Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020117\
Data File : BE092716.D
Acq On : 1 Feb 2017 15:44
Operator : SJ/MA
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 02 09:53:01 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	154	0.00
2		1,4-Dioxane	0.400	0.385	3.8	143	0.00
3		n-Nitrosodimethylamine	0.400	0.344	14.0	152	0.00
4	S	2-Fluorophenol	0.400	0.430	-7.5	195	-0.01
5	S	Phenol-d6	0.400	0.479	-19.7	225	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.363	9.3	167	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	161	0.00
8	S	Nitrobenzene-d5	0.400	0.386	3.5	187	-0.02
9		Naphthalene	0.400	0.402	-0.5	175	-0.02
10		Hexachlorobutadiene	0.400	0.431	-7.7	183	-0.02
11		2-Methylnaphthalene	0.400	0.402	-0.5	181	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	156	0.00
13	S	2,4,6-Tribromophenol	0.400	0.506	-26.5#	219	-0.01
14	S	2-Fluorobiphenyl	0.400	0.412	-3.0	173	-0.02
15		Acenaphthylene	0.400	0.421	-5.2	185	-0.02
16		Acenaphthene	0.400	0.430	-7.5	177	-0.02
17		Fluorene	0.400	0.409	-2.2	178	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	158	0.00
19		Hexachlorobenzene	0.400	0.327	18.3	135	0.00
20		Pentachlorophenol	0.400	0.488	-22.0	211	-0.02
21		Phenanthrene	0.400	0.379	5.3	155	0.00
22		Anthracene	0.400	0.432	-8.0	192	-0.02
23		Fluoranthene	0.400	0.445	-11.2	194	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	133	0.00
25		Pyrene	0.400	0.482	-20.5	191	-0.02
26	S	Terphenyl-d14	0.400	0.461	-15.3	178	0.00
27		Benzo(a)anthracene	0.400	0.471	-17.7	181	0.00
28		Chrysene	0.400	0.414	-3.5	147	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.816	-104.0#	313	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	162	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	138	0.00
32		Benzo(b)fluoranthene	0.400	0.422	-5.5	158	-0.01
33		Benzo(k)fluoranthene	0.400	0.425	-6.2	171	-0.01
34	C	Benzo(a)pyrene	0.400	0.428	-7.0	171	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.422	-5.5	164	-0.02
36		Benzo(g,h,i)perylene	0.400	0.419	-4.7	160	-0.02

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