

Method Path : Z:\HPCHEM1\BNA M\METHODS\
 Method File : SOM-EPA-SIM-BM070716.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Thu Jul 07 05:51:09 2016
 Response Via : Initial Calibration

Calibration Files

0.1 =BM006280.D 0.2 =BM006281.D 0.4 =BM006299.D
 0.8 =BM006283.D 1.6 =BM006284.D 3.2 =BM006285.D

	Compound	0.1	0.2	0.4	0.8	1.6	3.2	Avg	%RSD

1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2) SURR	1,4-Dioxane-d8	0.169	0.163	0.185	0.160	0.169		0.169	5.85
3)	1,4-Dioxane	0.236	0.217	0.234	0.198	0.209		0.219	7.34
4) I	Naphthalene-d8	-----ISTD-----							
5)	Naphthalene	0.885	0.907	1.017	0.901	0.955		0.933	5.77
6) SURR	2-Methylnaphthale	0.408	0.450	0.510	0.471	0.500		0.468	8.75
7)	2-Methylnaphthale	0.539	0.602	0.682	0.631	0.669		0.625	9.15
8) I	Acenaphthene-d10	-----ISTD-----							
9)	Acenaphthylene	1.238	1.303	1.439	1.308	1.450		1.347	6.89
10) C	Acenaphthene	1.250	1.300	1.457	1.288	1.390		1.337	6.32
11)	Fluorene	1.366	1.449	1.603	1.488	1.596		1.500	6.69
12) I	Phenanthrene-d10	-----ISTD-----							
13)	Pentachlorophenol		0.048	0.057	0.054	0.068	0.081	0.062	21.27
14)	Phenanthrene	1.066	1.072	1.188	1.056	1.150		1.106	5.33
15)	Anthracene	0.953	0.958	1.108	1.008	1.101		1.025	7.33
16) SURR	Fluoranthene-d10	0.711	0.832	0.905	0.876	0.891		0.843	9.34
17) C	Fluoranthene	1.034	1.199	1.301	1.267	1.298		1.220	9.16
18) I	Chrysene-d12	-----ISTD-----							
19)	Pyrene	1.617	1.666	1.754	1.592	1.746		1.675	4.37
20)	Benzo(a)anthracen	1.236	1.254	1.406	1.247	1.331		1.295	5.61
21)	Chrysene	1.241	1.244	1.381	1.226	1.253		1.269	5.01
22) I	Perylene-d12	-----ISTD-----							
23)	Benzo(b)fluoranth	1.421	1.283	1.487	1.494	1.477		1.432	6.19
24)	Benzo(k)fluoranth	1.174	1.338	1.501	1.271	1.378		1.332	9.15
25) C	Benzo(a)pyrene	1.276	1.292	1.454	1.278	1.363		1.333	5.74
26)	Indeno(1,2,3-cd)p	1.468	1.427	1.672	1.313	1.516		1.479	8.87
27)	Dibenzo(a,h)anthr	1.109	1.132	1.300	1.040	1.189		1.154	8.45
28)	Benzo(q,h,i)peryl	1.206	1.233	1.409	1.120	1.286		1.251	8.55

(#) = Out of Range