Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN100719\
Data File : BN008056.D
Acq On : 07 Oct 2019 19:27
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 08 00:24:15 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 03 18:39:49 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	40	-0.03
2		1,4-Dioxane	0.400	0.382	4.5	45	-0.04
3		n-Nitrosodimethylamine	0.400	0.742	-85.5#	49	-0.05
4	S	2-Fluorophenol	0.400	0.362	9.5	42	-0.02
5	S	Phenol-d6	0.400	0.372	7.0	41	0.00
6		bis(2-Chloroethyl)ether	0.400	0.420	-5.0	48	-0.04
7	I	Naphthalene-d8	0.400	0.400	0.0	42	-0.04
8	S	Nitrobenzene-d5	0.400	0.382	4.5	42	-0.03
9		Naphthalene	0.400	0.412	-3.0	44	-0.04
10		Hexachlorobutadiene	0.400	0.447	-11.7	47	-0.05
11	SURR	2-Methylnaphthalene-d10	0.400	0.415	-3.7	45	-0.03
12		2-Methylnaphthalene	0.400	0.421	-5.2	46	-0.04
13	I	Acenaphthene-d10	0.400	0.400	0.0	42	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.379	5.3	47	-0.03
15	S	2-Fluorobiphenyl	0.400	0.371	7.3	42	-0.03
16		Acenaphthylene	0.400	0.378	5.5	44	-0.04
17		Acenaphthene	0.400	0.366	8.5	42	-0.04
18		Fluorene	0.400	0.370	7.5	42	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	39	-0.03
20		4-Bromophenyl-phenylether	0.400	0.382	4.5	38	-0.04
21		Hexachlorobenzene	0.400	0.426	-6.5	42	-0.03
22		Pentachlorophenol	0.400	0.454	-13.5	51	-0.03
23		Phenanthrene	0.400	0.400	0.0	40	-0.03
24		Anthracene	0.400	0.412	-3.0	42	-0.03
25	SURR	Fluoranthene-d10	0.400	0.426	-6.5	42	-0.03
26		Fluoranthene	0.400	0.437	-9.2	43	-0.03
27	I	Chrysene-d12	0.400	0.400	0.0	52	-0.03
28		Pyrene	0.400	0.338	15.5	45	-0.03
29	S	Terphenyl-d14	0.400	0.349	12.8	47	-0.03
30		Benzo(a)anthracene	0.400	0.341	14.8	46	-0.02
31		Chrysene	0.400	0.370	7.5	49	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.650	-62.5#	94	-0.04
33		Indeno(1,2,3-cd)pyrene	0.400	0.356	11.0	47	-0.05
34	I	Perylene-d12	0.400	0.400	0.0	45	-0.04
35		Benzo(b)fluoranthene	0.400	0.417	-4.2	50	-0.03
36		Benzo(k)fluoranthene	0.400	0.399	0.3	47	-0.04
37	C	Benzo(a)pyrene	0.400	0.389	2.8	46	-0.04
38		Dibenzo(a,h)anthracene	0.400	0.395	1.3	47	-0.05
39		Benzo(g,h,i)perylene	0.400	0.405	-1.3	48	-0.05

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

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