

Method Path : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\
 Method File : SOM-EPA-SIM-BM091019.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Tue Sep 10 15:06:45 2019
 Response Via : Initial Calibration

Calibration Files

0.1 =BM022614.D 0.2 =BM022615.D 0.4 =BM022616.D 0.8 =BM022617.D 1.6 =BM022618.D 3.2 =BM022619.D

Compound	0.1	0.2	0.4	0.8	1.6	3.2	Avg	%RSD

1) I 1,4-Dichlorobenzen...	-----ISTD-----							
2) I Naphthalene-d8	-----ISTD-----							
3) Naphthalene	1.189	1.150	1.152	1.081	1.127		1.140	3.49
4) SURR2-Methylnaphth...	0.688	0.648	0.665	0.620	0.658		0.656	3.80
5) 2-Methylnaphth...	0.820	0.773	0.794	0.737	0.783		0.781	3.88
6) I Acenaphthene-d10	-----ISTD-----							
7) Acenaphthylene	1.878	1.821	1.822	1.774	1.964		1.852	3.94
8) C Acenaphthene	1.596	1.563	1.548	1.478	1.566		1.550	2.85
9) Fluorene	1.813	1.740	1.767	1.702	1.857		1.776	3.42
10) I Phenanthrene-d10	-----ISTD-----							
11) Pentachlorophenol		0.071	0.079	0.075	0.088	0.093	0.081	11.16
12) Phenanthrene	1.318	1.269	1.269	1.222	1.293		1.274	2.78
13) Anthracene	1.103	1.047	1.075	1.063	1.179		1.093	4.75
14) SURRFluoranthene-d10	1.142	1.059	1.057	1.067	1.165		1.098	4.68
15) C Fluoranthene	1.421	1.319	1.329	1.346	1.461		1.375	4.53
16) I Chrysene-d12	-----ISTD-----							
17) Pyrene	1.865	1.853	1.894	1.695	1.761		1.813	4.58
18) Benzo(a)anthra...	1.578	1.455	1.504	1.487	1.630		1.531	4.66
19) Chrysene	1.701	1.603	1.633	1.521	1.596		1.611	4.04
20) I Perylene-d12	-----ISTD-----							
21) Benzo(b)fluora...	1.723	1.691	1.752	1.694	1.831		1.738	3.30
22) Benzo(k)fluora...	1.810	1.786	1.779	1.687	1.803		1.773	2.80
23) C Benzo(a)pyrene	1.540	1.481	1.517	1.467	1.599		1.521	3.43
24) Indeno(1,2,3-c...	1.876	1.815	1.971	1.777	1.836		1.855	3.99
25) Dibenzo(a,h)an...	1.496	1.470	1.597	1.441	1.491		1.499	3.92
26) Benzo(g,h,i)pe...	1.480	1.451	1.533	1.429	1.497		1.478	2.71

(#) = Out of Range