Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092916\
Data File : BM007632.D
Acq On : 29 Sep 2016 18:11
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.462

Quant Time: Sep 29 18:48:55 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 23 04:00:23 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	152#	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	151#	0.00
3		Naphthalene	0.931	1.040	-11.7	169#	0.00
4	SURR	2-Methylnaphthalene-d10	0.466	0.540	-15.9	174#	0.00
5		2-Methylnaphthalene	0.630	0.743	-17.9	178#	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	154#	0.00
7		Acenaphthylene	1.693	1.938	-14.5	178#	0.00
8	C	Acenaphthene	1.320	1.479	-12.0	174#	0.00
9		Fluorene	1.618	1.829	-13.0	170#	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	151#	0.00
11		Pentachlorophenol	0.109	0.120	-10.1	182#	0.00
12		Phenanthrene	1.082	1.192	-10.2	167#	0.00
13		Anthracene	0.981	1.089	-11.0	170#	0.00
14	SURR	Fluoranthene-d10	0.932	0.992	-6.4	162#	0.00
15	C	Fluoranthene	1.364	1.458	-6.9	164#	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	127	0.00
17		Pyrene	1.322	1.670	-26.3#	161#	0.00
18		Benzo(a)anthracene	1.182	1.348	-14.0	147	0.00
19		Chrysene	1.262	1.363	-8.0	134	0.00
20	I	Perylene-d12	1.000	1.000	0.0	109	0.00
21		Benzo(b)fluoranthene	1.335	1.701	-27.4#	137	0.00
22		Benzo(k)fluoranthene	1.349	1.501	-11.3	123	0.00
23	C	Benzo(a)pyrene	1.241	1.432	-15.4	127	0.00
24		Indeno(1,2,3-cd)pyrene	1.278	1.445	-13.1	125	0.00
25		Dibenzo(a,h)anthracene	0.976	1.144	-17.2	130	0.00
26		Benzo(g,h,i)perylene	1.148	1.270	-10.6	121	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0