

Data Path : Z:\HPCHEM1\BNA M\DATA\BM070816\
 Data File : BM006322.D
 Acq On : 07 Jul 2016 16:45
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.444

Quant Time: Jul 08 03:05:57 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM070716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 08 01:14:08 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	81	0.00
2 SURR	1,4-Dioxane-d8	0.169	0.186	-10.1	81	0.00
3	1,4-Dioxane	0.219	0.233	-6.4	80	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	83	0.00
5	Naphthalene	0.933	1.014	-8.7	83	0.01
6 SURR	2-Methylnaphthalene-d10	0.468	0.545	-16.5	88	0.00
7	2-Methylnaphthalene	0.625	0.726	-16.2	88	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
9	Acenaphthylene	1.347	1.478	-9.7	98	0.00
10 C	Acenaphthene	1.337	1.461	-9.3	96	0.00
11	Fluorene	1.500	1.725	-15.0	103	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
13	Pentachlorophenol	0.062	0.064	-3.2	130	0.02
14	Phenanthrene	1.106	1.182	-6.9	115	0.00
15	Anthracene	1.025	1.117	-9.0	117	0.00
16 SURR	Fluoranthene-d10	0.843	1.032	-22.4#	132	0.00
17 C	Fluoranthene	1.220	1.488	-22.0#	132	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	130	0.00
19	Pyrene	1.675	1.811	-8.1	134	0.00
20	Benzo(a)anthracene	1.295	1.437	-11.0	133	0.00
21	Chrysene	1.269	1.392	-9.7	131	0.00
22 I	Perylene-d12	1.000	1.000	0.0	109	0.00
23	Benzo(b)fluoranthene	1.432	1.554	-8.5	114	0.00
24	Benzo(k)fluoranthene	1.332	1.566	-17.6	113	0.00
25 C	Benzo(a)pyrene	1.333	1.450	-8.8	108	0.00
26	Indeno(1,2,3-cd)pyrene	1.479	1.539	-4.1	100	0.00
27	Dibenzo(a,h)anthracene	1.154	1.211	-4.9	101	0.00
28	Benzo(q,h,i)perylene	1.251	1.276	-2.0	98	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1