Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE120916\
Data File : BE092589.D
Acq On : 9 Dec 2016 10:14
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 09 23:21:09 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 16 15:43:18 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	119	-0.02
2		1,4-Dioxane	0.400	0.434	-8.5	145	0.00
3		n-Nitrosodimethylamine	0.400	0.442	-10.5	122	0.01
4	S	2-Fluorophenol	0.400	0.331	17.3	99	0.00
5	S	Phenol-d6	0.400	0.394	1.5	118	0.00
6		bis(2-Chloroethyl)ether	0.400	0.330	17.5	112	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	119	0.00
8	S	Nitrobenzene-d5	0.400	0.392	2.0	116	0.00
9		Naphthalene	0.400	0.409	-2.2	122	-0.02
10		Hexachlorobutadiene	0.400	0.415	-3.7	123	0.00
11		2-Methylnaphthalene	0.400	0.392	2.0	123	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	125	0.00
13	S	2,4,6-Tribromophenol	0.400	0.330	17.5	110	0.00
14	S	2-Fluorobiphenyl	0.400	0.396	1.0	123	0.00
15		Acenaphthylene	0.400	0.382	4.5	124	0.00
16		Acenaphthene	0.400	0.405	-1.3	127	0.00
17		Fluorene	0.400	0.397	0.8	125	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	118	0.00
19		Hexachlorobenzene	0.400	0.419	-4.7	138	0.00
20		Pentachlorophenol	0.400	0.406	-1.5	128	0.00
21		Phenanthrene	0.400	0.427	-6.7	130	-0.02
22		Anthracene	0.400	0.395	1.3	124	0.00
23		Fluoranthene	0.400	0.395	1.3	124	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	115	0.00
25		Pyrene	0.400	0.393	1.8	121	0.00
26	S	Terphenyl-d14	0.400	0.396	1.0	123	0.00
27		Benzo(a)anthracene	0.400	0.444	-11.0	120	0.00
28		Chrysene	0.400	0.432	-8.0	120	-0.02
29		Bis(2-ethylhexyl)phthalate	0.400	0.452	-13.0	124	-0.02
30		Indeno(1,2,3-cd)pyrene	0.400	0.397	0.8	105	0.00
31	I	Perylene-d12	5.000	5.000	0.0	110	-0.01
32		Benzo(b)fluoranthene	0.400	0.353	11.8	101	-0.01
33		Benzo(k)fluoranthene	0.400	0.399	0.3	116	-0.01
34	C	Benzo(a)pyrene	0.400	0.377	5.8	112	0.00
35		Dibenzo(a,h)anthracene	0.400	0.354	11.5	104	-0.02
36		Benzo(g,h,i)perylene	0.400	0.364	9.0	104	0.00

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