Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112116\
Data File : BM007958.D
Acq On : 21 Nov 2016 15:10
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.444

Quant Time: Nov 21 15:49:27 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 21 15:11:37 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	92	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
3		Naphthalene	1.119	1.093	2.3	91	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.547	3.4	90	0.00
5		2-Methylnaphthalene	0.721	0.699	3.1	91	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
7		Acenaphthylene	1.810	1.822	-0.7	92	0.00
8	C	Acenaphthene	1.377	1.374	0.2	93	0.00
9		Fluorene	1.583	1.537	2.9	90	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
11		Pentachlorophenol	0.140	0.126	10.0	85	0.00
12		Phenanthrene	1.246	1.242	0.3	92	0.00
13		Anthracene	1.170	1.144	2.2	89	0.00
14	SURR	Fluoranthene-d10	1.132	1.101	2.7	90	0.00
15	C	Fluoranthene	1.608	1.522	5.3	89	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
17		Pyrene	1.389	1.375	1.0	88	0.00
18		Benzo(a)anthracene	1.228	1.227	0.1	92	0.00
19		Chrysene	1.527	1.425	6.7	86	0.00
20	I	Perylene-d12	1.000	1.000	0.0	86	0.00
21		Benzo(b)fluoranthene	1.596	1.496	6.3	82	0.00
22		Benzo(k)fluoranthene	1.647	1.693	-2.8	92	0.00
23	C	Benzo(a)pyrene	1.495	1.473	1.5	87	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.798	2.7	84	0.00
25		Dibenzo(a,h)anthracene	1.481	1.432	3.3	84	0.00
26		Benzo(g,h,i)perylene	1.640	1.610	1.8	84	0.00

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