

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
 Data File : BM009200.D
 Acq On : 11 Mar 2017 06:05
 Operator : SJ/MA
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.446

Quant Time: Mar 13 12:39:33 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 07 03:37:41 2017
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	277#	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	303#	0.00
3	Naphthalene	1.153	1.152	0.1	307#	0.00
4 SURR	2-Methylnaphthalene-d10	0.664	0.639	3.8	298#	0.00
5	2-Methylnaphthalene	0.829	0.794	4.2	299#	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	288#	0.00
7	Acenaphthylene	2.082	2.296	-10.3	330#	0.00
8 C	Acenaphthene	1.598	1.564	2.1	295#	0.00
9	Fluorene	1.869	1.834	1.9	279#	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	276#	0.00
11	Pentachlorophenol	0.194	0.150	22.7#	202#	0.00
12	Phenanthrene	1.256	1.290	-2.7	290#	0.00
13	Anthracene	1.241	1.236	0.4	276#	0.00
14 SURR	Fluoranthene-d10	1.251	1.230	1.7	272#	0.00
15 C	Fluoranthene	1.638	1.555	5.1	261#	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	242#	0.00
17	Pyrene	1.376	1.457	-5.9	270#	0.00
18	Benzo(a)anthracene	1.329	1.488	-12.0	267#	0.00
19	Chrysene	1.261	1.373	-8.9	268#	0.00
20 I	Perylene-d12	1.000	1.000	0.0	233#	0.00
21	Benzo(b)fluoranthene	1.571	1.590	-1.2	246#	0.00
22	Benzo(k)fluoranthene	1.389	1.392	-0.2	247#	0.00
23 C	Benzo(a)pyrene	1.414	1.444	-2.1	242#	0.00
24	Indeno(1,2,3-cd)pyrene	1.798	1.630	9.3	206#	0.00
25	Dibenzo(a,h)anthracene	1.510	1.316	12.8	194#	0.00
26	Benzo(g,h,i)perylene	1.534	1.331	13.2	194#	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0