Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022817\
Data File : BE092772.D
Acq On : 28 Feb 2017 20:48
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4EC

Quant Time: Mar 01 08:21:09 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	99	-0.02
2		1,4-Dioxane	0.400	0.454	-13.5	109	0.00
3		n-Nitrosodimethylamine	0.400	0.497	-24.2	105	-0.02
4	S	2-Fluorophenol	0.400	0.372	7.0	95	-0.02
5	S	Phenol-d6	0.400	0.371	7.3	95	0.00
6		bis(2-Chloroethyl)ether	0.400	0.360	10.0	97	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	94	-0.02
8	S	Nitrobenzene-d5	0.400	0.417	-4.2	101	-0.02
9		Naphthalene	0.400	0.402	-0.5	94	0.00
10		Hexachlorobutadiene	0.400	0.397	0.8	90	0.00
11		2-Methylnaphthalene	0.400	0.399	0.3	96	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	108	-0.02
13	S	2,4,6-Tribromophenol	0.400	0.451	-12.7	124	-0.02
14	S	2-Fluorobiphenyl	0.400	0.370	7.5	98	-0.01
15		Acenaphthylene	0.400	0.387	3.3	105	-0.02
16		Acenaphthene	0.400	0.387	3.3	103	-0.02
17		Fluorene	0.400	0.412	-3.0	112	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
19		Hexachlorobenzene	0.400	0.378	5.5	103	-0.02
20		Pentachlorophenol	0.400	0.506	-26.5#	124	-0.09
21		Phenanthrene	0.400	0.367	8.3	105	0.00
22		Anthracene	0.400	0.412	-3.0	127	0.00
23		Fluoranthene	0.400	0.430	-7.5	129	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	138	-0.02
25		Pyrene	0.400	0.358	10.5	127	0.00
26	S	Terphenyl-d14	0.400	0.370	7.5	132	-0.02
27		Benzo(a)anthracene	0.400	0.375	6.3	132	-0.02
28		Chrysene	0.400	0.330	17.5	114	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.283	29.3#	90	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.332	17.0	96	-0.03
31	I	Perylene-d12	5.000	5.000	0.0	117	-0.02
32		Benzo(b)fluoranthene	0.400	0.399	0.3	104	-0.02
33		Benzo(k)fluoranthene	0.400	0.394	1.5	118	-0.02
34	C	Benzo(a)pyrene	0.400	0.359	10.3	112	-0.02
35		Dibenzo(a,h)anthracene	0.400	0.348	13.0	95	-0.03
36		Benzo(g,h,i)perylene	0.400	0.372	7.0	100	-0.03

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