

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
 Data File : BM008089.D
 Acq On : 28 Nov 2016 16:43
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.457

Quant Time: Nov 28 17:19:18 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 25 01:27:01 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	121	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	137	0.00
3	Naphthalene	1.119	1.097	2.0	133	0.00
4 SURR	2-Methylnaphthalene-d10	0.566	0.574	-1.4	139	0.00
5	2-Methylnaphthalene	0.721	0.737	-2.2	140	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	139	0.00
7	Acenaphthylene	1.810	1.867	-3.1	144	0.00
8 C	Acenaphthene	1.377	1.379	-0.1	143	0.00
9	Fluorene	1.583	1.587	-0.3	142	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	136	0.00
11	Pentachlorophenol	0.140	0.147	-5.0	148	0.00
12	Phenanthrene	1.246	1.196	4.0	132	0.00
13	Anthracene	1.170	1.226	-4.8	144	0.00
14 SURR	Fluoranthene-d10	1.132	1.148	-1.4	140	0.00
15 C	Fluoranthene	1.608	1.784	-10.9	156#	0.00
16 I	Chrysene-d12	1.000	1.000	0.0	122	0.00
17	Pyrene	1.389	1.636	-17.8	144	0.00
18	Benzo(a)anthracene	1.228	1.267	-3.2	130	0.00
19	Chrysene	1.527	1.182	22.6#	97	0.00
20 I	Perylene-d12	1.000	1.000	0.0	129	0.00
21	Benzo(b)fluoranthene	1.596	1.452	9.0	120	0.00
22	Benzo(k)fluoranthene	1.647	1.400	15.0	114	0.00
23 C	Benzo(a)pyrene	1.495	1.241	17.0	109	0.00
24	Indeno(1,2,3-cd)pyrene	1.848	1.507	18.5	105	0.00
25	Dibenzo(a,h)anthracene	1.481	1.296	12.5	114	0.00
26	Benzo(a,h,i)perylene	1.640	1.347	17.9	106	0.00

(#) = Out of Range

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