Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
Data File : BM008080.D
Acq On : 24 Nov 2016 18:38
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.455

Quant Time: Nov 25 01:30:42 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 25 01:27:01 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	122	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	136	0.00
3		Naphthalene	1.119	1.106	1.2	133	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.580	-2.5	139	0.00
5		2-Methylnaphthalene	0.721	0.738	-2.4	139	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	140	0.00
7		Acenaphthylene	1.810	1.889	-4.4	147	0.00
8	C	Acenaphthene	1.377	1.349	2.0	141	0.00
9		Fluorene	1.583	1.552	2.0	140	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	139	0.00
11		Pentachlorophenol	0.140	0.135	3.6	139	0.00
12		Phenanthrene	1.246	1.232	1.1	139	0.00
13		Anthracene	1.170	1.186	-1.4	141	-0.02
14	SURR	Fluoranthene-d10	1.132	1.080	4.6	134	0.00
15	C	Fluoranthene	1.608	1.501	6.7	134	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	113	0.00
17		Pyrene	1.389	1.661	-19.6	136	0.00
18		Benzo(a)anthracene	1.228	1.453	-18.3	138	0.00
19		Chrysene	1.527	1.421	6.9	108	0.00
20	I	Perylene-d12	1.000	1.000	0.0	119	0.00
21		Benzo(b)fluoranthene	1.596	1.524	4.5	116	0.00
22		Benzo(k)fluoranthene	1.647	1.535	6.8	115	0.00
23	C	Benzo(a)pyrene	1.495	1.473	1.5	120	-0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.755	5.0	113	0.00
25		Dibenzo(a,h)anthracene	1.481	1.375	7.2	111	0.00
26		Benzo(g,h,i)perylene	1.640	1.507	8.1	109	0.00

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