Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112316\
Data File : BM008034.D
Acq On : 23 Nov 2016 15:11
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.452

Quant Time: Nov 24 02:56:49 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 23 01:26:48 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	93	0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	103	0.01
3		Naphthalene	1.119	1.120	-0.1	103	0.01
4	SURR	2-Methylnaphthalene-d10	0.566	0.572	-1.1	104	0.01
5		2-Methylnaphthalene	0.721	0.727	-0.8	104	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
7		Acenaphthylene	1.810	1.846	-2.0	109	0.02
8	C	Acenaphthene	1.377	1.358	1.4	107	0.02
9		Fluorene	1.583	1.577	0.4	107	0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
11		Pentachlorophenol	0.140	0.119	15.0	97	0.02
12		Phenanthrene	1.246	1.214	2.6	108	0.02
13		Anthracene	1.170	1.124	3.9	106	0.00
14	SURR	Fluoranthene-d10	1.132	1.049	7.3	103	0.01
15	C	Fluoranthene	1.608	1.453	9.6	103	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	90	0.01
17		Pyrene	1.389	1.590	-14.5	104	0.01
18		Benzo(a)anthracene	1.228	1.334	-8.6	101	0.01
19		Chrysene	1.527	1.433	6.2	87	0.01
20	I	Perylene-d12	1.000	1.000	0.0	91	0.01
21		Benzo(b)fluoranthene	1.596	1.523	4.6	88	0.01
22		Benzo(k)fluoranthene	1.647	1.672	-1.5	96	0.01
23	C	Benzo(a)pyrene	1.495	1.503	-0.5	93	0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.825	1.2	90	0.02
25		Dibenzo(a,h)anthracene	1.481	1.454	1.8	90	0.02
26		Benzo(g,h,i)perylene	1.640	1.617	1.4	89	0.02

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