

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032416\
 Data File : BM004756.D
 Acq On : 24 Mar 2016 18:56
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.434

Quant Time: Apr 05 18:05:47 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 16:23:17 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	86	0.00
2 SURR	1,4-Dioxane-d8	0.194	0.219	-12.9	94	0.00
3	1,4-Dioxane	0.236	0.263	-11.4	95	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
5	Naphthalene	0.943	1.013	-7.4	93	0.00
6 SURR	2-Methylnaphthalene-d10	0.456	0.501	-9.9	95	0.00
7	2-Methylnaphthalene	0.599	0.656	-9.5	95	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
9	Acenaphthylene	1.472	1.624	-10.3	102	0.00
10 C	Acenaphthene	1.320	1.402	-6.2	96	0.00
11	Fluorene	1.395	1.503	-7.7	99	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
13	Pentachlorophenol	0.054	0.066	-22.2#	140	0.00
14	Phenanthrene	1.103	1.154	-4.6	102	0.00
15	Anthracene	0.898	1.067	-18.8	120	0.00
16 SURR	Fluoranthene-d10	0.807	0.846	-4.8	105	0.00
17 C	Fluoranthene	1.134	1.208	-6.5	108	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
19	Pyrene	1.395	1.481	-6.2	108	0.00
20	Benzo(a)anthracene	0.947	1.029	-8.7	110	0.00
21	Chrysene	1.301	1.349	-3.7	111	0.00
22 I	Perylene-d12	1.000	1.000	0.0	105	0.00
23	Benzo(b)fluoranthene	1.431	1.502	-5.0	111	0.00
24	Benzo(k)fluoranthene	1.449	1.501	-3.6	109	0.00
25 C	Benzo(a)pyrene	1.204	1.276	-6.0	112	0.00
26	Indeno(1,2,3-cd)pyrene	1.322	1.379	-4.3	109	0.00
27	Dibenzo(a,h)anthracene	1.029	1.074	-4.4	109	0.00
28	Benzo(q,h,i)perylene	1.258	1.270	-1.0	106	0.00

(#) = Out of Range

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