

Method Path : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\
 Method File : 8270-SIM-BN110420.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Nov 05 05:46:47 2020
 Response Via : Initial Calibration

Calibration Files

0.1 =BN012491.D 0.2 =BN012492.D 0.4 =BN012493.D 0.8 =BN012494.D 1.6 =BN012495.D 3.2 =BN012496.D 5 =BN012497.D

Compound	0.1	0.2	0.4	0.8	1.6	3.2	5	Avg	%RSD

1) I 1,4-Dichlorobenzen...	-----ISTD-----								
2) 1,4-Dioxane	0.804	0.595	0.591	0.618	0.622	0.598	0.550	0.626	13.11
3) n-Nitrosodimet...	1.722	1.543	1.697	1.841	1.805	1.645	1.709		6.33
4) S 2-Fluorophenol	1.607	1.520	1.311	1.343	1.398	1.344	1.265	1.398	8.75
5) S Phenol-d6	1.486	1.627	1.556	1.641	1.769	1.733	1.640	1.636	5.90
6) bis(2-Chloroet...	1.019	0.944	0.976	1.143	1.286	1.251	1.189	1.115	12.23
7) I Naphthalene-d8	-----ISTD-----								
8) S Nitrobenzene-d5	0.399	0.454	0.422	0.488	0.527	0.522	0.501	0.473	10.48
9) Naphthalene	1.187	1.120	1.001	1.133	1.188	1.147	1.080	1.122	5.83
10) Hexachlorobuta...	0.288	0.248	0.223	0.245	0.249	0.239	0.223	0.245	8.98
11) SURR2-Methylnaphth...	0.658	0.627	0.569	0.653	0.703	0.688	0.647	0.649	6.75
12) 2-Methylnaphth...	0.685	0.686	0.642	0.739	0.797	0.781	0.735	0.724	7.73
13) I Acenaphthene-d10	-----ISTD-----								
14) S 2,4,6-Tribromo...	0.299	0.308	0.249	0.282	0.304	0.305	0.286	0.290	7.11
15) S 2-Fluorobiphenyl	1.553	1.534	1.412	1.600	1.672	1.640	1.506	1.560	5.62
16) Acenaphthylene	2.093	1.932	1.737	2.020	2.143	2.130	1.983	2.005	7.06
17) Acenaphthene	1.404	1.242	1.107	1.282	1.340	1.321	1.237	1.276	7.41
18) Fluorene	1.703	1.582	1.416	1.653	1.713	1.726	1.597	1.627	6.68
19) I Phenanthrene-d10	-----ISTD-----								
20) 4,6-Dinitro-2-...	0.087	0.072	0.080	0.101	0.113	0.116	0.112	0.097	18.29
21) 4-Bromophenyl-...	0.236	0.214	0.202	0.235	0.254	0.247	0.232	0.231	7.90
22) Hexachlorobenzene	0.321	0.307	0.273	0.302	0.318	0.301	0.286	0.301	5.71
23) Atrazine	0.256	0.235	0.210	0.235	0.257	0.251	0.236	0.240	6.84
24) Pentachlorophenol	0.141	0.117	0.137	0.153	0.152	0.146	0.141		9.45
25) Phenanthrene	1.179	1.147	1.051	1.174	1.289	1.260	1.185	1.184	6.55
26) Anthracene	1.086	0.992	0.955	1.080	1.211	1.201	1.145	1.096	8.95
27) SURRFluoranthene-d10	1.312	1.167	1.087	1.224	1.332	1.314	1.244	1.240	7.23
28) Fluoranthene	1.463	1.373	1.235	1.384	1.514	1.502	1.417	1.413	6.77
29) I Chrysene-d12	-----ISTD-----								
30) Pyrene	1.465	1.323	1.202	1.348	1.428	1.374	1.270	1.344	6.69
31) S Terphenyl-d14	1.010	0.969	0.866	0.958	1.010	0.975	0.887	0.954	5.90
32) Benzo(a)anthra...	1.233	1.233	1.148	1.311	1.355	1.353	1.281	1.273	5.87
33) Chrysene	1.394	1.353	1.196	1.307	1.421	1.386	1.277	1.333	5.92
34) Bis(2-ethylhex...	1.035	0.867	0.751	0.818	0.866	0.875	0.834	0.864	10.06
35) Indeno(1,2,3-c...	2.091	2.095	1.936	2.003	2.083	2.002	1.830	2.006	4.85

Method Path : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\

Method File : 8270-SIM-BN110420.M

36)	I	Perylene-d12	-----ISTD-----								
37)		Benzo(b)fluora...	1.234	1.200	1.115	1.303	1.396	1.355	1.327	1.276	7.64
38)		Benzo(k)fluora...	1.524	1.261	1.220	1.361	1.466	1.387	1.300	1.360	8.05
39)	C	Benzo(a)pyrene	1.337	1.204	1.106	1.248	1.329	1.283	1.234	1.249	6.35
40)		Dibenzo(a,h)an...	1.393	1.358	1.331	1.428	1.541	1.463	1.398	1.416	4.94
41)		Benzo(g,h,i)pe...	1.522	1.495	1.374	1.468	1.531	1.442	1.376	1.458	4.42

(#) = Out of Range