Response Factor Report gcms1

Method Path : Z:\HPCHEM1\BNA M\METHODS\
Method File : SOM-EPA-SIM-BM121916.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
Last Update : Tue Dec 20 04:48:59 2016
Response Via : Initial Calibration

Calibration Files
0.1 =BM008440.D 0.2 =BM008441.D 0.4 =BM008460.D
0.8 =BM008443.D 1.6 =BM008444.D 3.2 =BM008445.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	1.226	1.145	1.099	1.090	1.067		1.125	5.58
4)	SURR	2-Methylnaphthale	0.683	0.634	0.583	0.591	0.576		0.613	7.34
5)		2-Methylnaphthale	0.824	0.767	0.730	0.740	0.727		0.757	5.30
6)	I	Acenaphthene-d10	ISTD							
7)		Acenaphthylene	1.873	1.818	1.744	1.854	1.759		1.809	3.15
8)	C	Acenaphthene	1.381	1.365	1.332	1.385	1.341		1.361	1.73
9)		Fluorene	1.612	1.547	1.531	1.575	1.502		1.553	2.70
10)	I	Phenanthrene-d10	ISTD							
11)		Pentachlorophenol		0.117	0.135	0.114	0.119	0.137	0.124	8.70
12)		Phenanthrene	1.227	1.223	1.209	1.227	1.221		1.221	0.61
13)		Anthracene	1.203	1.191	1.137	1.196	1.204		1.186	2.35
14)	SURR	Fluoranthene-d10	1.152	1.091	1.094	1.089	1.080		1.101	2.62
15)	C	Fluoranthene	1.570	1.510	1.466	1.475	1.466		1.498	2.97
16)	I	Chrysene-d12	ISTD							
17)		Pyrene	1.616	1.535	1.521	1.547	1.512		1.546	2.66
18)		Benzo(a)anthracen	1.474	1.239	1.247	1.339	1.268		1.313	7.47
19)		Chrysene	1.525	1.484	1.460	1.349	1.372		1.438	5.20
20)	I	Perylene-d12	ISTD							
21)		Benzo(b)fluoranth	1.779	1.579	1.544	1.436	1.542		1.576	7.97
22)		Benzo(k)fluoranth	1.514	1.629	1.548	1.553	1.439		1.537	4.49
23)	C	Benzo(a)pyrene	1.544	1.470	1.447	1.421	1.420		1.461	3.50
24)		Indeno(1,2,3-cd)p	1.779	1.744	1.768	1.689	1.724		1.741	2.05
25)		Dibenzo(a,h)anthr	1.422	1.398	1.414	1.344	1.376		1.391	2.29
26)		Benzo(g,h,i)peryl	1.485	1.515	1.527	1.454	1.483		1.493	1.93

(#) = Out of Range