Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE033017\
Data File : BE092876.D
Acq On : 30 Mar 2017 15:51
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 31 01:34:13 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 29 17:47:56 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	105	0.00
2		1,4-Dioxane	0.400	0.423	-5.7	108	-0.01
3		n-Nitrosodimethylamine	0.400	0.407	-1.7	115	0.00
4	S	2-Fluorophenol	0.400	0.348	13.0	93	0.00
5	S	Phenol-d6	0.400	0.360	10.0	101	0.00
6		bis(2-Chloroethyl)ether	0.400	0.414	-3.5	110	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.400	0.395	1.3	116	0.00
9		Naphthalene	0.400	0.413	-3.2	109	0.00
10		Hexachlorobutadiene	0.400	0.395	1.3	105	0.00
11		2-Methylnaphthalene	0.400	0.394	1.5	106	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	0.400	0.324	19.0	97	0.00
14	S	2-Fluorobiphenyl	0.400	0.407	-1.7	108	-0.01
15		Acenaphthylene	0.400	0.341	14.8	102	0.00
16		Acenaphthene	0.400	0.381	4.8	106	0.00
17		Fluorene	0.400	0.384	4.0	104	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	107	-0.02
19		Hexachlorobenzene	0.400	0.424	-6.0	116	0.00
20		Pentachlorophenol	0.400	0.462	-15.5	95	0.00
21		Phenanthrene	0.400	0.465	-16.3	121	0.00
22		Anthracene	0.400	0.412	-3.0	119	0.00
23		Fluoranthene	0.400	0.349	12.8	101	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	102	-0.02
25		Pyrene	0.400	0.364	9.0	99	0.00
26	S	Terphenyl-d14	0.400	0.373	6.8	99	-0.02
27		Benzo(a)anthracene	0.400	0.355	11.3	99	0.00
28		Chrysene	0.400	0.415	-3.7	107	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.306	23.5	80	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.328	18.0	89	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	101	0.00
32		Benzo(b)fluoranthene	0.400	0.330	17.5	87	-0.01
33		Benzo(k)fluoranthene	0.400	0.368	8.0	99	-0.01
34	C	Benzo(a)pyrene	0.400	0.316	21.0#	87	0.00
35		Dibenzo(a,h)anthracene	0.400	0.340	15.0	91	-0.02
36		Benzo(g,h,i)perylene	0.400	0.337	15.8	88	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1