

Method Path : Z:\HPCHEM1\BNA_M\METHODS\
 Method File : SOM01.2-EPA-SIM-BM030315.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Wed Mar 04 02:50:23 2015
 Response Via : Initial Calibration

Calibration Files

0.1 =BM000617.D 0.2 =BM000618.D 0.4 =BM000619.D
 0.8 =BM000620.D 1 =BM000621.D

Compound		0.1	0.2	0.4	0.8	1	Avg	%RSD

1) I	1,4-Dichlorobenzene-d	-----ISTD-----						
2) SURR1	1,4-Dioxane-d8	0.344	0.304	0.333	0.240	0.237	0.292	17.40
3)	1,4-Dioxane	0.558	0.516	0.530	0.327	0.330	0.452	25.15
4) I	Naphthalene-d8	-----ISTD-----						
5)	Naphthalene	1.086	1.055	1.196	0.864	0.982	1.037	11.90
6) SURR2	Methylnaphthalene	0.510	0.496	0.585	0.422	0.483	0.499	11.70
7)	2-Methylnaphthalene	0.695	0.675	0.801	0.589	0.678	0.688	11.02
8) I	Acenaphthene-d10	-----ISTD-----						
9)	Acenaphthylene	1.884	1.927	2.406	1.677	2.142	2.007	13.83
10) C	Acenaphthene	1.547	1.569	1.924	1.317	1.623	1.596	13.63
11)	Fluorene	1.663	1.720	2.270	1.509	1.795	1.791	16.05
12) I	Phenanthrene-d10	-----ISTD-----						
13)	Pentachlorophenol		0.020	0.035	0.040	0.054	0.037	37.93
14)	Phenanthrene	1.213	1.186	1.385	0.994	1.121	1.180	12.08
15)	Anthracene	1.057	1.001	1.178	0.875	1.029	1.028	10.60
16) SURR	Fluoranthene-d10	0.817	0.692	0.891	0.604	0.669	0.735	15.85
17) C	Fluoranthene	1.289	1.061	1.419	0.979	1.149	1.179	14.96
18) I	Chrysene-d12	-----ISTD-----						
19)	Pyrene	1.699	1.741	2.226	1.436	1.713	1.763	16.25
20)	Benzo(a)anthracene	1.204	1.135	1.343	0.964	1.133	1.156	11.87
21)	Chrysene	1.521	1.367	1.642	1.081	1.228	1.368	16.36
22) I	Perylene-d12	-----ISTD-----						
23)	Benzo(b)fluoranthene	1.363	1.292	1.633	1.103	1.394	1.357	14.12
24)	Benzo(k)fluoranthene	1.551	1.517	1.831	1.186	1.441	1.505	15.40
25) C	Benzo(a)pyrene	1.430	1.412	1.547	1.050	1.331	1.354	13.79
26)	Indeno(1,2,3-cd)pyr	1.585	1.422	1.677	1.168	1.397	1.450	13.51
27)	Dibenzo(a,h)anthrac	1.274	1.112	1.321	0.928	1.121	1.151	13.47
28)	Benzo(g,h,i)perylene	1.406	1.217	1.490	1.010	1.213	1.267	14.81

(#) = Out of Range