Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE011917\
Data File : BE092658.D
Acq On : 19 Jan 2017 13:44
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 19 14:51:53 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	116	0.00
2		1,4-Dioxane	0.400	0.443	-10.7	122	0.01
3		n-Nitrosodimethylamine	0.400	0.350	12.5	117	0.02
4	S	2-Fluorophenol	0.400	0.340	15.0	116	0.02
5	S	Phenol-d6	0.400	0.337	15.8	119	0.00
6		bis(2-Chloroethyl)ether	0.400	0.344	14.0	119	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	117	0.00
8	S	Nitrobenzene-d5	0.400	0.346	13.5	122	0.00
9		Naphthalene	0.400	0.397	0.8	126	0.00
10		Hexachlorobutadiene	0.400	0.409	-2.2	126	0.00
11		2-Methylnaphthalene	0.400	0.382	4.5	125	0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	116	0.00
13	S	2,4,6-Tribromophenol	0.400	0.447	-11.7	137	0.00
14	S	2-Fluorobiphenyl	0.400	0.410	-2.5	128	0.00
15		Acenaphthylene	0.400	0.396	1.0	129	0.00
16		Acenaphthene	0.400	0.416	-4.0	127	0.00
17		Fluorene	0.400	0.391	2.3	126	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	115	0.00
19		Hexachlorobenzene	0.400	0.432	-8.0	123	0.00
20		Pentachlorophenol	0.400	0.325	18.8	103	0.00
21		Phenanthrene	0.400	0.427	-6.7	127	0.00
22		Anthracene	0.400	0.433	-8.2	140	0.00
23		Fluoranthene	0.400	0.401	-0.3	127	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	115	0.00
25		Pyrene	0.400	0.367	8.3	126	0.00
26	S	Terphenyl-d14	0.400	0.375	6.3	125	0.00
27		Benzo(a)anthracene	0.400	0.371	7.3	123	0.00
28		Chrysene	0.400	0.435	-8.7	134	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.431	-7.7	100	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.364	9.0	119	0.00
31	I	Perylene-d12	5.000	5.000	0.0	112	0.00
32		Benzo(b)fluoranthene	0.400	0.387	3.3	118	0.00
33		Benzo(k)fluoranthene	0.400	0.394	1.5	129	0.00
34	C	Benzo(a)pyrene	0.400	0.379	5.3	123	0.00
35		Dibenzo(a,h)anthracene	0.400	0.376	6.0	119	0.00
36		Benzo(g,h,i)perylene	0.400	0.382	4.5	119	0.00

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