Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122719\
Data File : BE100999.D
Acq On : 27 Dec 2019 18:44
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 27 23:40:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 17:39:05 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	0.214	46.5#	82	0.00
3		n-Nitrosodimethylamine	0.400	0.580	-45.0#	184	-0.01
4	S	2-Fluorophenol	0.400	0.405	-1.3	123	0.00
5	S	Phenol-d6	0.400	0.449	-12.2	123	-0.03
6		bis(2-Chloroethyl)ether	0.400	0.287	28.3#	93	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	111	0.00
8	S	Nitrobenzene-d5	0.400	0.445	-11.2	116	0.00
9		Naphthalene	0.400	0.413	-3.2	114	0.00
10		Hexachlorobutadiene	0.400	0.422	-5.5	115	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.436	-9.0	118	0.00
12		2-Methylnaphthalene	0.400	0.374	6.5	102	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	107	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.427	-6.7	107	-0.01
15	S	2-Fluorobiphenyl	0.400	0.432	-8.0	114	0.00
16		Acenaphthylene	0.400	0.349	12.8	101	0.00
17		Acenaphthene	0.400	0.413	-3.2	110	0.00
18		Fluorene	0.400	0.397	0.8	110	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	102	-0.01
20		4-Bromophenyl-phenylether	0.400	0.376	6.0	109	0.00
21		Hexachlorobenzene	0.400	0.397	0.8	105	0.00
22		Pentachlorophenol	0.400	0.316	21.0	68	-0.04
23		Phenanthrene	0.400	0.394	1.5	113	0.00
24		Anthracene	0.400	0.357	10.8	101	0.00
25	SURR	Fluoranthene-d10	0.400	0.370	7.5	102	0.00
26		Fluoranthene	0.400	0.367	8.3	104	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	92	0.00
28		Pyrene	0.400	0.449	-12.2	102	0.00
29	S	Terphenyl-d14	0.400	0.463	-15.8	108	-0.01
30		Benzo(a)anthracene	0.400	0.410	-2.5	109	0.00
31		Chrysene	0.400	0.421	-5.2	87	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.377	5.8	89	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.416	-4.0	100	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	88	0.00
35		Benzo(b)fluoranthene	0.400	0.449	-12.2	101	0.00
36		Benzo(k)fluoranthene	0.400	0.415	-3.7	90	0.00
37	C	Benzo(a)pyrene	0.400	0.392	2.0	88	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.356	11.0	94	-0.01
39		Benzo(g,h,i)perylene	0.400	0.442	-10.5	97	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE122719\
Data File : BE100999.D
Acq On : 27 Dec 2019 18:44
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 27 23:40:37 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 27 17:39:05 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