

Method Path : Z:\svoasrv\HPCHEM1\BNA_N\Methods\
 Method File : 8270-SIM-BN041224.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Fri Apr 12 15:55:57 2024
 Response Via : Initial Calibration

Calibration Files

0.1 =BN030857.D 0.2 =BN030858.D 0.4 =BN030859.D 0.8 =BN030860.D 1.6 =BN030861.D 3.2 =BN030862.D 5.0 =BN030863.D

	Compound	0.1	0.2	0.4	0.8	1.6	3.2	5.0	Avg	%RSD

1) I	1,4-Dichlorobenzen...	-----ISTD-----								
2)	1,4-Dioxane	0.536	0.522	0.505	0.461	0.473	0.450	0.436	0.483	7.84
3)	n-Nitrosodimet...	0.411	0.469	0.465	0.456	0.496	0.483	0.471	0.464	5.78
4) S	2-Fluorophenol	0.989	0.985	0.917	0.917	0.977	0.962	0.970	0.960	3.16
5) S	Phenol-d6	1.139	1.125	1.081	1.101	1.211	1.228	1.251	1.162	5.77
6)	bis(2-Chloroet...	1.042	1.093	1.039	1.066	1.144	1.115	1.097	1.085	3.55
7) I	Naphthalene-d8	-----ISTD-----								
8) S	Nitrobenzene-d5	0.264	0.276	0.267	0.273	0.303	0.305	0.311	0.286	7.03
9)	Naphthalene	1.076	1.124	1.074	1.077	1.157	1.124	1.113	1.106	2.87
10)	Hexachlorobuta...	0.216	0.227	0.219	0.217	0.231	0.220	0.216	0.221	2.68
11) SURR	2-Methylnaphth...	0.597	0.612	0.589	0.594	0.642	0.629	0.628	0.613	3.37
12)	2-Methylnaphth...	0.702	0.721	0.699	0.700	0.762	0.745	0.737	0.724	3.47
13) I	Acenaphthene-d10	-----ISTD-----								
14) S	2,4,6-Tribromo...	0.159	0.150	0.143	0.145	0.167	0.179	0.191	0.162	11.11
15) S	2-Fluorobiphenyl	1.419	1.473	1.363	1.389	1.469	1.424	1.387	1.418	2.95
16)	Acenaphthylene	1.545	1.634	1.600	1.661	1.876	1.941	1.977	1.748	10.18
17)	Acenaphthene	1.167	1.249	1.176	1.194	1.301	1.286	1.270	1.235	4.44
18)	Fluorene	1.543	1.613	1.543	1.582	1.718	1.705	1.659	1.623	4.47
19) I	Phenanthrene-d10	-----ISTD-----								
20)	4,6-Dinitro-2-...	0.042	0.042	0.047	0.058	0.063		0.050		19.20
21)	4-Bromophenyl-...	0.224	0.231	0.224	0.227	0.247	0.245	0.247	0.235	4.69
22)	Hexachlorobenzene	0.266	0.273	0.263	0.268	0.292	0.280	0.280	0.274	3.72
23)	Atrazine	0.174	0.163	0.160	0.163	0.185	0.188	0.195	0.175	8.04
24)	Pentachlorophenol	0.072	0.082	0.087	0.105	0.117		0.093		19.64
25)	Phenanthrene	1.119	1.148	1.123	1.138	1.235	1.220	1.211	1.171	4.24
26)	Anthracene	0.888	0.900	0.901	0.926	1.049	1.075	1.110	0.978	9.75
27) SURR	Fluoranthene-d10	1.060	1.077	1.059	1.084	1.165	1.165	1.173	1.112	4.76
28)	Fluoranthene	1.403	1.397	1.385	1.407	1.538	1.521	1.492	1.449	4.52
29) I	Chrysene-d12	-----ISTD-----								
30)	Pyrene	1.539	1.502	1.436	1.441	1.574	1.501	1.534	1.504	3.38
31) S	Terphenyl-d14	0.988	0.919	0.859	0.880	0.938	0.883	0.898	0.909	4.76
32)	Benzo(a)anthra...	1.435	1.305	1.313	1.309	1.449	1.439	1.456	1.387	5.25
33)	Chrysene	1.584	1.482	1.412	1.443	1.534	1.483	1.464	1.486	3.85
34)	Bis(2-ethylhex...	0.642	0.548	0.522	0.582	0.635	0.702	0.605		11.04
35) I	Perylene-d12	-----ISTD-----								

Method Path : Z:\svoasrv\HPCHEM1\BNA_N\Methods\

Method File : 8270-SIM-BN041224.M

36)	Indeno(1,2,3-c...	1.426	1.401	1.421	1.529	1.697	1.767	1.710	1.564	10.02
37)	Benzo(b)fluora...	1.875	1.794	1.657	1.700	1.833	1.772	1.770	1.771	4.19
38)	Benzo(k)fluora...	1.689	1.620	1.570	1.619	1.777	1.712	1.706	1.670	4.23
39) C	Benzo(a)pyrene	1.322	1.284	1.284	1.301	1.439	1.465	1.452	1.364	6.14
40)	Dibenzo(a,h)an...	1.116	1.095	1.118	1.224	1.359	1.419	1.367	1.243	11.08
41)	Benzo(g,h,i)pe...	1.253	1.237	1.229	1.336	1.442	1.489	1.436	1.346	8.14

(#) = Out of Range