Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM092916\
Data File : BM007621.D
Acq On : 29 Sep 2016 10:29
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.461

Quant Time: Sep 29 14:46:56 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	139	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	137	0.01
3		Naphthalene	0.931	1.048	-12.6	155#	0.00
4	SURR	2-Methylnaphthalene-d10	0.466	0.544	-16.7	159#	0.00
5		2-Methylnaphthalene	0.630	0.751	-19.2	163#	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	139	0.00
7		Acenaphthylene	1.693	1.964	-16.0	163#	0.00
8	C	Acenaphthene	1.320	1.505	-14.0	160#	0.00
9		Fluorene	1.618	1.889	-16.7	160#	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	138	0.00
11		Pentachlorophenol	0.109	0.111	-1.8	155#	0.02
12		Phenanthrene	1.082	1.192	-10.2	153#	0.00
13		Anthracene	0.981	1.094	-11.5	156#	0.02
14	SURR	Fluoranthene-d10	0.932	0.970	-4.1	145	0.00
15	C	Fluoranthene	1.364	1.440	-5.6	148	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	110	0.00
17		Pyrene	1.322	1.743	-31.8#	145	0.00
18		Benzo(a)anthracene	1.182	1.296	-9.6	122	0.00
19		Chrysene	1.262	1.414	-12.0	120	0.00
20	I	Perylene-d12	1.000	1.000	0.0	94	0.00
21		Benzo(b)fluoranthene	1.335	1.588	-19.0	110	0.00
22		Benzo(k)fluoranthene	1.349	1.608	-19.2	113	0.00
23	C	Benzo(a)pyrene	1.241	1.408	-13.5	107	0.00
24		Indeno(1,2,3-cd)pyrene	1.278	1.451	-13.5	108	0.00
25		Dibenzo(a,h)anthracene	0.976	1.126	-15.4	111	0.00
26		Benzo(g,h,i)perylene	1.148	1.296	-12.9	107	0.00

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