Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101119\
Data File : BE100639.D
Acq On : 11 Oct 2019 17:21
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 11 18:35:35 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 10 14:01:52 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	79	-0.01
2		1,4-Dioxane	0.400	0.399	0.3	80	0.00
3		n-Nitrosodimethylamine	0.400	0.408	-2.0	88	-0.01
4	S	2-Fluorophenol	0.400	0.491	-22.7	102	0.00
5	S	Phenol-d6	0.400	0.507	-26.7#	106	0.00
6		bis(2-Chloroethyl)ether	0.400	0.388	3.0	82	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	80	0.00
8	S	Nitrobenzene-d5	0.400	0.547	-36.8#	124	0.00
9		Naphthalene	0.400	0.395	1.3	81	0.00
10		Hexachlorobutadiene	0.400	0.414	-3.5	86	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.380	5.0	79	0.00
12		2-Methylnaphthalene	0.400	0.390	2.5	81	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	79	0.00
14	S	2,4,6-Tribromophenol	0.400	0.540	-35.0#	121	0.00
15	S	2-Fluorobiphenyl	0.400	0.517	-29.3#	104	0.00
16		Acenaphthylene	0.400	0.360	10.0	74	0.00
17		Acenaphthene	0.400	0.404	-1.0	83	0.00
18		Fluorene	0.400	0.381	4.8	78	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	81	0.00
20		4-Bromophenyl-phenylether	0.400	0.375	6.3	79	-0.01
21		Hexachlorobenzene	0.400	0.398	0.5	85	0.00
22		Pentachlorophenol	0.400	0.496	-24.0	113	0.00
23		Phenanthrene	0.400	0.407	-1.7	82	0.00
24		Anthracene	0.400	0.366	8.5	77	0.00
25	SURR	Fluoranthene-d10	0.400	0.391	2.3	82	0.00
26		Fluoranthene	0.400	0.419	-4.7	85	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	91	0.00
28		Pyrene	0.400	0.370	7.5	85	0.00
29	S	Terphenyl-d14	0.400	0.478	-19.5	115	0.00
30		Benzo(a)anthracene	0.400	0.377	5.8	87	0.00
31		Chrysene	0.400	0.397	0.8	91	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.278	30.5#	63	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.346	13.5	82	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	85	0.00
35		Benzo(b)fluoranthene	0.400	0.366	8.5	84	0.00
36		Benzo(k)fluoranthene	0.400	0.375	6.3	85	0.00
37	C	Benzo(a)pyrene	0.400	0.344	14.0	79	0.00
38		Dibenzo(a,h)anthracene	0.400	0.345	13.8	81	-0.01
39		Benzo(g,h,i)perylene	0.400	0.366	8.5	83	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101119\
Data File : BE100639.D
Acq On : 11 Oct 2019 17:21
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 11 18:35:35 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 10 14:01:52 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