Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM112216\
Data File : BM007986.D
Acq On : 22 Nov 2016 09:22
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.447

Quant Time: Nov 22 15:04:56 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 22 15:01:04 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	100	0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	103	0.01
3		Naphthalene	1.119	1.133	-1.3	104	0.01
4	SURR	2-Methylnaphthalene-d10	0.566	0.559	1.2	102	0.01
5		2-Methylnaphthalene	0.721	0.730	-1.2	105	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.02
7		Acenaphthylene	1.810	1.830	-1.1	107	0.00
8	C	Acenaphthene	1.377	1.394	-1.2	109	0.00
9		Fluorene	1.583	1.587	-0.3	107	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
11		Pentachlorophenol	0.140	0.118	15.7	91	0.02
12		Phenanthrene	1.246	1.234	1.0	105	0.02
13		Anthracene	1.170	1.146	2.1	103	0.02
14	SURR	Fluoranthene-d10	1.132	1.073	5.2	100	0.00
15	C	Fluoranthene	1.608	1.477	8.1	99	0.01
16	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
17		Pyrene	1.389	1.529	-10.1	102	0.01
18		Benzo(a)anthracene	1.228	1.258	-2.4	97	0.01
19		Chrysene	1.527	1.504	1.5	93	0.00
20	I	Perylene-d12	1.000	1.000	0.0	92	0.01
21		Benzo(b)fluoranthene	1.596	1.656	-3.8	97	0.01
22		Benzo(k)fluoranthene	1.647	1.552	5.8	90	0.01
23	C	Benzo(a)pyrene	1.495	1.476	1.3	93	0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.816	1.7	90	0.02
25		Dibenzo(a,h)anthracene	1.481	1.449	2.2	91	0.02
26		Benzo(g,h,i)perylene	1.640	1.603	2.3	90	0.01

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