Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE032219\
Data File : BE099101.D
Acq On : 23 Mar 2019 4:18
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 23 07:08:46 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 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	115	0.00
2		1,4-Dioxane	0.400	0.329	17.8	109	0.00
3		n-Nitrosodimethylamine	0.400	0.296	26.0#	92	-0.01
4	S	2-Fluorophenol	0.400	0.337	15.8	109	0.00
5	S	Phenol-d6	0.400	0.345	13.8	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.352	12.0	113	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	124	0.00
8	S	Nitrobenzene-d5	0.400	0.529	-32.3#	206	-0.01
9		Naphthalene	0.400	0.318	20.5	111	0.00
10		Hexachlorobutadiene	0.400	0.322	19.5	111	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.330	17.5	117	-0.02
12		2-Methylnaphthalene	0.400	0.337	15.8	121	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	147	0.00
14	S	2,4,6-Tribromophenol	0.400	0.493	-23.2	234	0.00
15	S	2-Fluorobiphenyl	0.400	0.299	25.3#	120	0.00
16		Acenaphthylene	0.400	0.302	24.5	131	-0.02
17		Acenaphthene	0.400	0.301	24.8	130	0.00
18		Fluorene	0.400	0.307	23.3	131	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	147	0.00
20		4-Bromophenyl-phenylether	0.400	0.322	19.5	139	0.00
21		Hexachlorobenzene	0.400	0.324	19.0	142	0.00
22		Pentachlorophenol	0.400	0.556	-39.0#	226	0.00
23		Phenanthrene	0.400	0.308	23.0	134	0.00
24		Anthracene	0.400	0.299	25.3#	132	0.00
25	SURR	Fluoranthene-d10	0.400	0.289	27.8#	123	0.00
26		Fluoranthene	0.400	0.277	30.8#	122	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	128	0.00
28		Pyrene	0.400	0.321	19.8	121	0.00
29	S	Terphenyl-d14	0.400	0.334	16.5	121	0.00
30		Benzo(a)anthracene	0.400	0.306	23.5	115	0.00
31		Chrysene	0.400	0.317	20.8	121	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.354	11.5	160	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.317	20.8	120	0.00
34	I	Perylene-d12	0.400	0.400	0.0	134	0.00
35		Benzo(b)fluoranthene	0.400	0.296	26.0#	120	0.00
36		Benzo(k)fluoranthene	0.400	0.300	25.0#	120	0.00
37	C	Benzo(a)pyrene	0.400	0.298	25.5#	121	0.00
38		Dibenzo(a,h)anthracene	0.400	0.301	24.8	123	-0.01
39		Benzo(g,h,i)perylene	0.400	0.297	25.8#	119	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE032219\
Data File : BE099101.D
Acq On : 23 Mar 2019 4:18
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Mar 23 07:08:46 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 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 = 1