	0.331	0.337	0.348	0.321	0.334	0.333	0.309	0.330	3.72
43) C	2,4,6-Trichlor...	0.434	0.427	0.435	0.415	0.433	0.438	0.421	0.429	1.98
44)	2,4,5-Trichlor...	0.505	0.523	0.536	0.502	0.518	0.535	0.505	0.518	2.76
45) S	2-Fluorobiphenyl	1.403	1.408	1.418	1.296	1.361	1.367	1.287	1.363	3.90
46)	1,1'-Biphenyl	1.341	1.357	1.407	1.307	1.384	1.395	1.327	1.360	2.73
47)	2-Chloronaphth...	0.997	1.024	1.085	1.012	1.064	1.072	1.031	1.041	3.19
48)	2-Nitroaniline	0.382	0.403	0.445	0.420	0.441	0.463	0.435	0.427	6.47
49)	Acenaphthylene	1.722	1.784	1.817	1.702	1.773	1.796	1.726	1.760	2.47
50)	Dimethylphthalate	1.588	1.620	1.672	1.529	1.587	1.608	1.505	1.587	3.52
51)	2,6-Dinitrotol...	0.294	0.306	0.343	0.321	0.333	0.343	0.325	0.323	5.70
52) C	Acenaphthene	1.062	1.080	1.119	1.025	1.051	1.096	1.038	1.067	3.11
53)	3-Nitroaniline	0.337	0.350	0.382	0.348	0.360	0.367	0.341	0.355	4.44
54) P	2,4-Dinitrophenol		0.127	0.157	0.179	0.195	0.211	0.206	0.179	17.93
55)	Dibenzofuran	1.803	1.832	1.903	1.741	1.800	1.835	1.708	1.803	3.56
56) P	4-Nitrophenol	0.245	0.257	0.277	0.275	0.284	0.290	0.276	0.272	5.77
57)	2,4-Dinitrotol...	0.393	0.439	0.495	0.458	0.486	0.497	0.467	0.462	8.03
58)	Fluorene	1.500	1.539	1.606	1.454	1.475	1.492	1.403	1.496	4.30
59)	2,3,4,6-Tetrac...	0.453	0.483	0.502	0.468	0.481	0.483	0.459	0.475	3.57
60)	Diethylphthalate	1.581	1.595	1.689	1.548	1.578	1.626	1.511	1.590	3.58
61)	4-Chlorophenyl...	0.864	0.837	0.889	0.805	0.818	0.819	0.771	0.829	4.70
62)	4-Nitroaniline	0.372	0.393	0.406	0.363	0.386	0.386	0.366	0.382	4.03
63)	Azobenzene	1.464	1.504	1.588	1.452	1.506	1.538	1.436	1.498	3.54
64) I	Phenanthrene-d10	-----ISTD-----								
65)	4,6-Dinitro-2-...		0.089	0.107	0.109	0.117	0.125	0.123	0.112	11.64
66) c	n-Nitrosodiphe...	0.551	0.548	0.575	0.519	0.554	0.559	0.537	0.549	3.21
67)	4-Bromophenyl-...	0.225	0.218	0.233	0.215	0.227	0.231	0.220	0.224	3.06
68)	Hexachlorobenzene	0.242	0.241	0.253	0.234	0.245	0.252	0.240	0.244	2.80
69)	Atrazine	0.212	0.210	0.213	0.194	0.191	0.190	0.172	0.197	7.73
70) C	Pentachlorophenol	0.130	0.152	0.169	0.168	0.178	0.182	0.176	0.165	11.02
71)	Phenanthrene	1.035	1.025	1.079	0.958	1.010	1.017	0.963	1.012	4.13
72)	Anthracene	1.047	1.055	1.114	0.995	1.041	1.043	0.985	1.040	4.08
73)	Carbazole	0.937	0.936	0.960	0.858	0.897	0.905	0.850	0.906	4.56
74)	Di-n-butylphth...	1.114	1.125	1.170	1.037	1.077	1.097	1.015	1.091	4.88
75) C	Fluoranthene	1.381	1.353	1.391	1.199	1.260	1.266	1.169	1.289	6.85
76) I	Chrysene-d12	-----ISTD-----								
77)	Benzidine	0.497	0.467	0.403	0.467	0.417	0.417	0.419	0.441	8.02
78)	Pyrene	1.381	1.386	1.481	1.329	1.358	1.380	1.289	1.372	4.33
79) S	Terphenyl-d14	1.228	1.232	1.276	1.107	1.098	1.097	0.983	1.146	9.01
80)	Butylbenzylpht...	0.493	0.493	0.532	0.484	0.509	0.519	0.498	0.504	3.36
81)	Benzo(a)anthra...	1.438	1.409	1.496	1.355	1.402	1.418	1.368	1.412	3.29
82)	3,3'-Dichlorob...	0.519	0.503	0.534	0.488	0.501	0.498	0.478	0.503	3.70
83)	Chrysene	1.325	1.368	1.420	1.300	1.358	1.374	1.294	1.348	3.32
84)	Bis(2-ethylhex...	0.736	0.740	0.798	0.712	0.744	0.763	0.724	0.745	3.77
85) c	Di-n-octyl pht...	1.274	1.275	1.357	1.231	1.291	1.318	1.254	1.286	3.25

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86) I	Perylene-d12	-----ISTD-----								
87)	Indeno(1,2,3-c...	1.299	1.350	1.458	1.328	1.403	1.447	1.396	1.383	4.31
88)	Benzo(b)fluora...	1.102	1.144	1.182	1.090	1.123	1.186	1.094	1.132	3.56
89)	Benzo(k)fluora...	1.143	1.147	1.211	1.101	1.179	1.194	1.129	1.158	3.34
90) C	Benzo(a)pyrene	1.125	1.108	1.166	1.059	1.112	1.146	1.093	1.116	3.11
91)	Dibenzo(a,h)an...	1.015	1.113	1.209	1.107	1.169	1.195	1.161	1.138	5.84
92)	Benzo(g,h,i)pe...	1.076	1.093	1.134	1.071	1.130	1.167	1.120	1.113	3.12

(#) = Out of Range

2,4-Dinitrophenol



Response = $2.217 \times 10^{-1} \times \text{Amt} - 6.026 \times 10^{-2}$
 Coef of Det (r^2) = 0.996043 Curve Fit: Linear
 Method Name: Z:\svoasrv\HPCHEM1\BNA G\Methods\8270-BG041724.M
 Calibration Table Last Updated: Thu Apr 18 02:02:19 2024