Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060217\
Data File : BM010393.D
Acq On : 02 Jun 2017 18:36
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.417

Quant Time: Jun 02 19:08:53 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM053017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 02 17:41:09 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	102	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
3		Naphthalene	1.055	1.034	2.0	103	0.00
4	SURR	2-Methylnaphthalene-d10	0.581	0.586	-0.9	105	0.00
5		2-Methylnaphthalene	0.744	0.750	-0.8	106	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
7		Acenaphthylene	1.997	2.028	-1.6	106	0.00
8	C	Acenaphthene	1.510	1.514	-0.3	106	0.00
9		Fluorene	1.808	1.767	2.3	103	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
11		Pentachlorophenol	0.153	0.170	-11.1	112	0.00
12		Phenanthrene	1.286	1.249	2.9	99	0.00
13		Anthracene	1.204	1.208	-0.3	102	0.00
14	SURR	Fluoranthene-d10	1.061	1.063	-0.2	99	0.00
15	C	Fluoranthene	1.595	1.579	1.0	100	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
17		Pyrene	1.704	1.673	1.8	100	0.00
18		Benzo(a)anthracene	1.626	1.636	-0.6	102	0.00
19		Chrysene	1.573	1.541	2.0	98	0.00
20	I	Perylene-d12	1.000	1.000	0.0	96	0.00
21		Benzo(b)fluoranthene	1.751	1.668	4.7	93	0.00
22		Benzo(k)fluoranthene	1.665	1.648	1.0	97	0.00
23	C	Benzo(a)pyrene	1.657	1.678	-1.3	99	0.00
24		Indeno(1,2,3-cd)pyrene	2.178	2.096	3.8	94	0.00
25		Dibenzo(a,h)anthracene	1.735	1.669	3.8	94	0.00
26		Benzo(g,h,i)perylene	1.878	1.772	5.6	92	0.00

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