Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021017\
Data File : BE092718.D
Acq On : 10 Feb 2017 12:21
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 10 14:46:07 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE011617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 17 10:14:01 2017
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	98	-0.02
2		1,4-Dioxane	0.400	0.370	7.5	88	0.00
3		n-Nitrosodimethylamine	0.400	0.296	26.0#	83	0.00
4	S	2-Fluorophenol	0.400	0.403	-0.8	116	-0.01
5	S	Phenol-d6	0.400	0.423	-5.7	126	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.330	17.5	96	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	93	-0.02
8	S	Nitrobenzene-d5	0.400	0.394	1.5	111	-0.02
9		Naphthalene	0.400	0.430	-7.5	108	-0.02
10		Hexachlorobutadiene	0.400	0.441	-10.2	108	-0.02
11		2-Methylnaphthalene	0.400	0.412	-3.0	107	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	91	-0.02
13	S	2,4,6-Tribromophenol	0.400	0.490	-22.5	123	-0.01
14	S	2-Fluorobiphenyl	0.400	0.430	-7.5	106	-0.02
15		Acenaphthylene	0.400	0.426	-6.5	110	-0.02
16		Acenaphthene	0.400	0.431	-7.7	104	-0.02
17		Fluorene	0.400	0.410	-2.5	105	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	81	0.00
19		Hexachlorobenzene	0.400	0.444	-11.0	88	-0.02
20		Pentachlorophenol	0.400	0.389	2.8	86	0.05
21		Phenanthrene	0.400	0.452	-13.0	95	-0.02
22		Anthracene	0.400	0.441	-10.2	100	-0.02
23		Fluoranthene	0.400	0.474	-18.5	106	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	95	-0.02
25		Pyrene	0.400	0.357	10.8	102	-0.02
26	S	Terphenyl-d14	0.400	0.414	-3.5	115	-0.02
27		Benzo(a)anthracene	0.400	0.405	-1.3	112	-0.02
28		Chrysene	0.400	0.370	7.5	95	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.465	-16.3	96	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.380	5.0	104	-0.03
31	I	Perylene-d12	5.000	5.000	0.0	83	-0.02
32		Benzo(b)fluoranthene	0.400	0.424	-6.0	96	-0.02
33		Benzo(k)fluoranthene	0.400	0.419	-4.7	101	-0.02
34	C	Benzo(a)pyrene	0.400	0.423	-5.7	102	-0.02
35		Dibenzo(a,h)anthracene	0.400	0.444	-11.0	104	-0.05
36		Benzo(g,h,i)perylene	0.400	0.449	-12.2	103	-0.05

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