

Method Path : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\
 Method File : 8270-SIM-BN100421.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Mon Oct 04 17:52:33 2021
 Response Via : Initial Calibration

Calibration Files

0.1 =BN016708.D 0.2 =BN016709.D 0.4 =BN016710.D 0.8 =BN016712.D 1.6 =BN016711.D 3.2 =BN016713.D 5 =BN016714.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	0.179	0.174	0.173	0.205	0.208	0.193	0.176	0.187	7.96
3)	n-Nitrosodimet...	0.260	0.281	0.280	0.320	0.340	0.325	0.304	0.301	9.56
4) S	2-Fluorophenol	1.051	1.009	0.968	1.049	1.097	1.022	0.948	1.021	5.03
5) S	Phenol-d6	1.110	1.082	1.061	1.174	1.238	1.171	1.089	1.132	5.61
6)	bis(2-Chloroet...	1.106	1.097	1.060	1.140	1.168	1.060	0.976	1.087	5.76
7) I	Naphthalene-d8	-----ISTD-----								
8) S	Nitrobenzene-d5	0.264	0.264	0.240	0.277	0.309	0.308	0.301	0.280	9.38
9)	Naphthalene	1.146	1.114	1.038	1.104	1.153	1.083	1.022	1.094	4.59
10)	Hexachlorobuta...	0.206	0.204	0.196	0.208	0.218	0.206	0.195	0.205	3.77
11) SURR2	Methylnaphth...	0.593	0.588	0.567	0.609	0.642	0.601	0.570	0.596	4.27
12)	2-Methylnaphth...	0.708	0.704	0.671	0.714	0.741	0.684	0.645	0.695	4.50
13) I	Acenaphthene-d10	-----ISTD-----								
14) S	2,4,6-Tribromo...	0.133	0.127	0.141	0.179	0.210	0.223	0.225	0.177	24.52
15) S	2-Fluorobiphenyl	1.665	1.667	1.489	1.616	1.677	1.586	1.499	1.600	4.95
16)	Acenaphthylene	1.571	1.573	1.595	1.850	2.001	1.924	1.849	1.766	10.31
17)	Acenaphthene	1.137	1.141	1.093	1.197	1.263	1.192	1.145	1.167	4.73
18)	Fluorene	1.374	1.353	1.306	1.449	1.528	1.450	1.392	1.407	5.23
19) I	Phenanthrene-d10	-----ISTD-----								
20)	4,6-Dinitro-2-...	0.002	0.002	0.004	0.019	0.034	0.058	0.074	0.028	105.77
21)	4-Bromophenyl-...	0.217	0.228	0.221	0.254	0.266	0.264	0.256	0.244	8.58
22)	Hexachlorobenzene	0.260	0.269	0.259	0.284	0.298	0.289	0.273	0.276	5.40
23)	Atrazine	0.173	0.160	0.159	0.192	0.225	0.223	0.220	0.193	15.39
24)	Pentachlorophenol	0.022	0.033	0.033	0.066	0.091	0.118	0.128	0.076	57.60
25)	Phenanthrene	1.127	1.158	1.108	1.210	1.261	1.218	1.149	1.176	4.70
26)	Anthracene	0.978	0.962	0.937	1.080	1.172	1.138	1.083	1.050	8.72
27) SURR	Fluoranthene-d10	1.053	1.062	1.008	1.122	1.170	1.151	1.095	1.094	5.27
28)	Fluoranthene	1.270	1.291	1.233	1.356	1.408	1.326	1.249	1.305	4.79
29) I	Chrysene-d12	-----ISTD-----								
30)	Pyrene	1.467	1.394	1.344	1.425	0.779	1.439	1.359	1.315	18.28
31) S	Terphenyl-d14	1.024	0.996	0.907	0.953	0.507	0.961	0.917	0.895	19.68
32)	Benzo(a)anthra...	1.281	1.267	1.232	1.335	0.815	1.406	1.349	1.241	15.82
33)	Chrysene	1.344	1.329	1.290	1.364	0.799	1.375	1.310	1.259	16.28
34)	Bis(2-ethylhex...	0.916	0.735	0.655	0.696	0.480	0.887	0.898	0.752	21.22
35) I	Perylene-d12	-----ISTD-----								

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

Method File : 8270-SIM-BN100421.M

36)	Indeno(1,2,3-c...	1.160	1.204	1.202	1.442	0.894	1.524	1.463	1.270	17.43
37)	Benzo(b)fluora...	1.324	1.371	1.333	1.413	0.869	1.410	1.334	1.293	14.73
38)	Benzo(k)fluora...	1.328	1.286	1.259	1.387	0.808	1.359	1.272	1.243	15.87
39) C	Benzo(a)pyrene	1.184	1.177	1.168	1.274	0.770	1.299	1.230	1.157	15.38
40)	Dibenzo(a,h)an...	0.889	0.932	0.940	1.148	0.714	1.236	1.195	1.008	18.88
41)	Benzo(g,h,i)pe...	1.048	1.088	1.090	1.261	0.782	1.337	1.284	1.127	16.78

(#) = Out of Range