

Method Path : Z:\HPCHEM1\BNA_E\METHODS\
 Method File : 8270-PAH-BE080216.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 Last Update : Tue Aug 02 16:59:37 2016
 Response Via : Initial Calibration

Calibration Files

0.1 =BE092106.D 0.2 =BE092107.D 0.4 =BE092108.D 0.8 =BE092109.D 1.6 =BE092110.D 3.2 =BE092111.D 5 =BE092112.D
 10 =BE092113.D

	Compound		0.1	0.2	0.4	0.8	1.6	3.2	5	10	Avg	%RSD

1) I	1,4-Dichlorobenzen...	-----ISTD-----										
2)	1,4-Dioxane		0.833	0.783	0.705	0.696	0.646	0.575	0.623	0.581	0.680	13.62
3)	n-Nitrosodimet...		1.517	0.788	0.758	0.817	0.882	0.861	0.956	0.959	0.942	25.82
4) S	2-Fluorophenol		1.670	1.234	1.135	1.187	1.225	1.173	1.323	1.326	1.284	13.24
5) S	Phenol-d6		1.802	1.573	1.520	1.632	1.636	1.881	1.932	1.711		9.35
6) I	Naphthalene-d8	-----ISTD-----										
7) S	Nitrobenzene-d5		0.506	0.451	0.376	0.337	0.356	0.347	0.391	0.393	0.395	14.50
8)	Naphthalene		1.572	1.212	1.121	1.107	1.083	1.005	1.101	1.071	1.159	15.24
9)	2-Methylnaphth...		1.014	0.724	0.691	0.698	0.706	0.666	0.730	0.711	0.743	15.03
10) I	Acenaphthene-d10	-----ISTD-----										
11) S	2,4,6-Tribromo...		0.149	0.136	0.154	0.166	0.170	0.199		0.162		13.33
12) S	2-Fluorobiphenyl		1.952	1.596	1.515	1.550	1.535	1.448	1.573	1.507	1.584	9.80
13)	Acenaphthylene		1.047	0.835	0.830	0.859	0.872	0.825	0.906	0.883	0.882	E1 8.22
14)	Acenaphthene		1.710	1.486	1.367	1.362	1.325	1.235	1.358	1.404	1.406	10.06
15)	Fluorene		2.409	1.745	1.681	1.721	1.719	1.587	1.739	1.663	1.783	14.47
16) I	Phenanthrene-d10	-----ISTD-----										
17)	Hexachlorobenzene		0.296	0.244	0.247	0.249	0.228	0.225	0.236	0.232	0.245	9.23
18)	Pentachlorophenol		0.058	0.055	0.061	0.075	0.091			0.068		21.77
19)	Phenanthrene		1.996	1.455	1.389	1.378	1.371	1.318	1.439	1.339	1.460	15.13
20)	Anthracene		1.564	1.169	1.246	1.216	1.302	1.340	1.423	1.388	1.331	9.57
21)	Fluoranthene		7.575	5.913	5.875	5.869	5.772	5.757	6.200	5.964	6.116	9.90
22) I	Chrysene-d12	-----ISTD-----										
23)	Pyrene		6.384	5.270	4.836	5.016	5.017	4.883	5.267	5.155	5.228	9.44
24) S	Terphenyl-d14		0.733	0.657	0.707	0.714	0.692	0.772	0.751	0.718		5.31
25)	Benzo(a)anthra...		7.095	5.474	5.131	5.163	5.261	5.025	5.475	5.345	5.496	12.11
26)	Chrysene		7.328	5.601	5.223	5.389	5.248	4.967	5.216	5.070	5.505	13.82
27)	Indeno(1,2,3-c...		7.396	6.068	5.530	5.746	5.786	5.489	5.955	5.856	5.978	10.12
28) I	Perylene-d12	-----ISTD-----										
29)	Benzo(b)fluora...		1.793	1.466	1.394	1.233	1.256	1.192	1.340	1.267	1.368	14.21
30)	Benzo(k)fluora...		1.892	1.255	1.190	1.347	1.349	1.259	1.362	1.298	1.369	16.02
31) C	Benzo(a)pyrene		1.448	1.247	1.208	1.198	1.214	1.189	1.309	1.247	1.257	6.83
32)	Dibenzo(a,h)an...		1.407	1.164	1.132	1.144	1.183	1.126	1.248	1.193	1.200	7.73

Method Path : Z:\HPCHEM1\BNA_E\METHODS\

Method File : 8270-PAH-BE080216.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

33)	Benzo(g,h,i)pe...	6.258	5.121	4.885	4.928	4.977	4.687	5.176	4.946	5.122	9.42
-----	-------------------	-------	-------	-------	-------	-------	-------	-------	-------	-------	------

(#) = Out of Range