

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122716\
 Data File : BM008684.D
 Acq On : 27 Dec 2016 21:41
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.437

Quant Time: Dec 28 03:29:13 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 28 03:24:04 2016
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)

1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	-0.03
2 I	Naphthalene-d8	1.000	1.000	0.0	91	-0.01
3	Naphthalene	1.125	1.147	-2.0	95	-0.01
4 SURR	2-Methylnaphthalene-d10	0.613	0.636	-3.8	100	-0.02
5	2-Methylnaphthalene	0.757	0.759	-0.3	95	-0.02
6 I	Acenaphthene-d10	1.000	1.000	0.0	100	-0.02
7	Acenaphthylene	1.809	1.944	-7.5	111	0.00
8 C	Acenaphthene	1.361	1.387	-1.9	104	0.00
9	Fluorene	1.553	1.535	1.2	100	-0.02
10 I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
11	Pentachlorophenol	0.124	0.153	-23.4#	121	-0.02
12	Phenanthrene	1.221	1.166	4.5	103	-0.02
13	Anthracene	1.186	1.195	-0.8	112	-0.02
14 SURR	Fluoranthene-d10	1.101	1.129	-2.5	110	0.00
15 C	Fluoranthene	1.498	1.560	-4.1	114	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	116	-0.01
17	Pyrene	1.546	1.593	-3.0	122	-0.01
18	Benzo(a)anthracene	1.313	1.254	4.5	117	0.00
19	Chrysene	1.438	1.364	5.1	109	-0.01
20 I	Perylene-d12	1.000	1.000	0.0	108	0.00
21	Benzo(b)fluoranthene	1.576	1.432	9.1	100	0.00
22	Benzo(k)fluoranthene	1.537	1.537	0.0	107	0.00
23 C	Benzo(a)pyrene	1.461	1.476	-1.0	110	0.00
24	Indeno(1,2,3-cd)pyrene	1.741	1.747	-0.3	107	0.00
25	Dibenzo(a,h)anthracene	1.391	1.384	0.5	106	-0.01
26	Benzo(a,h,i)perylene	1.493	1.464	1.9	103	-0.01

(#) = Out of Range

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