Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112816\
Data File : BM008082.D
Acq On : 28 Nov 2016 10:16
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.456

Quant Time: Nov 28 14:45:54 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 28 14:44:03 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	90	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
3		Naphthalene	1.119	1.091	2.5	95	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.576	-1.8	100	0.01
5		2-Methylnaphthalene	0.721	0.725	-0.6	99	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
7		Acenaphthylene	1.810	1.743	3.7	96	0.00
8	C	Acenaphthene	1.377	1.354	1.7	100	0.00
9		Fluorene	1.583	1.539	2.8	98	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
11		Pentachlorophenol	0.140	0.119	15.0	88	0.00
12		Phenanthrene	1.246	1.186	4.8	96	0.00
13		Anthracene	1.170	1.104	5.6	94	0.02
14	SURR	Fluoranthene-d10	1.132	1.054	6.9	94	-0.01
15	C	Fluoranthene	1.608	1.441	10.4	92	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	80	0.00
17		Pyrene	1.389	1.585	-14.1	91	0.00
18		Benzo(a)anthracene	1.228	1.288	-4.9	86	-0.01
19		Chrysene	1.527	1.391	8.9	75	0.00
20	I	Perylene-d12	1.000	1.000	0.0	82	-0.01
21		Benzo(b)fluoranthene	1.596	1.487	6.8	78	-0.01
22		Benzo(k)fluoranthene	1.647	1.511	8.3	78	-0.01
23	C	Benzo(a)pyrene	1.495	1.419	5.1	79	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.624	12.1	72	-0.01
25		Dibenzo(a,h)anthracene	1.481	1.230	16.9	69	-0.01
26		Benzo(g,h,i)perylene	1.640	1.495	8.8	74	-0.02

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