Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011117\
Data File : BM008768.D
Acq On : 10 Jan 2017 13:30
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.466

Quant Time: Jan 10 18:12:18 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM010917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 09 13:24:07 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	76	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	78	-0.01
3		Naphthalene	1.097	1.080	1.5	77	-0.01
4	SURR	2-Methylnaphthalene-d10	0.663	0.683	-3.0	80	0.00
5		2-Methylnaphthalene	0.796	0.803	-0.9	77	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
7		Acenaphthylene	1.763	1.772	-0.5	81	0.00
8	C	Acenaphthene	1.260	1.238	1.7	78	0.00
9		Fluorene	1.577	1.541	2.3	78	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	82	-0.02
11		Pentachlorophenol	0.184	0.191	-3.8	82	0.00
12		Phenanthrene	1.221	1.196	2.0	80	0.00
13		Anthracene	1.191	1.172	1.6	81	0.00
14	SURR	Fluoranthene-d10	1.180	1.224	-3.7	84	-0.01
15	C	Fluoranthene	1.529	1.521	0.5	81	-0.01
16	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
17		Pyrene	1.594	1.502	5.8	80	-0.01
18		Benzo(a)anthracene	1.483	1.431	3.5	83	-0.01
19		Chrysene	1.405	1.347	4.1	83	-0.01
20	I	Perylene-d12	1.000	1.000	0.0	88	-0.02
21		Benzo(b)fluoranthene	1.713	1.524	11.0	77	-0.01
22		Benzo(k)fluoranthene	1.495	1.489	0.4	90	-0.01
23	C	Benzo(a)pyrene	1.560	1.431	8.3	81	-0.01
24		Indeno(1,2,3-cd)pyrene	1.821	1.773	2.6	86	-0.02
25		Dibenzo(a,h)anthracene	1.449	1.428	1.4	85	-0.02
26		Benzo(g,h,i)perylene	1.588	1.517	4.5	85	-0.02

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