Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121916\
Data File : BM008475.D
Acq On : 20 Dec 2016 11:58
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.449

Quant Time: Dec 20 12:31:15 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 19 14:18:36 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	142	-0.05
2	I	Naphthalene-d8	1.000	1.000	0.0	141	-0.04
3		Naphthalene	1.125	1.096	2.6	141	-0.04
4	SURR	2-Methylnaphthalene-d10	0.613	0.569	7.2	138	-0.05
5		2-Methylnaphthalene	0.757	0.712	5.9	138	-0.05
6	I	Acenaphthene-d10	1.000	1.000	0.0	131	-0.02
7		Acenaphthylene	1.809	1.836	-1.5	138	-0.02
8	C	Acenaphthene	1.361	1.371	-0.7	134	-0.02
9		Fluorene	1.553	1.534	1.2	131	-0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	127	-0.03
11		Pentachlorophenol	0.124	0.175	-41.1#	164#	-0.05
12		Phenanthrene	1.221	1.189	2.6	125	-0.02
13		Anthracene	1.186	1.276	-7.6	142	-0.03
14	SURR	Fluoranthene-d10	1.101	1.127	-2.4	131	-0.02
15	C	Fluoranthene	1.498	1.488	0.7	129	-0.03
16	I	Chrysene-d12	1.000	1.000	0.0	144	-0.01
17		Pyrene	1.546	1.424	7.9	134	-0.03
18		Benzo(a)anthracene	1.313	1.220	7.1	141	-0.02
19		Chrysene	1.438	1.125	21.8#	111	-0.01
20	I	Perylene-d12	1.000	1.000	0.0	131	-0.02
21		Benzo(b)fluoranthene	1.576	1.452	7.9	123	-0.01
22		Benzo(k)fluoranthene	1.537	1.461	4.9	124	-0.02
23	C	Benzo(a)pyrene	1.461	1.420	2.8	129	-0.02
24		Indeno(1,2,3-cd)pyrene	1.741	1.753	-0.7	130	-0.04
25		Dibenzo(a,h)anthracene	1.391	1.422	-2.2	132	-0.03
26		Benzo(g,h,i)perylene	1.493	1.520	-1.8	131	-0.03

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