Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE121916\
Data File : BE092615.D
Acq On : 19 Dec 2016 19:46
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 20 01:01:23 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 19 16:16:32 2016
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	104	0.00
2		1,4-Dioxane	0.400	0.389	2.8	113	0.00
3		n-Nitrosodimethylamine	0.400	0.450	-12.5	121	0.00
4	S	2-Fluorophenol	0.400	0.427	-6.7	110	0.00
5	S	Phenol-d6	0.400	0.416	-4.0	112	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.360	10.0	108	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	102	0.00
8	S	Nitrobenzene-d5	0.400	0.448	-12.0	110	0.00
9		Naphthalene	0.400	0.376	6.0	104	0.00
10		Hexachlorobutadiene	0.400	0.377	5.8	102	0.00
11		2-Methylnaphthalene	0.400	0.351	12.3	100	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	0.400	0.416	-4.0	99	0.00
14	S	2-Fluorobiphenyl	0.400	0.370	7.5	101	0.00
15		Acenaphthylene	0.400	0.362	9.5	101	0.00
16		Acenaphthene	0.400	0.377	5.8	101	0.00
17		Fluorene	0.400	0.364	9.0	100	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
19		Hexachlorobenzene	0.400	0.373	6.8	100	0.00
20		Pentachlorophenol	0.400	0.477	-19.2	133	0.00
21		Phenanthrene	0.400	0.357	10.8	100	0.00
22		Anthracene	0.400	0.344	14.0	101	0.00
23		Fluoranthene	0.400	0.356	11.0	101	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
25		Pyrene	0.400	0.370	7.5	100	0.00
26	S	Terphenyl-d14	0.400	0.374	6.5	100	-0.02
27		Benzo(a)anthracene	0.400	0.368	8.0	102	0.00
28		Chrysene	0.400	0.389	2.8	101	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.421	-5.2	111	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.364	9.0	101	0.00
31	I	Perylene-d12	5.000	5.000	0.0	100	0.00
32		Benzo(b)fluoranthene	0.400	0.354	11.5	101	0.00
33		Benzo(k)fluoranthene	0.400	0.351	12.3	101	0.00
34	C	Benzo(a)pyrene	0.400	0.337	15.8	100	0.00
35		Dibenzo(a,h)anthracene	0.400	0.345	13.8	102	0.00
36		Benzo(g,h,i)perylene	0.400	0.357	10.8	101	-0.02

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