

Method Path : Z:\HPCHEM1\BNA_E\METHODS\
 Method File : 8270-SIM-BE021517.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Wed Feb 15 17:05:20 2017
 Response Via : Initial Calibration

Calibration Files

0.1 =BE092728.D 0.2 =BE092729.D 0.4 =BE092730.D 0.8 =BE092731.D 1.6 =BE092732.D 3.2 =BE092733.D 5 =BE092734.D
 10 =BE092735.D

	Compound		0.1	0.2	0.4	0.8	1.6	3.2	5	10	Avg	%RSD

1) I	1,4-Dichlorobenzen...	-----ISTD-----										
2)	1,4-Dioxane		0.919	0.912	0.702	0.657	0.652	0.607	0.612	0.620	0.710	18.35
3)	n-Nitrosodimet...		0.537	0.584	0.648	0.704	0.782	0.820	0.882	0.962	0.740	19.96
4) S	2-Fluorophenol		0.858	0.839	0.934	0.960	1.001	1.055	1.162		0.973	11.58
5) S	Phenol-d6		1.025	1.020	1.152	1.229	1.298	1.438			1.194	13.62
6)	bis(2-Chloroet...		1.196	1.096	1.216	1.338	1.494	1.585			1.321	14.24
7) I	Naphthalene-d8	-----ISTD-----										
8) S	Nitrobenzene-d5		0.220	0.235	0.269	0.290	0.320	0.337		0.279		16.61
9)	Naphthalene		1.164	1.071	1.091	1.087	1.061	1.047	1.074	1.064	1.082	3.31
10)	Hexachlorobuta...		0.175	0.148	0.165	0.161	0.157	0.154	0.155	0.158	0.159	5.11
11)	2-Methylnaphth...		0.633	0.625	0.662	0.680	0.693	0.695	0.715	0.735	0.680	5.61
12) I	Acenaphthene-d10	-----ISTD-----										
13) S	2,4,6-Tribromo...		0.094	0.093	0.104	0.101	0.114	0.131		0.106		13.56
14) S	2-Fluorobiphenyl		1.206	1.152	1.236	1.239	1.215	1.205	1.239	1.262	1.219	2.75
15)	Acenaphthylene		7.363	6.982	7.490	7.560	7.376	7.668	7.800	8.115	7.544	4.44
16)	Acenaphthene		1.010	1.017	1.107	1.133	1.096	1.104	1.144	1.161	1.096	5.07
17)	Fluorene		1.338	1.295	1.393	1.433	1.404	1.422	1.473	1.499	1.407	4.74
18) I	Phenanthrene-d10	-----ISTD-----										
19)	Hexachlorobenzene		0.224	0.209	0.212	0.194	0.185	0.179	0.174	0.169	0.193	10.41
20)	Pentachlorophenol			0.033	0.036	0.036	0.042	0.051	0.058		0.042	23.00
21)	Phenanthrene		1.235	1.176	1.251	1.182	1.202	1.221	1.140	1.098	1.188	4.26
22)	Anthracene		1.020	0.963	1.070	1.137	1.146	1.176	1.136	1.103	1.094	6.58
23)	Fluoranthene		5.266	4.943	5.106	4.999	5.062	5.086	5.178	5.130	5.096	1.98
24) I	Chrysene-d12	-----ISTD-----										
25)	Pyrene		4.451	4.123	4.620	4.485	5.002	4.738	5.264	5.403	4.761	9.13
26) S	Terphenyl-d14		0.596	0.549	0.619	0.583	0.662	0.635	0.722	0.739	0.638	10.43
27)	Benzo(a)anthra...		5.283	4.501	4.747	4.584	4.903	4.666	5.066	5.040	4.849	5.56
28)	Chrysene		6.404	5.292	5.757	5.504	5.848	5.476	5.810	5.921	5.752	5.94
29)	Bis(2-ethylhex...		1.099	0.557	0.582	0.492	0.635	0.677			0.674	32.34
30)	Indeno(1,2,3-c...		9.647	6.799	6.531	5.517	5.944	5.476	5.929	6.044	6.486	20.89
31) I	Perylene-d12	-----ISTD-----										
32)	Benzo(b)fluora...		2.157	1.364	1.370	1.312	1.230	1.244	1.300	1.330	1.414	21.54

Method Path : Z:\HPCHEM1\BNA_E\METHODS\

Method File : 8270-SIM-BE021517.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

33)	Benzo(k)fluora...	2.123	1.558	1.363	1.363	1.383	1.369	1.396	1.402	1.495	17.52
34) C	Benzo(a)pyrene	1.604	1.222	1.210	1.192	1.200	1.253	1.282	1.319	1.285	10.59
35)	Dibenzo(a,h)an...	1.859	1.416	1.333	1.248	1.227	1.234	1.252	1.272	1.355	15.73
36)	Benzo(g,h,i)pe...	8.144	6.052	5.676	5.307	5.160	5.034	5.172	5.280	5.728	17.99

(#) = Out of Range