Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121319\
Data File : BE100895.D
Acq On : 13 Dec 2019 20:10
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 14 03:07:44 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 13 10:42:38 2019
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	87	0.00
2		1,4-Dioxane	0.400	0.442	-10.5	96	0.00
3		n-Nitrosodimethylamine	0.400	0.411	-2.7	103	0.00
4	S	2-Fluorophenol	0.400	0.487	-21.7	113	-0.01
5	S	Phenol-d6	0.400	0.471	-17.7	111	0.00
6		bis(2-Chloroethyl)ether	0.400	0.447	-11.7	103	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	89	0.00
8	S	Nitrobenzene-d5	0.400	0.598	-49.5#	153	0.00
9		Naphthalene	0.400	0.442	-10.5	102	0.00
10		Hexachlorobutadiene	0.400	0.434	-8.5	99	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.442	-10.5	104	0.00
12		2-Methylnaphthalene	0.400	0.444	-11.0	105	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	96	0.00
14	S	2,4,6-Tribromophenol	0.400	0.706	-76.5#	194	-0.01
15	S	2-Fluorobiphenyl	0.400	0.397	0.8	100	0.00
16		Acenaphthylene	0.400	0.444	-11.0	124	0.00
17		Acenaphthene	0.400	0.437	-9.2	115	0.00
18		Fluorene	0.400	0.435	-8.7	117	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	96	0.00
20		4-Bromophenyl-phenylether	0.400	0.415	-3.7	110	0.00
21		Hexachlorobenzene	0.400	0.423	-5.7	108	0.00
22		Pentachlorophenol	0.400	0.775	-93.8#	959	-0.01
23		Phenanthrene	0.400	0.419	-4.7	106	0.00
24		Anthracene	0.400	0.411	-2.7	115	0.00
25	SURR	Fluoranthene-d10	0.400	0.366	8.5	97	0.00
26		Fluoranthene	0.400	0.378	5.5	100	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	79	0.00
28		Pyrene	0.400	0.477	-19.2	100	0.00
29	S	Terphenyl-d14	0.400	0.423	-5.7	87	0.00
30		Benzo(a)anthracene	0.400	0.461	-15.3	102	0.00
31		Chrysene	0.400	0.449	-12.2	91	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.745	-86.3#	155	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.452	-13.0	101	0.00
34	I	Perylene-d12	0.400	0.400	0.0	85	0.00
35		Benzo(b)fluoranthene	0.400	0.415	-3.7	94	0.00
36		Benzo(k)fluoranthene	0.400	0.419	-4.7	94	0.00
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	96	0.00
38		Dibenzo(a,h)anthracene	0.400	0.413	-3.2	99	0.00
39		Benzo(g,h,i)perylene	0.400	0.412	-3.0	93	0.00

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

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