Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE022017\
Data File : BE092753.D
Acq On : 20 Feb 2017 22:52
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 100 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 21 00:54:24 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE021517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 15 17:05:20 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	185	0.00
2		1,4-Dioxane	0.400	0.374	6.5	172	0.00
3		n-Nitrosodimethylamine	0.400	0.499	-24.7	199	-0.01
4	S	2-Fluorophenol	0.400	0.483	-20.7	232	-0.02
5	S	Phenol-d6	0.400	0.588	-47.0#	281	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.434	-8.5	218	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	195	-0.02
8	S	Nitrobenzene-d5	0.400	0.456	-14.0	234	-0.02
9		Naphthalene	0.400	0.408	-2.0	197	0.00
10		Hexachlorobutadiene	0.400	0.432	-8.0	202	0.00
11		2-Methylnaphthalene	0.400	0.400	0.0	200	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	201	0.00
13	S	2,4,6-Tribromophenol	0.400	0.518	-29.5#	266	-0.02
14	S	2-Fluorobiphenyl	0.400	0.426	-6.5	211	-0.02
15		Acenaphthylene	0.400	0.397	0.8	201	0.00
16		Acenaphthene	0.400	0.411	-2.7	205	-0.02
17		Fluorene	0.400	0.397	0.8	201	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	202	0.00
19		Hexachlorobenzene	0.400	0.383	4.3	176	0.00
20		Pentachlorophenol	0.400	0.553	-38.3#	246	-0.09
21		Phenanthrene	0.400	0.368	8.0	176	0.00
22		Anthracene	0.400	0.399	0.3	206	0.00
23		Fluoranthene	0.400	0.378	5.5	190	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	154	0.00
25		Pyrene	0.400	0.487	-21.7	193	0.00
26	S	Terphenyl-d14	0.400	0.517	-29.3#	205	-0.02
27		Benzo(a)anthracene	0.400	0.446	-11.5	175	0.00
28		Chrysene	0.400	0.415	-3.7	159	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.702	-75.5#	288	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.443	-10.7	146	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	158	-0.01
32		Benzo(b)fluoranthene	0.400	0.422	-5.5	148	-0.01
33		Benzo(k)fluoranthene	0.400	0.398	0.5	160	-0.01
34	C	Benzo(a)pyrene	0.400	0.390	2.5	163	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.389	2.8	143	-0.02
36		Benzo(g,h,i)perylene	0.400	0.383	4.3	138	-0.02

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