Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN032420\
Data File : BN010347.D
Acq On : 25 Mar 2020 03:40
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Mar 25 04:12:06 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 19 15:26:45 2020
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	107	0.00
2		1,4-Dioxane	0.400	0.383	4.3	100	0.00
3		n-Nitrosodimethylamine	0.400	0.405	-1.3	114	0.02
4	S	2-Fluorophenol	0.400	0.404	-1.0	109	0.00
5	S	Phenol-d6	0.400	0.402	-0.5	110	0.00
6		bis(2-Chloroethyl)ether	0.400	0.404	-1.0	106	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	108	0.00
8	S	Nitrobenzene-d5	0.400	0.387	3.3	109	0.00
9		Naphthalene	0.400	0.405	-1.3	109	-0.01
10		Hexachlorobutadiene	0.400	0.411	-2.7	112	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.449	-12.2	124	0.00
12		2-Methylnaphthalene	0.400	0.402	-0.5	110	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	107	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.346	13.5	104	0.00
15	S	2-Fluorobiphenyl	0.400	0.409	-2.2	108	-0.01
16		Acenaphthylene	0.400	0.370	7.5	107	-0.01
17		Acenaphthene	0.400	0.408	-2.0	110	0.00
18		Fluorene	0.400	0.394	1.5	109	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	104	0.00
20		4-Bromophenyl-phenylether	0.400	0.381	4.8	101	0.00
21		Hexachlorobenzene	0.400	0.429	-7.2	113	0.00
22		Pentachlorophenol	0.400	0.317	20.8	97	0.00
23		Phenanthrene	0.400	0.406	-1.5	108	-0.01
24		Anthracene	0.400	0.374	6.5	104	0.00
25	SURR	Fluoranthene-d10	0.400	0.434	-8.5	119	0.00
26		Fluoranthene	0.400	0.392	2.0	107	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	101	0.00
28		Pyrene	0.400	0.421	-5.2	107	0.00
29	S	Terphenyl-d14	0.400	0.410	-2.5	103	0.00
30		Benzo(a)anthracene	0.400	0.376	6.0	99	-0.01
31		Chrysene	0.400	0.408	-2.0	103	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.311	22.3	88	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.417	-4.2	98	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	95	-0.01
35		Benzo(b)fluoranthene	0.400	0.398	0.5	104	-0.01
36		Benzo(k)fluoranthene	0.400	0.415	-3.7	101	-0.01
37	C	Benzo(a)pyrene	0.400	0.395	1.3	103	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.409	-2.2	96	-0.02
39		Benzo(g,h,i)perylene	0.400	0.413	-3.2	103	-0.02

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

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