Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120914\
Data File : BE089189.D
Acq On : 9 Dec 2014 14:39
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Dec 09 15:35:35 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-ACID-BE120814.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 09 08:57:21 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	109	0.00
2	S	2-Fluorophenol	1.000	1.088	-8.8	109	0.00
3	S	Phenol-d6	1.000	1.069	-6.9	110	0.00
4		2-Chlorophenol	1.000	1.055	-5.5	109	0.00
5	C	Phenol	1.000	1.080	-8.0	111	0.00
6		2-Methylphenol	1.000	1.032	-3.2	104	0.00
7		3+4-Methylphenols	1.000	1.101	-10.1	114	0.01
8	I	Naphthalene-d8	5.000	5.000	0.0	111	0.00
9	C	2-Nitrophenol	1.000	1.036	-3.6	110	0.00
10		2,4-Dimethylphenol	1.000	1.018	-1.8	110	0.00
11	C	2,4-Dichlorophenol	1.000	0.995	0.5	109	0.00
12	C	4-Chloro-3-methylphenol	1.000	1.014	-1.4	110	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	112	0.00
14	S	2,4,6-Tribromophenol	1.000	1.022	-2.2	108	0.00
15	C	2,4,6-Trichlorophenol	1.000	1.006	-0.6	111	0.00
16		2,4,5-Trichlorophenol	1.000	1.033	-3.3	110	0.00
17	P	2,4-Dinitrophenol	1.000	0.900	10.0	118	0.01
18	P	4-Nitrophenol	1.000	1.078	-7.8	112	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
20		4,6-Dinitro-2-methylphenol	1.000	0.940	6.0	114	0.00
21	C	Pentachlorophenol	1.000	0.961	3.9	114	0.00
22	I	Chrysene-d12	5.000	5.000	0.0	108	0.00
23	I	Perylene-d12	5.000	5.000	0.0	106	0.00

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