Response Factor Report gcms1

Method Path : Z:\HPCHEM1\BNA_M\METHODS\
Method File : SOM01.2-EPA-SIM-BM021115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
Last Update : Wed Feb 11 14:38:48 2015
Response Via : Initial Calibration

Calibration Files
0.1 =BM000476.D 0.2 =BM000477.D 0.4 =BM000478.D
0.8 =BM000479.D 1 =BM000480.D

		Compound	0.1	0.2	0.4	0.8	1	Avg	%RSD
1)	I	1,4-Dichlorobenzene-d	----------ISTD----------						
2)	I	Naphthalene-d8	----------ISTD----------						
3)		Naphthalene	1.148	0.936	0.931	0.945	0.923	0.977	9.82
4)	SUR	2-Methylnaphthalene	0.590	0.518	0.525	0.531	0.514	0.536	5.78
5)		2-Methylnaphthalene	0.795	0.646	0.657	0.669	0.649	0.683	9.25
6)	I	Acenaphthene-d10	----------ISTD----------						
7)		Acenaphthylene	1.950	1.483	1.440	1.607	1.505	1.597	12.92
8)	C	Acenaphthene	1.480	1.143	1.123	1.229	1.143	1.224	12.18
9)		Fluorene	1.817	1.390	1.382	1.480	1.398	1.493	12.38
10)	I	Phenanthrene-d10	----------ISTD----------						
11)		Pentachlorophenol		0.032	0.035	0.041	0.040	0.037	11.18
12)		Phenanthrene	1.330	1.066	1.085	1.095	1.069	1.129	10.02
13)		Anthracene	1.150	0.951	0.978	0.999	0.991	1.014	7.72
14)	SUR	Fluoranthene-d10	0.988	0.861	0.852	0.874	0.857	0.886	6.51
15)	C	Fluoranthene	1.529	1.226	1.222	1.275	1.261	1.303	9.86
16)	I	Chrysene-d12	----------ISTD----------						
17)		Pyrene	1.723	1.391	1.436	1.406	1.381	1.468	9.85
18)		Benzo(a)anthracene	1.417	1.098	1.101	1.130	1.130	1.175	11.57
19)		Chrysene	1.517	1.325	1.325	1.309	1.301	1.356	6.70
20)	I	Perylene-d12	----------ISTD----------						
21)		Benzo(b)fluoranthen	1.539	1.283	1.343	1.377	1.346	1.377	7.01
22)		Benzo(k)fluoranthen	1.666	1.393	1.441	1.428	1.392	1.464	7.85
23)	C	Benzo(a)pyrene	1.496	1.230	1.253	1.261	1.239	1.296	8.67
24)		Indeno(1,2,3-cd)pyr	1.603	1.331	1.270	1.294	1.315	1.363	10.00
25)		Dibenzo(a,h)anthrac	1.254	1.060	0.996	1.030	1.050	1.078	9.41
26)		Benzo(g,h,i)perylen	1.390	1.174	1.102	1.111	1.128	1.181	10.18

(#) = Out of Range