Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112519\
Data File : BN008750.D
Acq On : 25 Nov 2019 23:03
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Nov 26 02:55:11 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN112219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 25 10:49:40 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	109	0.00
2		1,4-Dioxane	0.400	0.450	-12.5	123	0.00
3		n-Nitrosodimethylamine	0.400	0.399	0.3	117	0.00
4	S	2-Fluorophenol	0.400	0.463	-15.8	126	0.00
5	S	Phenol-d6	0.400	0.460	-15.0	124	0.00
6		bis(2-Chloroethyl)ether	0.400	0.452	-13.0	121	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	110	0.00
8	S	Nitrobenzene-d5	0.400	0.452	-13.0	125	-0.01
9		Naphthalene	0.400	0.455	-13.7	123	-0.01
10		Hexachlorobutadiene	0.400	0.449	-12.2	121	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.448	-12.0	123	0.00
12		2-Methylnaphthalene	0.400	0.449	-12.2	123	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	109	0.00
14	S	2,4,6-Tribromophenol	0.400	0.440	-10.0	125	0.00
15	S	2-Fluorobiphenyl	0.400	0.461	-15.3	124	-0.01
16		Acenaphthylene	0.400	0.457	-14.2	127	-0.01
17		Acenaphthene	0.400	0.439	-9.7	122	-0.01
18		Fluorene	0.400	0.457	-14.2	125	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	109	-0.01
20		4-Bromophenyl-phenylether	0.400	0.457	-14.2	127	-0.01
21		Hexachlorobenzene	0.400	0.457	-14.2	125	-0.01
22		Pentachlorophenol	0.400	0.368	8.0	118	-0.01
23		Phenanthrene	0.400	0.454	-13.5	125	-0.01
24		Anthracene	0.400	0.457	-14.2	128	0.00
25	SURR	Fluoranthene-d10	0.400	0.454	-13.5	125	0.00
26		Fluoranthene	0.400	0.452	-13.0	125	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	109	-0.01
28		Pyrene	0.400	0.461	-15.3	124	0.00
29	S	Terphenyl-d14	0.400	0.469	-17.2	126	0.00
30		Benzo(a)anthracene	0.400	0.440	-10.0	120	-0.01
31		Chrysene	0.400	0.463	-15.8	124	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.346	13.5	102	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.451	-12.7	121	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	106	-0.02
35		Benzo(b)fluoranthene	0.400	0.469	-17.2	124	-0.01
36		Benzo(k)fluoranthene	0.400	0.462	-15.5	121	-0.01
37	C	Benzo(a)pyrene	0.400	0.458	-14.5	122	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.460	-15.0	121	-0.02
39		Benzo(g,h,i)perylene	0.400	0.465	-16.3	123	-0.02

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

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