Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
Data File : BM008065.D
Acq On : 24 Nov 2016 09:43
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.454

Quant Time: Nov 25 00:23:50 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 24 04:11:15 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	121	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
3		Naphthalene	1.119	1.101	1.6	130	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.587	-3.7	138	0.00
5		2-Methylnaphthalene	0.721	0.750	-4.0	139	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	139	0.00
7		Acenaphthylene	1.810	1.897	-4.8	147	0.00
8	C	Acenaphthene	1.377	1.348	2.1	140	0.00
9		Fluorene	1.583	1.567	1.0	140	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	139	0.00
11		Pentachlorophenol	0.140	0.134	4.3	137	0.00
12		Phenanthrene	1.246	1.217	2.3	137	0.00
13		Anthracene	1.170	1.175	-0.4	140	0.00
14	SURR	Fluoranthene-d10	1.132	1.094	3.4	136	0.00
15	C	Fluoranthene	1.608	1.491	7.3	133	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	116	0.00
17		Pyrene	1.389	1.611	-16.0	135	0.00
18		Benzo(a)anthracene	1.228	1.402	-14.2	137	0.00
19		Chrysene	1.527	1.401	8.3	110	0.00
20	I	Perylene-d12	1.000	1.000	0.0	121	0.00
21		Benzo(b)fluoranthene	1.596	1.504	5.8	116	0.00
22		Benzo(k)fluoranthene	1.647	1.535	6.8	117	0.00
23	C	Benzo(a)pyrene	1.495	1.496	-0.1	124	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.741	5.8	114	0.00
25		Dibenzo(a,h)anthracene	1.481	1.374	7.2	113	0.00
26		Benzo(g,h,i)perylene	1.640	1.499	8.6	110	0.00

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