Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM072916\
Data File : BM006734.D
Acq On : 29 Jul 2016 11:32
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.456

Quant Time: Jul 29 12:44:26 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM072816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 28 15:44:03 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	106	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
3		Naphthalene	1.007	1.023	-1.6	106	0.00
4	SURR	2-Methylnaphthalene-d10	0.507	0.524	-3.4	108	0.00
5		2-Methylnaphthalene	0.683	0.689	-0.9	106	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
7		Acenaphthylene	2.296	2.295	0.0	105	0.00
8	C	Acenaphthene	1.390	1.432	-3.0	107	0.00
9		Fluorene	1.630	1.663	-2.0	105	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
11		Pentachlorophenol	0.050	0.041	18.0	103	0.00
12		Phenanthrene	1.112	1.144	-2.9	105	0.00
13		Anthracene	1.063	1.072	-0.8	105	0.00
14	SURR	Fluoranthene-d10	1.063	1.109	-4.3	111	0.00
15	C	Fluoranthene	1.549	1.552	-0.2	106	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
17		Pyrene	1.210	1.417	-17.1	108	0.00
18		Benzo(a)anthracene	1.097	1.267	-15.5	106	0.00
19		Chrysene	1.098	1.358	-23.7#	116	0.00
20	I	Perylene-d12	1.000	1.000	0.0	115	0.00
21		Benzo(b)fluoranthene	1.368	1.413	-3.3	114	0.00
22		Benzo(k)fluoranthene	1.301	1.447	-11.2	110	0.00
23	C	Benzo(a)pyrene	1.267	1.395	-10.1	114	0.00
24		Indeno(1,2,3-cd)pyrene	1.598	1.655	-3.6	107	0.00
25		Dibenzo(a,h)anthracene	1.198	1.386	-15.7	118	0.00
26		Benzo(g,h,i)perylene	1.287	1.411	-9.6	113	-0.01

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