Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013117\
Data File : BE092689.D
Acq On : 31 Jan 2017 16:27
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 01 10:34:49 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	156	0.00
2		1,4-Dioxane	0.400	0.354	11.5	134	0.00
3		n-Nitrosodimethylamine	0.400	0.251	37.3#	112	0.02
4	S	2-Fluorophenol	0.400	0.360	10.0	165	0.00
5	S	Phenol-d6	0.400	0.391	2.3	185	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.355	11.3	164	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	162	0.00
8	S	Nitrobenzene-d5	0.400	0.349	12.8	170	-0.02
9		Naphthalene	0.400	0.388	3.0	170	-0.02
10		Hexachlorobutadiene	0.400	0.416	-4.0	177	-0.02
11		2-Methylnaphthalene	0.400	0.381	4.8	172	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	159	0.00
13	S	2,4,6-Tribromophenol	0.400	0.458	-14.5	195	0.02
14	S	2-Fluorobiphenyl	0.400	0.396	1.0	170	-0.01
15		Acenaphthylene	0.400	0.401	-0.3	180	-0.02
16		Acenaphthene	0.400	0.403	-0.8	169	0.00
17		Fluorene	0.400	0.388	3.0	172	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	166	0.00
19		Hexachlorobenzene	0.400	0.330	17.5	143	0.00
20		Pentachlorophenol	0.400	0.297	25.8#	135	0.28
21		Phenanthrene	0.400	0.375	6.3	161	0.00
22		Anthracene	0.400	0.389	2.8	181	-0.02
23		Fluoranthene	0.400	0.408	-2.0	187	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	151	0.00
25		Pyrene	0.400	0.421	-5.2	191	-0.02
26	S	Terphenyl-d14	0.400	0.408	-2.0	179	0.00
27		Benzo(a)anthracene	0.400	0.425	-6.2	186	0.00
28		Chrysene	0.400	0.436	-9.0	177	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.643	-60.7#	256	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.439	-9.7	189	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	162	0.00
32		Benzo(b)fluoranthene	0.400	0.400	0.0	176	-0.01
33		Benzo(k)fluoranthene	0.400	0.421	-5.2	198	-0.01
34	C	Benzo(a)pyrene	0.400	0.417	-4.2	195	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.414	-3.5	189	-0.03
36		Benzo(g,h,i)perylene	0.400	0.412	-3.0	184	-0.02

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