Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061319\
Data File : BE100010.D
Acq On : 14 Jun 2019 3:24
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 14 04:47:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 10 14:57:20 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	106	-0.01
2		1,4-Dioxane	0.400	0.392	2.0	111	0.00
3		n-Nitrosodimethylamine	0.400	0.360	10.0	95	0.01
4	S	2-Fluorophenol	0.400	0.424	-6.0	112	0.00
5	S	Phenol-d6	0.400	0.451	-12.7	128	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.412	-3.0	115	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	111	-0.01
8	S	Nitrobenzene-d5	0.400	0.397	0.8	117	0.00
9		Naphthalene	0.400	0.418	-4.5	117	-0.01
10		Hexachlorobutadiene	0.400	0.445	-11.2	121	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.427	-6.7	121	0.00
12		2-Methylnaphthalene	0.400	0.426	-6.5	120	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	120	0.00
14	S	2,4,6-Tribromophenol	0.400	0.338	15.5	113	0.00
15	S	2-Fluorobiphenyl	0.400	0.401	-0.3	127	-0.01
16		Acenaphthylene	0.400	0.404	-1.0	126	-0.01
17		Acenaphthene	0.400	0.410	-2.5	128	0.00
18		Fluorene	0.400	0.418	-4.5	132	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	126	0.00
20		4-Bromophenyl-phenylether	0.400	0.385	3.8	132	-0.01
21		Hexachlorobenzene	0.400	0.408	-2.0	136	-0.01
22		Pentachlorophenol	0.400	0.528	-32.0#	126	-0.03
23		Phenanthrene	0.400	0.392	2.0	132	0.00
24		Anthracene	0.400	0.386	3.5	129	0.00
25	SURR	Fluoranthene-d10	0.400	0.369	7.8	119	0.00
26		Fluoranthene	0.400	0.379	5.3	122	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	110	-0.01
28		Pyrene	0.400	0.479	-19.7	124	0.00
29	S	Terphenyl-d14	0.400	0.446	-11.5	117	0.00
30		Benzo(a)anthracene	0.400	0.429	-7.2	114	0.00
31		Chrysene	0.400	0.473	-18.2	120	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.405	-1.3	109	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.435	-8.7	113	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	107	0.00
35		Benzo(b)fluoranthene	0.400	0.425	-6.2	119	0.00
36		Benzo(k)fluoranthene	0.400	0.429	-7.2	115	0.00
37	C	Benzo(a)pyrene	0.400	0.426	-6.5	117	0.00
38		Dibenzo(a,h)anthracene	0.400	0.406	-1.5	114	0.00
39		Benzo(g,h,i)perylene	0.400	0.416	-4.0	114	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061319\
Data File : BE100010.D
Acq On : 14 Jun 2019 3:24
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jun 14 04:47:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 10 14:57:20 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