Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031017\
Data File : BM009191.D
Acq On : 11 Mar 2017 00:44
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.456

Quant Time: Mar 11 03:22:17 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM030317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 11 00:54:34 2017
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	253#	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	275#	0.00
3		Naphthalene	1.153	1.153	0.0	279#	0.00
4	SURR	2-Methylnaphthalene-d10	0.664	0.639	3.8	271#	0.00
5		2-Methylnaphthalene	0.829	0.804	3.0	275#	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	266#	0.00
7		Acenaphthylene	2.082	2.174	-4.4	290#	0.00
8	C	Acenaphthene	1.598	1.505	5.8	263#	0.00
9		Fluorene	1.869	1.745	6.6	246#	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	246#	0.00
11		Pentachlorophenol	0.194	0.144	25.8#	173#	0.00
12		Phenanthrene	1.256	1.232	1.9	247#	0.00
13		Anthracene	1.241	1.236	0.4	245#	0.00
14	SURR	Fluoranthene-d10	1.251	1.145	8.5	226#	0.00
15	C	Fluoranthene	1.638	1.442	12.0	216#	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	177#	0.00
17		Pyrene	1.376	1.638	-19.0	222#	0.00
18		Benzo(a)anthracene	1.329	1.438	-8.2	189#	0.00
19		Chrysene	1.261	1.352	-7.2	193#	0.00
20	I	Perylene-d12	1.000	1.000	0.0	162#	0.00
21		Benzo(b)fluoranthene	1.571	1.562	0.6	169#	0.00
22		Benzo(k)fluoranthene	1.389	1.427	-2.7	176#	0.00
23	C	Benzo(a)pyrene	1.414	1.468	-3.8	172#	0.00
24		Indeno(1,2,3-cd)pyrene	1.798	1.735	3.5	153#	0.00
25		Dibenzo(a,h)anthracene	1.510	1.435	5.0	148	0.00
26		Benzo(g,h,i)perylene	1.534	1.441	6.1	147	0.00

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