

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020116\
 Data File : BM004163.D
 Acq On : 01 Feb 2016 17:02
 Operator : SJ/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.455

Quant Time: Feb 02 01:26:39 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM012916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 02 01:18:17 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	0.00
2 SURR	1,4-Dioxane-d8	0.201	0.209	-4.0	119	0.00
3	1,4-Dioxane	0.265	0.263	0.8	116	0.00
4 I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
5	Naphthalene	0.962	1.021	-6.1	121	0.00
6 SURR	2-Methylnaphthalene-d10	0.523	0.546	-4.4	119	-0.01
7	2-Methylnaphthalene	0.685	0.717	-4.7	119	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
9	Acenaphthylene	1.435	1.532	-6.8	118	-0.02
10 C	Acenaphthene	1.317	1.410	-7.1	117	0.00
11	Fluorene	1.594	1.667	-4.6	113	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
13	Pentachlorophenol	0.033	0.045	-36.4#	182#	0.00
14	Phenanthrene	1.091	1.165	-6.8	111	0.00
15	Anthracene	0.922	1.052	-14.1	120	0.00
16 SURR	Fluoranthene-d10	1.008	1.162	-15.3	120	0.00
17 C	Fluoranthene	1.365	1.551	-13.6	120	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
19	Pyrene	1.248	1.292	-3.5	109	0.00
20	Benzo(a)anthracene	1.020	1.097	-7.5	111	0.00
21	Chrysene	1.238	1.271	-2.7	108	0.00
22 I	Perylene-d12	1.000	1.000	0.0	101	0.00
23	Benzo(b)fluoranthene	1.327	1.399	-5.4	99	0.00
24	Benzo(k)fluoranthene	1.255	1.317	-4.9	110	0.00
25 C	Benzo(a)pyrene	1.164	1.247	-7.1	107	0.00
26	Indeno(1,2,3-cd)pyrene	1.278	1.316	-3.0	102	0.00
27	Dibenzo(a,h)anthracene	0.963	1.028	-6.7	106	0.00
28	Benzo(q,h,i)perylene	1.199	1.178	1.8	98	0.00

(#) = Out of Range

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