Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM031017\
Data File : BM009167.D
Acq On : 10 Mar 2017 09:48
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.454

Quant Time: Mar 10 16:19:11 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 10 15:19:24 2017
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	171#	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	187#	0.00
3		Naphthalene	1.153	1.155	-0.2	190#	0.00
4	SURR	2-Methylnaphthalene-d10	0.664	0.614	7.5	177#	0.00
5		2-Methylnaphthalene	0.829	0.768	7.4	178#	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	166#	0.00
7		Acenaphthylene	2.082	2.211	-6.2	184#	0.00
8	C	Acenaphthene	1.598	1.639	-2.6	179#	0.00
9		Fluorene	1.869	1.853	0.9	163#	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	161#	0.00
11		Pentachlorophenol	0.194	0.168	13.4	132	0.00
12		Phenanthrene	1.256	1.220	2.9	160#	0.00
13		Anthracene	1.241	1.204	3.0	157#	0.00
14	SURR	Fluoranthene-d10	1.251	1.173	6.2	151#	0.00
15	C	Fluoranthene	1.638	1.508	7.9	148	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	134	0.00
17		Pyrene	1.376	1.447	-5.2	148	0.00
18		Benzo(a)anthracene	1.329	1.373	-3.3	136	0.00
19		Chrysene	1.261	1.315	-4.3	142	0.00
20	I	Perylene-d12	1.000	1.000	0.0	126	0.00
21		Benzo(b)fluoranthene	1.571	1.551	1.3	130	0.00
22		Benzo(k)fluoranthene	1.389	1.434	-3.2	138	0.00
23	C	Benzo(a)pyrene	1.414	1.453	-2.8	132	0.00
24		Indeno(1,2,3-cd)pyrene	1.798	1.781	0.9	122	0.00
25		Dibenzo(a,h)anthracene	1.510	1.485	1.7	119	0.00
26		Benzo(g,h,i)perylene	1.534	1.533	0.1	122	0.00

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