

Data Path : Z:\HPCHEM1\BNA_M\Data\BM030817\
 Data File : BM009106.D
 Acq On : 08 Mar 2017 13:17
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.448

Quant Time: Mar 08 15:08:05 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 08 05:48:55 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	137	0.02
2 I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
3	Naphthalene	1.153	1.146	0.6	138	0.01
4 SURR	2-Methylnaphthalene-d10	0.664	0.638	3.9	135	0.01
5	2-Methylnaphthalene	0.829	0.764	7.8	130	0.01
6 I	Acenaphthene-d10	1.000	1.000	0.0	130	0.00
7	Acenaphthylene	2.082	2.118	-1.7	138	0.00
8 C	Acenaphthene	1.598	1.547	3.2	132	0.00
9	Fluorene	1.869	1.780	4.8	123	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
11	Pentachlorophenol	0.194	0.167	13.9	103	0.00
12	Phenanthrene	1.256	1.238	1.4	128	0.00
13	Anthracene	1.241	1.187	4.4	121	0.00
14 SURR	Fluoranthene-d10	1.251	1.173	6.2	119	0.00
15 C	Fluoranthene	1.638	1.525	6.9	117	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	107	0.00
17	Pyrene	1.376	1.456	-5.8	119	0.00
18	Benzo(a)anthracene	1.329	1.312	1.3	104	0.00
19	Chrysene	1.261	1.255	0.5	108	0.00
20 I	Perylene-d12	1.000	1.000	0.0	96	0.00
21	Benzo(b)fluoranthene	1.571	1.471	6.4	94	0.00
22	Benzo(k)fluoranthene	1.389	1.427	-2.7	104	0.00
23 C	Benzo(a)pyrene	1.414	1.344	5.0	93	0.00
24	Indeno(1,2,3-cd)pyrene	1.798	1.610	10.5	84	0.00
25	Dibenzo(a,h)anthracene	1.510	1.319	12.6	81	0.00
26	Benzo(g,h,i)perylene	1.534	1.341	12.6	81	0.00

(#) = Out of Range

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