Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112216\
Data File : BM008025.D
Acq On : 23 Nov 2016 09:28
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.450

Quant Time: Nov 23 12:08:05 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 23 01:26:48 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	100	0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
3		Naphthalene	1.119	1.123	-0.4	107	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.573	-1.2	109	0.01
5		2-Methylnaphthalene	0.721	0.741	-2.8	111	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.02
7		Acenaphthylene	1.810	1.867	-3.1	116	0.00
8	C	Acenaphthene	1.377	1.363	1.0	114	0.00
9		Fluorene	1.583	1.582	0.1	114	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
11		Pentachlorophenol	0.140	0.122	12.9	103	0.02
12		Phenanthrene	1.246	1.220	2.1	114	0.02
13		Anthracene	1.170	1.120	4.3	110	0.00
14	SURR	Fluoranthene-d10	1.132	1.063	6.1	109	0.00
15	C	Fluoranthene	1.608	1.563	2.8	115	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
17		Pyrene	1.389	1.595	-14.8	110	0.00
18		Benzo(a)anthracene	1.228	1.308	-6.5	105	0.00
19		Chrysene	1.527	1.510	1.1	97	0.00
20	I	Perylene-d12	1.000	1.000	0.0	97	0.00
21		Benzo(b)fluoranthene	1.596	1.624	-1.8	101	0.00
22		Benzo(k)fluoranthene	1.647	1.566	4.9	96	0.00
23	C	Benzo(a)pyrene	1.495	1.499	-0.3	99	0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.791	3.1	94	0.00
25		Dibenzo(a,h)anthracene	1.481	1.420	4.1	94	0.00
26		Benzo(g,h,i)perylene	1.640	1.585	3.4	93	0.00

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