Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121016\
Data File : BM008238.D
Acq On : 10 Dec 2016 18:41
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.445

Quant Time: Dec 12 03:52:41 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 Dec 12 02:45:24 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	105	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
3		Naphthalene	1.119	1.084	3.1	108	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.558	1.4	110	0.00
5		2-Methylnaphthalene	0.721	0.683	5.3	106	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
7		Acenaphthylene	1.810	1.812	-0.1	104	0.00
8	C	Acenaphthene	1.377	1.320	4.1	102	0.00
9		Fluorene	1.583	1.406	11.2	94	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
11		Pentachlorophenol	0.140	0.133	5.0	88	0.00
12		Phenanthrene	1.246	1.206	3.2	87	0.00
13		Anthracene	1.170	1.137	2.8	87	0.00
14	SURR	Fluoranthene-d10	1.132	1.045	7.7	83	0.00
15	C	Fluoranthene	1.608	1.430	11.1	82	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	71	0.00
17		Pyrene	1.389	1.632	-17.5	84	0.00
18		Benzo(a)anthracene	1.228	1.337	-8.9	80	0.00
19		Chrysene	1.527	1.449	5.1	70	0.00
20	I	Perylene-d12	1.000	1.000	0.0	80	0.00
21		Benzo(b)fluoranthene	1.596	1.395	12.6	71	0.00
22		Benzo(k)fluoranthene	1.647	1.533	6.9	77	0.00
23	C	Benzo(a)pyrene	1.495	1.491	0.3	81	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.793	3.0	77	0.00
25		Dibenzo(a,h)anthracene	1.481	1.418	4.3	77	0.00
26		Benzo(g,h,i)perylene	1.640	1.530	6.7	74	0.00

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