Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122016\
Data File : BM008490.D
Acq On : 20 Dec 2016 21:42
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.451

Quant Time: Dec 21 12:51:37 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 02:43:00 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	117	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
3		Naphthalene	1.125	1.110	1.3	124	0.00
4	SURR	2-Methylnaphthalene-d10	0.613	0.572	6.7	121	0.00
5		2-Methylnaphthalene	0.757	0.705	6.9	119	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
7		Acenaphthylene	1.809	1.890	-4.5	125	0.00
8	C	Acenaphthene	1.361	1.373	-0.9	119	0.00
9		Fluorene	1.553	1.506	3.0	113	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
11		Pentachlorophenol	0.124	0.139	-12.1	113	0.00
12		Phenanthrene	1.221	1.175	3.8	107	0.00
13		Anthracene	1.186	1.174	1.0	114	0.00
14	SURR	Fluoranthene-d10	1.101	1.058	3.9	107	0.00
15	C	Fluoranthene	1.498	1.534	-2.4	115	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
17		Pyrene	1.546	1.686	-9.1	111	0.00
18		Benzo(a)anthracene	1.313	1.326	-1.0	106	0.00
19		Chrysene	1.438	1.440	-0.1	99	0.00
20	I	Perylene-d12	1.000	1.000	0.0	106	0.00
21		Benzo(b)fluoranthene	1.576	1.366	13.3	93	0.00
22		Benzo(k)fluoranthene	1.537	1.597	-3.9	109	0.00
23	C	Benzo(a)pyrene	1.461	1.469	-0.5	107	0.00
24		Indeno(1,2,3-cd)pyrene	1.741	1.801	-3.4	108	0.00
25		Dibenzo(a,h)anthracene	1.391	1.430	-2.8	107	0.00
26		Benzo(g,h,i)perylene	1.493	1.595	-6.8	110	0.00

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