Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100864.D
Acq On : 12 Dec 2019 9:10
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 12 10:11:04 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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	99	0.00
2		1,4-Dioxane	0.400	0.475	-18.7	118	0.00
3		n-Nitrosodimethylamine	0.400	0.477	-19.2	135	-0.02
4	S	2-Fluorophenol	0.400	0.432	-8.0	114	0.00
5	S	Phenol-d6	0.400	0.444	-11.0	119	0.00
6		bis(2-Chloroethyl)ether	0.400	0.467	-16.8	122	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	102	0.00
8	S	Nitrobenzene-d5	0.400	0.580	-45.0#	170	0.00
9		Naphthalene	0.400	0.459	-14.8	122	0.00
10		Hexachlorobutadiene	0.400	0.455	-13.7	120	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.460	-15.0	125	0.00
12		2-Methylnaphthalene	0.400	0.463	-15.8	126	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	111	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.704	-76.0#	224	-0.01
15	S	2-Fluorobiphenyl	0.400	0.418	-4.5	121	-0.01
16		Acenaphthylene	0.400	0.440	-10.0	142	0.00
17		Acenaphthene	0.400	0.451	-12.7	137	0.00
18		Fluorene	0.400	0.472	-18.0	147	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	126	-0.01
20		4-Bromophenyl-phenylether	0.400	0.405	-1.3	141	-0.01
21		Hexachlorobenzene	0.400	0.414	-3.5	139	0.00
22		Pentachlorophenol	0.400	0.767	-91.8#	1253	-0.01
23		Phenanthrene	0.400	0.429	-7.2	144	0.00
24		Anthracene	0.400	0.435	-8.7	160	0.00
25	SURR	Fluoranthene-d10	0.400	0.376	6.0	132	-0.01
26		Fluoranthene	0.400	0.384	4.0	134	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	97	-0.01
28		Pyrene	0.400	0.510	-27.5#	132	-0.01
29	S	Terphenyl-d14	0.400	0.462	-15.5	117	-0.01
30		Benzo(a)anthracene	0.400	0.465	-16.3	127	-0.01
31		Chrysene	0.400	0.472	-18.0	118	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.703	-75.7#	178	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.465	-16.3	127	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	107	-0.01
35		Benzo(b)fluoranthene	0.400	0.418	-4.5	118	-0.01
36		Benzo(k)fluoranthene	0.400	0.447	-11.7	126	-0.02
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	121	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.426	-6.5	128	-0.04
39		Benzo(g,h,i)perylene	0.400	0.427	-6.7	121	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100864.D
Acq On : 12 Dec 2019 9:10
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 12 10:11:04 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 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