

Method Path : Z:\HPCHEM1\BNA_E\METHODS\
 Method File : SIM-BE030316.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Fri Mar 04 11:58:27 2016
 Response Via : Initial Calibration

Calibration Files

0.1 =BE091423.D 0.2 =BE091424.D 0.4 =BE091425.D
 0.8 =BE091426.D 1.6 =BE091427.D 3.2 =BE091428.D

	Compound	0.1	0.2	0.4	0.8	1.6	3.2	Avg	%RSD
1) I	1,4-Dichlorobenzene-d	-----ISTD-----							
2)	1,4-Dioxane	0.456	0.403	0.390	0.431	0.483	0.494	0.443	9.57
3)	n-Nitrosodimethyl	0.038	0.234	0.427	0.514	0.617	0.953	0.464	68.27
4) I	Naphthalene-d8	-----ISTD-----							
5) S	Nitrobenzene-d5	0.166	0.202	0.233	0.244	0.290	0.311	0.241	22.33
6)	Naphthalene	1.164	1.145	1.234	1.099	1.186	1.158	1.164	3.84
7)	2-Methylnaphthale	0.647	0.621	0.669	0.652	0.729	0.747	0.677	7.34
8) I	Acenaphthene-d10	-----ISTD-----							
9) S	2-Fluorobiphenyl	1.068	1.111	1.181	1.208	1.314	1.402	1.214	10.30
10)	Acenaphthylene	0.765	0.737	0.749	0.778	0.907	1.098	0.839	E1 16.84
11)	Acenaphthene	1.636	1.441	1.376	1.384	1.512	1.698	1.508	8.88
12)	Fluorene	1.739	1.778	1.731	1.937	2.227	2.206	1.936	11.84
13) I	Phenanthrene-d10	-----ISTD-----							
14)	Hexachlorobenzene	0.385	0.328	0.271	0.232	0.255	0.307	0.296	18.81
15)	Pentachlorophenol	0.062	0.018	0.012	0.024	0.061	0.135	0.052	88.92
16)	Phenanthrene	1.863	1.690	1.687	1.789	1.835	1.691	1.759	4.55
17)	Anthracene	1.573	1.526	1.631	2.587	3.606	4.156	2.513	45.48
18)	Fluoranthene	1.805	1.927	1.987	2.617	3.427	3.595	2.559	30.89
19) I	Chrysene-d12	-----ISTD-----							
20)	Pyrene	1.339	1.318	1.396	1.322	1.553	1.620	1.425	9.15
21) S	Terphenyl-d14	0.729	0.722	0.748	0.751	0.776	0.838	0.761	5.57
22)	Benzo(a)anthracen	1.599	1.522	1.559	1.504	2.026	2.108	1.720	15.83
23)	Chrysene	1.929	1.872	1.901	1.784	1.769	1.864	1.853	3.44
24)	Indeno(1,2,3-cd)p	1.428	1.469	1.439	1.436	1.613	1.666	1.509	6.89
25) I	Perylene-d12	-----ISTD-----							
26)	Benzo(b)fluoranth	1.420	1.143	1.215	1.233	1.570	1.787	1.395	17.79
27)	Benzo(k)fluoranth	1.083	1.487	1.474	1.423	1.360	1.283	1.351	11.24
28) C	Benzo(a)pyrene	1.118	1.125	1.130	1.098	1.298	1.453	1.204	11.81
29)	Dibenzo(a,h)anthr	1.052	1.059	1.094	1.101	1.255	1.352	1.152	10.67
30)	Benzo(g,h,i)peryl	1.180	1.176	1.207	1.199	1.333	1.384	1.247	7.13

(#) = Out of Range