Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101917\
Data File : BE094455.D
Acq On : 19 Oct 2017 19:40
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 20 12:42:36 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 2017
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	88	-0.01
2		1,4-Dioxane	0.400	0.485	-21.2	101	0.00
3		n-Nitrosodimethylamine	0.400	0.411	-2.7	91	0.01
4	S	2-Fluorophenol	0.400	0.450	-12.5	98	0.00
5	S	Phenol-d6	0.400	0.440	-10.0	94	0.00
6		bis(2-Chloroethyl)ether	0.400	0.428	-7.0	91	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	90	-0.02
8	S	Nitrobenzene-d5	0.400	0.418	-4.5	95	0.00
9		Naphthalene	0.400	0.432	-8.0	97	-0.02
10		Hexachlorobutadiene	0.400	0.407	-1.7	92	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.445	-11.2	100	0.00
12		2-Methylnaphthalene	0.400	0.452	-13.0	103	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	101	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.425	-6.2	101	0.00
15	S	2-Fluorobiphenyl	0.400	0.412	-3.0	101	0.00
16		Acenaphthylene	0.400	0.419	-4.7	105	0.00
17		Acenaphthene	0.400	0.413	-3.2	104	0.00
18		Fluorene	0.400	0.419	-4.7	106	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	105	-0.03
20		4-Bromophenyl-phenylether	0.400	0.484	-21.0	141	-0.03
21		Hexachlorobenzene	0.400	0.259	35.3#	72	0.00
22		Pentachlorophenol	0.400	0.523	-30.8#	137	0.00
23		Phenanthrene	0.400	0.286	28.5#	72	0.00
24		Anthracene	0.400	0.397	0.8	107	-0.03
25	SURR	Fluoranthene-d10	0.400	0.368	8.0	100	0.00
26		Fluoranthene	0.400	0.366	8.5	101	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	100	0.00
28		Pyrene	0.400	0.397	0.8	100	0.00
29	S	Terphenyl-d14	0.400	0.403	-0.8	100	0.00
30		Benzo(a)anthracene	0.400	0.407	-1.7	103	-0.01
31		Chrysene	0.400	0.400	0.0	98	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.368	8.0	96	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	102	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	100	-0.02
35		Benzo(b)fluoranthene	0.400	0.412	-3.0	103	-0.01
36		Benzo(k)fluoranthene	0.400	0.417	-4.2	103	-0.01
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	103	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.412	-3.0	102	-0.03
39		Benzo(g,h,i)perylene	0.400	0.407	-1.7	101	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101917\
Data File : BE094455.D
Acq On : 19 Oct 2017 19:40
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 20 12:42:36 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 2017
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