

Method Path : Z:\HPCHEM1\BNA\_G\METHODS\

Method File : 8270-BG040615.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Last Update : Tue Apr 07 04:17:17 2015

Response Via : Initial Calibration

## Calibration Files

2.5 =BG016224.D 10 =BG016225.D 25 =BG016226.D 40 =BG016227.D 50 =BG016228.D 60 =BG016229.D 80 =BG016230.D

Compound	2.5	10	25	40	50	60	80	Avg	%RSD
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1) I	1,4-Dichlorobenzene	-----	ISTD-----						
2)	1,4-Dioxane	0.666	0.634	0.594	0.560	0.530	0.529	0.520	0.576
3)	Pyridine	1.722	1.679	1.784	1.719	1.708	1.711	1.729	1.722
4)	n-Nitrosodimethylamine	0.470	0.533	0.540	0.517	0.505	0.503	0.516	0.512
5) S	2-Fluorophenol	1.327	1.276	1.319	1.259	1.229	1.228	1.233	1.267
6)	Aniline	2.802	2.740	2.798	2.730	2.644	2.662	2.644	2.717
7) S	Phenol-d6	1.948	1.892	1.963	1.915	1.872	1.877	1.849	1.902
8)	2-Chlorophenol	1.468	1.534	1.579	1.520	1.507	1.520	1.525	1.522
9)	Benzaldehyde	1.306	1.241	1.260	1.066	1.036	1.001	0.891	1.114
10) C	Phenol	1.997	2.029	2.106	2.061	2.015	2.023	2.017	2.035
11)	bis(2-Chloroethane)	1.770	1.664	1.640	1.600	1.553	1.576	1.559	1.623
12)	1,3-Dichlorobenzene	1.651	1.576	1.556	1.510	1.478	1.488	1.499	1.537
13) C	1,4-Dichlorobenzene	1.799	1.638	1.600	1.528	1.526	1.536	1.532	1.594
14)	1,2-Dichlorobenzene	1.662	1.552	1.541	1.492	1.461	1.482	1.483	1.525
15)	Benzyl Alcohol	1.159	1.262	1.377	1.370	1.363	1.374	1.378	1.326
16)	2,2'-oxybis(1,6-hexamethylene)	1.939	1.875	1.861	1.822	1.744	1.783	1.771	1.828
17)	2-Methylphenol	1.326	1.327	1.403	1.389	1.367	1.377	1.365	1.365
18)	Hexachloroethane	0.702	0.615	0.603	0.592	0.571	0.587	0.599	0.610
19) P	n-Nitroso-di-n-butylamine	1.201	1.350	1.436	1.413	1.378	1.394	1.369	1.363
20)	3+4-Methylphenols	1.694	1.885	1.939	1.948	1.908	1.934	1.928	1.891
21) I	Naphthalene-d8	-----	ISTD-----						
22)	Acetophenone	0.471	0.471	0.478	0.453	0.458	0.456	0.460	0.464
23) S	Nitrobenzene-d5	0.336	0.348	0.355	0.342	0.347	0.342	0.347	0.345
24)	Nitrobenzene	0.365	0.375	0.382	0.363	0.364	0.365	0.370	0.369
25)	Isophorone	0.744	0.768	0.803	0.760	0.770	0.760	0.762	0.767
26) C	2-Nitrophenol	0.128	0.153	0.176	0.179	0.185	0.190	0.194	0.172
27)	2,4-Dimethylphenol	0.310	0.314	0.324	0.318	0.325	0.323	0.330	0.321
28)	bis(2-Chloroethane)	0.456	0.448	0.448	0.427	0.427	0.427	0.431	0.438
29) C	2,4-Dichlorophenol	0.220	0.244	0.263	0.257	0.265	0.261	0.266	0.254
30)	1,2,4-Trichlorobenzene	0.277	0.267	0.266	0.256	0.263	0.259	0.267	0.265
31)	Naphthalene	1.140	1.078	1.073	1.003	1.010	0.999	0.991	1.042
32)	Benzoic acid	0.088	0.159	0.186	0.207	0.219	0.233	0.182	29.10
33)	4-Chloroaniline	0.473	0.490	0.505	0.484	0.493	0.489	0.489	0.489
34) C	Hexachlorobutane	0.120	0.117	0.116	0.112	0.115	0.116	0.118	0.116
35)	Caprolactam	0.118	0.152	0.171	0.168	0.172	0.167	0.169	0.160
36) C	4-Chloro-3-methylphenol	0.312	0.322	0.346	0.338	0.345	0.340	0.343	0.335
37)	2-Methylnaphthalene	0.795	0.769	0.769	0.726	0.742	0.724	0.723	0.750



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83)	Bis(2-ethylhex...	0.797	0.902	0.987	0.941	0.967	0.949	0.942	0.926	6.78
84) c	Di-n-octyl pht...	1.160	1.465	1.642	1.577	1.609	1.576	1.535	1.509	10.87
85)	Indeno(1,2,3-c...	1.108	1.169	1.227	1.161	1.207	1.206	1.219	1.186	3.55
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86) I	Perylene-d12	-----ISTD-----								
87)	Benzo(b)fluora...	1.240	1.231	1.189	1.130	1.159	1.110	1.140	1.171	4.31
88)	Benzo(k)fluora...	1.195	1.211	1.217	1.073	1.099	1.146	1.083	1.146	5.41
89) C	Benzo(a)pyrene	1.099	1.143	1.127	1.055	1.080	1.089	1.091	1.098	2.66
90)	Dibenzo(a,h)an...	1.076	1.110	1.099	1.019	1.058	1.067	1.049	1.068	2.88
91)	Benzo(g,h,i)pe...	1.054	1.092	1.094	1.017	1.066	1.070	1.076	1.067	2.48

(#) = Out of Range