Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE103114\
Data File : BE088433.D
Acq On : 31 Oct 2014 21:29
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 31 23:47:57 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE103114.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Nov 01 00:42:07 2014
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	100	0.00
2		1,4-Dioxane	1.000	1.043	-4.3	100	0.00
3	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
4	S	Nitrobenzene-d5	1.000	1.038	-3.8	99	0.00
5		Naphthalene	1.000	1.028	-2.8	98	0.00
6		2-Methylnaphthalene	1.000	1.040	-4.0	98	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
8	S	2-Fluorobiphenyl	1.000	1.036	-3.6	97	0.00
9		Acenaphthylene	1.000	1.026	-2.6	98	0.00
10		Acenaphthene	1.000	1.009	-0.9	97	0.00
11		Fluorene	1.000	1.030	-3.0	95	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
13		Phenanthrene	1.000	1.133	-13.3	93	0.00
14		Anthracene	1.000	1.041	-4.1	95	0.00
15		Fluoranthene	1.000	1.032	-3.2	92	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
17		Pyrene	1.000	1.034	-3.4	91	0.00
18	S	Terphenyl-d14	1.000	1.029	-2.9	92	0.00
19		Benzo(a)anthracene	1.000	1.024	-2.4	93	0.00
20		Chrysene	1.000	1.040	-4.0	94	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	1.030	-3.0	95	0.00
22	I	Perylene-d12	5.000	5.000	0.0	96	0.00
23		Benzo(b)fluoranthene	1.000	1.036	-3.6	97	0.00
24		Benzo(k)fluoranthene	1.000	1.030	-3.0	94	0.00
25	C	Benzo(a)pyrene	1.000	1.031	-3.1	96	0.00
26		Dibenzo(a,h)anthracene	1.000	1.029	-2.9	96	0.00
27		Benzo(g,h,i)perylene	1.000	1.027	-2.7	97	0.00

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