Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013117\
Data File : BE092705.D
Acq On : 1 Feb 2017 2:27
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 01 03:24:19 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	195	0.00
2		1,4-Dioxane	0.400	0.330	17.5	159	0.01
3		n-Nitrosodimethylamine	0.400	0.277	30.8#	154	0.01
4	S	2-Fluorophenol	0.400	0.405	-1.3	232	0.00
5	S	Phenol-d6	0.400	0.463	-15.8	274	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.342	14.5	199	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	206	0.00
8	S	Nitrobenzene-d5	0.400	0.371	7.3	231	-0.02
9		Naphthalene	0.400	0.394	1.5	220	-0.02
10		Hexachlorobutadiene	0.400	0.416	-4.0	226	-0.02
11		2-Methylnaphthalene	0.400	0.402	-0.5	232	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	214	0.00
13	S	2,4,6-Tribromophenol	0.400	0.483	-20.7	282	-0.01
14	S	2-Fluorobiphenyl	0.400	0.386	3.5	223	-0.02
15		Acenaphthylene	0.400	0.409	-2.2	247	-0.02
16		Acenaphthene	0.400	0.402	-0.5	227	0.00
17		Fluorene	0.400	0.389	2.8	232	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	208	0.00
19		Hexachlorobenzene	0.400	0.363	9.3	193	0.00
20		Pentachlorophenol	0.400	0.408	-2.0	232	0.00
21		Phenanthrene	0.400	0.398	0.5	214	0.00
22		Anthracene	0.400	0.417	-4.2	243	-0.02
23		Fluoranthene	0.400	0.407	-1.7	233	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	171	0.00
25		Pyrene	0.400	0.457	-14.2	234	-0.02
26	S	Terphenyl-d14	0.400	0.440	-10.0	219	0.00
27		Benzo(a)anthracene	0.400	0.430	-7.5	213	0.00
28		Chrysene	0.400	0.421	-5.2	193	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.744	-86.0#	356	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.400	0.0	195	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	167	0.00
32		Benzo(b)fluoranthene	0.400	0.420	-5.0	191	-0.01
33		Benzo(k)fluoranthene	0.400	0.425	-6.2	207	-0.01
34	C	Benzo(a)pyrene	0.400	0.425	-6.2	206	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.417	-4.2	197	-0.03
36		Benzo(g,h,i)perylene	0.400	0.407	-1.7	188	-0.02

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