Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101917\
Data File : BE094465.D
Acq On : 20 Oct 2017 3:01
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4EC

Quant Time: Oct 20 14:02:04 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 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	0.400	0.400	0.0	90	-0.02
2		1,4-Dioxane	0.400	0.405	-1.3	89	0.00
3		n-Nitrosodimethylamine	0.400	0.421	-5.2	95	0.00
4	S	2-Fluorophenol	0.400	0.440	-10.0	98	0.00
5	S	Phenol-d6	0.400	0.438	-9.5	96	0.00
6		bis(2-Chloroethyl)ether	0.400	0.434	-8.5	95	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	92	-0.02
8	S	Nitrobenzene-d5	0.400	0.412	-3.0	95	0.00
9		Naphthalene	0.400	0.431	-7.7	99	-0.02
10		Hexachlorobutadiene	0.400	0.404	-1.0	94	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.440	-10.0	101	0.00
12		2-Methylnaphthalene	0.400	0.444	-11.0	103	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	99	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.425	-6.2	98	0.00
15	S	2-Fluorobiphenyl	0.400	0.417	-4.2	100	0.00
16		Acenaphthylene	0.400	0.414	-3.5	102	0.00
17		Acenaphthene	0.400	0.423	-5.7	105	0.00
18		Fluorene	0.400	0.414	-3.5	102	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	102	-0.03
20		4-Bromophenyl-phenylether	0.400	0.457	-14.2	130	-0.03
21		Hexachlorobenzene	0.400	0.270	32.5#	74	0.00
22		Pentachlorophenol	0.400	0.473	-18.2	115	0.00
23		Phenanthrene	0.400	0.287	28.3#	71	0.00
24		Anthracene	0.400	0.410	-2.5	108	-0.03
25	SURR	Fluoranthene-d10	0.400	0.375	6.3	100	0.00
26		Fluoranthene	0.400	0.374	6.5	100	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	98	0.00
28		Pyrene	0.400	0.407	-1.7	100	0.00
29	S	Terphenyl-d14	0.400	0.413	-3.2	100	0.00
30		Benzo(a)anthracene	0.400	0.407	-1.7	100	-0.01
31		Chrysene	0.400	0.429	-7.2	103	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.368	8.0	93	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	101	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	100	-0.02
35		Benzo(b)fluoranthene	0.400	0.396	1.0	99	-0.01
36		Benzo(k)fluoranthene	0.400	0.424	-6.0	105	-0.01
37	C	Benzo(a)pyrene	0.400	0.409	-2.2	102	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.412	-3.0	102	-0.02
39		Benzo(g,h,i)perylene	0.400	0.409	-2.2	101	-0.03

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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