Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122316\
Data File : BM008596.D
Acq On : 23 Dec 2016 15:32
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.462

Quant Time: Dec 23 16:36:50 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 23 15:11:17 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	129	-0.03
2	I	Naphthalene-d8	1.000	1.000	0.0	141	-0.01
3		Naphthalene	1.125	1.107	1.6	142	-0.01
4	SURR	2-Methylnaphthalene-d10	0.613	0.595	2.9	144	-0.02
5		2-Methylnaphthalene	0.757	0.720	4.9	139	-0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	139	-0.02
7		Acenaphthylene	1.809	1.867	-3.2	149	-0.02
8	C	Acenaphthene	1.361	1.371	-0.7	143	0.00
9		Fluorene	1.553	1.464	5.7	133	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
11		Pentachlorophenol	0.124	0.119	4.0	117	-0.02
12		Phenanthrene	1.221	1.197	2.0	132	0.00
13		Anthracene	1.186	1.260	-6.2	147	0.00
14	SURR	Fluoranthene-d10	1.101	1.121	-1.8	136	0.00
15	C	Fluoranthene	1.498	1.554	-3.7	141	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	134	0.00
17		Pyrene	1.546	1.608	-4.0	142	0.00
18		Benzo(a)anthracene	1.313	1.287	2.0	139	0.00
19		Chrysene	1.438	1.456	-1.3	134	0.00
20	I	Perylene-d12	1.000	1.000	0.0	140	0.00
21		Benzo(b)fluoranthene	1.576	1.455	7.7	132	-0.01
22		Benzo(k)fluoranthene	1.537	1.466	4.6	132	0.00
23	C	Benzo(a)pyrene	1.461	1.446	1.0	140	-0.01
24		Indeno(1,2,3-cd)pyrene	1.741	1.701	2.3	134	-0.01
25		Dibenzo(a,h)anthracene	1.391	1.351	2.9	133	-0.01
26		Benzo(g,h,i)perylene	1.493	1.482	0.7	135	-0.02

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