Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE020117\
Data File : BE092707.D
Acq On : 1 Feb 2017 9:58
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Feb 01 12:32:28 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	150	-0.02
2		1,4-Dioxane	0.400	0.375	6.3	136	0.00
3		n-Nitrosodimethylamine	0.400	0.391	2.3	168	-0.01
4	S	2-Fluorophenol	0.400	0.466	-16.5	206	-0.01
5	S	Phenol-d6	0.400	0.470	-17.5	215	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.394	1.5	176	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	145	0.00
8	S	Nitrobenzene-d5	0.400	0.418	-4.5	182	-0.02
9		Naphthalene	0.400	0.404	-1.0	158	-0.02
10		Hexachlorobutadiene	0.400	0.426	-6.5	162	-0.02
11		2-Methylnaphthalene	0.400	0.398	0.5	161	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	141	0.00
13	S	2,4,6-Tribromophenol	0.400	0.519	-29.8#	204	-0.01
14	S	2-Fluorobiphenyl	0.400	0.417	-4.2	158	-0.02
15		Acenaphthylene	0.400	0.410	-2.5	163	-0.02
16		Acenaphthene	0.400	0.407	-1.7	151	-0.02
17		Fluorene	0.400	0.396	1.0	155	-0.01
18	I	Phenanthrene-d10	5.000	5.000	0.0	135	0.00
19		Hexachlorobenzene	0.400	0.375	6.3	128	0.00
20		Pentachlorophenol	0.400	0.592	-48.0#	219	-0.02
21		Phenanthrene	0.400	0.388	3.0	135	0.00
22		Anthracene	0.400	0.424	-6.0	160	-0.02
23		Fluoranthene	0.400	0.420	-5.0	156	-0.02
24	I	Chrysene-d12	5.000	5.000	0.0	125	0.00
25		Pyrene	0.400	0.416	-4.0	156	-0.02
26	S	Terphenyl-d14	0.400	0.421	-5.2	154	0.00
27		Benzo(a)anthracene	0.400	0.458	-14.5	167	0.00
28		Chrysene	0.400	0.375	6.3	127	0.00
29		Bis(2-ethylhexyl)phthalate	0.400	0.815	-103.7#	296	0.00
30		Indeno(1,2,3-cd)pyrene	0.400	0.457	-14.2	164	-0.02
31	I	Perylene-d12	5.000	5.000	0.0	136	0.00
32		Benzo(b)fluoranthene	0.400	0.411	-2.7	152	-0.01
33		Benzo(k)fluoranthene	0.400	0.418	-4.5	165	-0.01
34	C	Benzo(a)pyrene	0.400	0.421	-5.2	165	-0.01
35		Dibenzo(a,h)anthracene	0.400	0.427	-6.7	164	-0.03
36		Benzo(g,h,i)perylene	0.400	0.422	-5.5	159	-0.02

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