

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121515\
 Data File : BM003572.D
 Acq On : 16 Dec 2015 06:04
 Operator : UM/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.454

Quant Time: Dec 16 18:31:41 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM121515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 16 05:49:57 2015
 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	67	0.00
2 SURR	1,4-Dioxane-d8	0.171	0.172	-0.6	67	-0.02
3	1,4-Dioxane	0.218	0.213	2.3	66	-0.02
4 I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
5	Naphthalene	0.901	0.926	-2.8	70	0.00
6 SURR	2-Methylnaphthalene-d10	0.479	0.504	-5.2	72	-0.01
7	2-Methylnaphthalene	0.605	0.650	-7.4	73	-0.01
8 I	Acenaphthene-d10	1.000	1.000	0.0	75	0.00
9	Acenaphthylene	1.528	1.565	-2.4	77	0.00
10 C	Acenaphthene	1.232	1.294	-5.0	77	0.00
11	Fluorene	1.365	1.534	-12.4	81	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
13	Pentachlorophenol	0.057	0.066	-15.8	103	0.00
14	Phenanthrene	1.032	1.091	-5.7	89	0.00
15	Anthracene	0.893	0.924	-3.5	88	0.00
16 SURR	Fluoranthene-d10	0.849	0.957	-12.7	94	0.00
17 C	Fluoranthene	1.120	1.295	-15.6	96	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	132	0.00
19	Pyrene	1.080	0.865	19.9	100	0.00
20	Benzo(a)anthracene	0.977	0.854	12.6	112	0.00
21	Chrysene	1.119	1.095	2.1	134	0.00
22 I	Perylene-d12	1.000	1.000	0.0	136	0.00
23	Benzo(b)fluoranthene	1.365	1.458	-6.8	141	0.00
24	Benzo(k)fluoranthene	1.237	1.269	-2.6	141	0.00
25 C	Benzo(a)pyrene	1.152	1.175	-2.0	138	0.00
26	Indeno(1,2,3-cd)pyrene	1.335	1.381	-3.4	150	0.00
27	Dibenzo(a,h)anthracene	1.075	1.099	-2.2	148	0.00
28	Benzo(q,h,i)perylene	1.162	1.206	-3.8	151#	0.00

(#) = Out of Range

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