Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071219\
Data File : BE100423.D
Acq On : 12 Jul 2019 17:17
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE071219

Quant Time: Jul 12 18:04:23 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 17:51:04 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	108	-0.01
2		1,4-Dioxane	0.400	0.432	-8.0	122	-0.01
3		n-Nitrosodimethylamine	0.400	0.415	-3.7	102	-0.01
4	S	2-Fluorophenol	0.400	0.416	-4.0	110	0.00
5	S	Phenol-d6	0.400	0.402	-0.5	106	0.00
6		bis(2-Chloroethyl)ether	0.400	0.411	-2.7	110	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	110	0.00
8	S	Nitrobenzene-d5	0.400	0.429	-7.2	108	0.00
9		Naphthalene	0.400	0.434	-8.5	110	0.00
10		Hexachlorobutadiene	0.400	0.415	-3.7	108	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.433	-8.2	111	0.00
12		2-Methylnaphthalene	0.400	0.435	-8.7	111	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	111	0.00
14	S	2,4,6-Tribromophenol	0.400	0.392	2.0	115	0.00
15	S	2-Fluorobiphenyl	0.400	0.434	-8.5	112	0.00
16		Acenaphthylene	0.400	0.427	-6.7	109	0.00
17		Acenaphthene	0.400	0.434	-8.5	111	0.01
18		Fluorene	0.400	0.433	-8.2	108	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	111	0.00
20		4-Bromophenyl-phenylether	0.400	0.429	-7.2	112	0.00
21		Hexachlorobenzene	0.400	0.428	-7.0	112	0.00
22		Pentachlorophenol	0.400	0.414	-3.5	117	0.00
23		Phenanthrene	0.400	0.449	-12.2	113	0.00
24		Anthracene	0.400	0.435	-8.7	110	0.01
25	SURR	Fluoranthene-d10	0.400	0.429	-7.2	106	0.00
26		Fluoranthene	0.400	0.436	-9.0	107	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	104	0.00
28		Pyrene	0.400	0.457	-14.2	105	0.00
29	S	Terphenyl-d14	0.400	0.450	-12.5	105	0.00
30		Benzo(a)anthracene	0.400	0.434	-8.5	101	0.00
31		Chrysene	0.400	0.438	-9.5	102	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.377	5.8	106	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	98	0.00
34	I	Perylene-d12	0.400	0.400	0.0	101	0.00
35		Benzo(b)fluoranthene	0.400	0.425	-6.2	96	0.00
36		Benzo(k)fluoranthene	0.400	0.459	-14.8	108	0.00
37	C	Benzo(a)pyrene	0.400	0.426	-6.5	98	0.00
38		Dibenzo(a,h)anthracene	0.400	0.420	-5.0	98	0.00
39		Benzo(g,h,i)perylene	0.400	0.432	-8.0	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071219\
Data File : BE100423.D
Acq On : 12 Jul 2019 17:17
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
ICVBE071219

Quant Time: Jul 12 18:04:23 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 12 17:51:04 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