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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	149	-0.02
2		1,4-Dioxane	0.635	0.686	-8.0	159#	-0.01
3		n-Nitrosodimethylamine	0.588	0.679	-15.5	190#	-0.02
4	S	2-Fluorophenol	0.999	1.005	-0.6	160#	0.00
5	S	Phenol-d6	1.388	1.324	4.6	163#	0.00
6		bis(2-Chloroethyl)ether	1.222	1.197	2.0	187#	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	140	0.00
8	S	Nitrobenzene-d5	0.299	0.311	-4.0	173#	0.00
9		Naphthalene	1.082	1.211	-11.9	158#	-0.02
10		Hexachlorobutadiene	0.170	0.182	-7.1	152#	-0.02
11		2-Methylnaphthalene	0.716	0.795	-11.0	158#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	135	0.00
13	S	2,4,6-Tribromophenol	0.144	0.127	11.8	139	0.00
14	S	2-Fluorobiphenyl	1.184	1.373	-16.0	159#	0.00
15		Acenaphthylene	7.240	7.843	-8.3	150#	0.00
16		Acenaphthene	1.134	1.276	-12.5	147	0.00
17		Fluorene	1.438	1.563	-8.7	149	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	120	0.02
19		Hexachlorobenzene	0.214	0.275	-28.5#	165#	0.00
20		Pentachlorophenol	0.050	0.055	-10.0	156#	0.00
21		Phenanthrene	1.149	1.484	-29.2#	164#	0.00
22		Anthracene	1.105	1.292	-16.9	154#	0.02
23		Fluoranthene	5.460	6.306	-15.5	151#	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	135	0.00
25		Pyrene	4.333	5.259	-21.4	170#	0.00
26	S	Terphenyl-d14	0.620	0.702	-13.2	155#	0.00
27		Benzo(a)anthracene	4.873	5.483	-12.5	155#	0.00
28		Chrysene	4.462	5.139	-15.2	145	0.00
29		Bis(2-ethylhexyl)phthalate	0.420	0.456	-8.6	145	0.00
30		Indeno(1,2,3-cd)pyrene	4.734	4.683	1.1	136	0.03
31	I	Perylene-d12	1.000	1.000	0.0	122	0.02
32		Benzo(b)fluoranthene	1.211	1.382	-14.1	140	0.02
33		Benzo(k)fluoranthene	1.267	1.436	-13.3	149	0.02
34	C	Benzo(a)pyrene	1.187	1.329	-12.0	145	0.02
35		Dibenzo(a,h)anthracene	1.088	1.170	-7.5	137	0.05
36		Benzo(g,h,i)perylene	4.673	5.007	-7.1	135	0.03

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