Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101716\
Data File : BE092462.D
Acq On : 17 Oct 2016 20:59
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 18 01:05:21 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	149	-0.02
2		1,4-Dioxane	0.400	0.408	-2.0	159	-0.01
3		n-Nitrosodimethylamine	0.400	0.521	-30.3#	190	-0.02
4	S	2-Fluorophenol	0.400	0.449	-12.2	160	0.00
5	S	Phenol-d6	0.400	0.396	1.0	163	0.00
6		bis(2-Chloroethyl)ether	0.400	0.446	-11.5	187	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	140	0.00
8	S	Nitrobenzene-d5	0.400	0.481	-20.2	173	0.00
9		Naphthalene	0.400	0.448	-12.0	158	-0.02
10		Hexachlorobutadiene	0.400	0.428	-7.0	152	-0.02
11		2-Methylnaphthalene	0.400	0.444	-11.0	158	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	135	0.00
13	S	2,4,6-Tribromophenol	0.400	0.442	-10.5	139	0.00
14	S	2-Fluorobiphenyl	0.400	0.464	-16.0	159	0.00
15		Acenaphthylene	0.400	0.433	-8.2	150	0.00
16		Acenaphthene	0.400	0.450	-12.5	147	0.00
17		Fluorene	0.400	0.435	-8.7	149	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.02
19		Hexachlorobenzene	0.400	0.513	-28.2#	165	0.00
20		Pentachlorophenol	0.400	0.450	-12.5	156	0.00
21		Phenanthrene	0.400	0.517	-29.3#	164	0.00
22		Anthracene	0.400	0.468	-17.0	154	0.02
23		Fluoranthene	0.400	0.462	-15.5	151	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	135	0.00
25		Pyrene	0.400	0.485	-21.2	170	0.00
26	S	Terphenyl-d14	0.400	0.453	-13.2	155	0.00
27		Benzo(a)anthracene	0.400	0.450	-12.5	155	0.00
28		Chrysene	0.400	0.461	-15.3	145	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.435	-8.7	145	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.396	1.0	136	0.03
31	I	Perylene-d12	5.000	5.000	0.0	122	0.02
32		Benzo(b)fluoranthene	0.400	0.457	-14.2	140	0.02
33		Benzo(k)fluoranthene	0.400	0.453	-13.2	149	0.02
34	C	Benzo(a)pyrene	0.400	0.448	-12.0	145	0.02
35		Dibenzo(a,h)anthracene	0.400	0.430	-7.5	137	0.05
36		Benzo(g,h,i)perylene	0.400	0.429	-7.2	135	0.03

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