

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004771.D
 Acq On : 25 Mar 2016 10:35
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.446

Quant Time: Mar 25 11:09:09 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 03:18:48 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	110	0.00
2 SURR	1,4-Dioxane-d8	2.253	2.468	-9.5	119	0.00
3	1,4-Dioxane	2.745	2.952	-7.5	116	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
5	Naphthalene	12.077	12.844	-6.4	118	0.00
6 SURR	2-Methylnaphthalene-d10	6.972	7.339	-5.3	117	0.00
7	2-Methylnaphthalene	8.652	9.188	-6.2	115	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
9	Acenaphthylene	23.000	23.973	-4.2	118	0.00
10 C	Acenaphthene	16.271	16.666	-2.4	115	0.00
11	Fluorene	19.428	19.858	-2.2	115	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
13	Pentachlorophenol	1.249	1.100	11.9	96	0.02
14	Phenanthrene	14.049	14.625	-4.1	113	0.00
15	Anthracene	13.273	13.948	-5.1	115	0.00
16 SURR	Fluoranthene-d10	11.819	12.271	-3.8	112	0.00
17 C	Fluoranthene	15.993	16.560	-3.5	113	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	103	0.00
19	Pyrene	17.122	18.650	-8.9	113	0.00
20	Benzo(a)anthracene	15.370	16.428	-6.9	110	0.00
21	Chrysene	14.847	15.500	-4.4	106	0.00
22 I	Perylene-d12	1.000	1.000	0.0	90	0.00
23	Benzo(b)fluoranthene	16.844	18.879	-12.1	103	0.00
24	Benzo(k)fluoranthene	15.612	16.405	-5.1	92	0.00
25 C	Benzo(a)pyrene	15.401	16.251	-5.5	94	0.00
26	Indeno(1,2,3-cd)pyrene	14.255	14.324	-0.5	90	0.00
27	Dibenzo(a,h)anthracene	11.221	11.271	-0.4	90	0.00
28	Benzo(q,h,i)perylene	11.798	11.730	0.6	89	0.00

(#) = Out of Range

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