Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092519\
Data File : BN007867.D
Acq On : 25 Sep 2019 23:04
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 26 05:25:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 18:24:40 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	176	0.00
2		1,4-Dioxane	0.400	0.395	1.3	205	0.00
3		n-Nitrosodimethylamine	0.400	0.542	-35.5#	157	0.00
4	S	2-Fluorophenol	0.400	0.368	8.0	186	0.00
5	S	Phenol-d6	0.400	0.385	3.8	186	0.00
6		bis(2-Chloroethyl)ether	0.400	0.472	-18.0	235	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	175	-0.01
8	S	Nitrobenzene-d5	0.400	0.426	-6.5	192	0.00
9		Naphthalene	0.400	0.427	-6.7	188	-0.01
10		Hexachlorobutadiene	0.400	0.446	-11.5	194	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.415	-3.7	187	-0.01
12		2-Methylnaphthalene	0.400	0.418	-4.5	188	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	154	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.305	23.8	140	-0.01
15	S	2-Fluorobiphenyl	0.400	0.419	-4.7	174	-0.01
16		Acenaphthylene	0.400	0.738	-84.5#	316	-0.01
17		Acenaphthene	0.400	0.420	-5.0	177	-0.01
18		Fluorene	0.400	0.434	-8.5	182	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	125	-0.01
20		4-Bromophenyl-phenylether	0.400	0.462	-15.5	146	-0.01
21		Hexachlorobenzene	0.400	0.451	-12.7	142	-0.01
22		Pentachlorophenol	0.400	0.404	-1.0	145	-0.01
23		Phenanthrene	0.400	0.427	-6.7	135	-0.01
24		Anthracene	0.400	0.450	-12.5	145	-0.01
25	SURR	Fluoranthene-d10	0.400	0.382	4.5	121	-0.02
26		Fluoranthene	0.400	0.385	3.8	121	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	102	-0.02
28		Pyrene	0.400	0.482	-20.5	128	-0.02
29	S	Terphenyl-d14	0.400	0.452	-13.0	120	-0.02
30		Benzo(a)anthracene	0.400	0.408	-2.0	109	-0.01
31		Chrysene	0.400	0.405	-1.3	106	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.519	-29.8#	148	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.397	0.8	104	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	99	-0.03
35		Benzo(b)fluoranthene	0.400	0.411	-2.7	108	-0.02
36		Benzo(k)fluoranthene	0.400	0.397	0.8	103	-0.02
37	C	Benzo(a)pyrene	0.400	0.414	-3.5	108	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.394	1.5	102	-0.03
39		Benzo(g,h,i)perylene	0.400	0.435	-8.7	113	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092519\
Data File : BN007867.D
Acq On : 25 Sep 2019 23:04
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 26 05:25:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 18:24:40 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)

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