Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE030617\
Data File : BE092803.D
Acq On : 7 Mar 2017 1:09
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 07 02:49:32 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE030617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 06 15:46:01 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	105	-0.02
2		1,4-Dioxane	0.400	0.388	3.0	118	-0.01
3		n-Nitrosodimethylamine	0.400	0.364	9.0	110	-0.02
4	S	2-Fluorophenol	0.400	0.371	7.3	117	-0.02
5	S	Phenol-d6	0.400	0.355	11.3	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.372	7.0	111	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	111	-0.02
8	S	Nitrobenzene-d5	0.400	0.367	8.3	124	-0.02
9		Naphthalene	0.400	0.415	-3.7	126	0.00
10		Hexachlorobutadiene	0.400	0.422	-5.5	123	-0.02
11		2-Methylnaphthalene	0.400	0.402	-0.5	125	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	128	-0.02
13	S	2,4,6-Tribromophenol	0.400	0.363	9.3	118	-0.02
14	S	2-Fluorobiphenyl	0.400	0.401	-0.3	128	-0.02
15		Acenaphthylene	0.400	0.387	3.3	129	-0.02
16		Acenaphthene	0.400	0.410	-2.5	130	-0.02
17		Fluorene	0.400	0.407	-1.7	130	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
19		Hexachlorobenzene	0.400	0.372	7.0	131	-0.02
20		Pentachlorophenol	0.400	0.324	19.0	99	-0.12
21		Phenanthrene	0.400	0.390	2.5	133	0.00
22		Anthracene	0.400	0.373	6.8	131	0.00
23		Fluoranthene	0.400	0.383	4.3	131	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	107	-0.02
25		Pyrene	0.400	0.435	-8.7	130	0.00
26	S	Terphenyl-d14	0.400	0.443	-10.7	129	-0.02
27		Benzo(a)anthracene	0.400	0.426	-6.5	125	-0.02
28		Chrysene	0.400	0.477	-19.2	123	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.489	-22.2	137	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.417	-4.2	132	-0.03
31	I	Perylene-d12	5.000	5.000	0.0	116	-0.02
32		Benzo(b)fluoranthene	0.400	0.380	5.0	132	-0.02
33		Benzo(k)fluoranthene	0.400	0.429	-7.2	129	-0.02
34	C	Benzo(a)pyrene	0.400	0.409	-2.2	137	-0.02
35		Dibenzo(a,h)anthracene	0.400	0.381	4.8	130	-0.02
36		Benzo(g,h,i)perylene	0.400	0.392	2.0	128	-0.02

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