Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE010715\
Data File : BE089348.D
Acq On : 8 Jan 2015 2:30
Operator : TP/IZ
Sample : SSTDCCC001
Misc : LOQ-WATER-1.0PPM-SIM
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Jan 08 15:40:42 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BASE-BE010715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 08 15:21:44 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	103	0.00
2	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
3	S	Nitrobenzene-d5	1.000	1.047	-4.7	103	0.00
4		Naphthalene	1.000	1.031	-3.1	102	0.00
5		2-Methylnaphthalene	1.000	1.017	-1.7	101	0.00
6	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
7	S	2-Fluorobiphenyl	1.000	1.039	-3.9	101	0.00
8		Acenaphthylene	1.000	1.024	-2.4	99	0.00
9	C	Acenaphthene	1.000	1.010	-1.0	101	0.00
10		Fluorene	1.000	1.013	-1.3	101	0.00
11	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
12		Phenanthrene	1.000	1.037	-3.7	100	0.00
13		Anthracene	1.000	1.046	-4.6	99	0.00
14	C	Fluoranthene	1.000	1.057	-5.7	101	0.00
15	I	Chrysene-d12	5.000	5.000	0.0	100	0.00
16		Pyrene	1.000	1.016	-1.6	101	0.00
17	S	Terphenyl-d14	1.000	0.993	0.7	101	0.00
18		Benzo(a)anthracene	1.000	1.011	-1.1	102	0.00
19		Chrysene	1.000	1.031	-3.1	99	0.00
20		Indeno(1,2,3-cd)pyrene	1.000	0.965	3.5	87	0.00
21	I	Perylene-d12	5.000	5.000	0.0	99	0.00
22		Benzo(b)fluoranthene	1.000	1.016	-1.6	98	0.00
23		Benzo(k)fluoranthene	1.000	1.046	-4.6	99	0.00
24	C	Benzo(a)pyrene	1.000	1.014	-1.4	99	0.00
25		Dibenzo(a,h)anthracene	1.000	1.013	-1.3	96	0.00
26		Benzo(g,h,i)perylene	1.000	1.018	-1.8	98	0.00

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