

Method Path : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\
 Method File : SOM-EPA-SIM-BE082918.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Aug 30 02:12:09 2018
 Response Via : Initial Calibration

Calibration Files

0.1 =BE097381.D 0.2 =BE097382.D 0.4 =BE097388.D
 0.8 =BE097384.D 1.6 =BE097385.D 3.2 =BE097386.D

	Compound	0.1	0.2	0.4	0.8	1.6	3.2	Avg	%RSD

1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2) I	Naphthalene-d8	-----ISTD-----							
3)	Naphthalene	0.981	0.988	0.967	0.956	0.985		0.975	1.40
4) SURR2	Methylnaphthale	0.583	0.624	0.627	0.621	0.654		0.622	4.05
5)	2-Methylnaphthale	0.574	0.602	0.616	0.627	0.663		0.616	5.29
6) I	Acenaphthene-d10	-----ISTD-----							
7)	Acenaphthylene	1.384	1.433	1.439	1.472	1.558		1.457	4.44
8) C	Acenaphthene	1.254	1.255	1.242	1.238	1.277		1.253	1.20
9)	Fluorene	1.356	1.413	1.430	1.463	1.514		1.435	4.10
10) I	Phenanthrene-d10	-----ISTD-----							
11)	Pentachlorophenol		0.033	0.032	0.037	0.045	0.050	0.039	20.12
12)	Phenanthrene	1.010	1.032	1.029	1.022	1.057		1.030	1.68
13)	Anthracene	0.795	0.827	0.851	0.886	0.970		0.866	7.75
14) SURR	Fluoranthene-d10	1.314	1.301	1.282	1.323	1.377		1.319	2.70
15) C	Fluoranthene	1.288	1.296	1.292	1.336	1.401		1.323	3.63
16) I	Chrysene-d12	-----ISTD-----							
17)	Pyrene	1.165	1.107	1.067	1.002	1.041		1.076	5.83
18)	Benzo(a)anthracen	1.000	1.011	1.004	1.016	1.108		1.028	4.40
19)	Chrysene	1.201	1.202	1.178	1.158	1.180		1.184	1.54
20) I	Perylene-d12	-----ISTD-----							
21)	Benzo(b)fluoranth	0.976	1.076	1.013	1.103	1.188		1.071	7.68
22)	Benzo(k)fluoranth	1.160	1.159	1.200	1.194	1.260		1.195	3.46
23) C	Benzo(a)pyrene	0.944	0.961	0.950	1.010	1.091		0.991	6.24
24)	Indeno(1,2,3-cd)p	1.244	1.250	1.276	1.281	1.371		1.284	3.96
25)	Dibenzo(a,h)anthr	1.041	1.056	1.087	1.091	1.166		1.088	4.45
26)	Benzo(g,h,i)peryl	1.054	1.074	1.076	1.104	1.168		1.095	4.06

(#) = Out of Range