Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
Data File : BE100067.D
Acq On : 17 Jun 2019 17:25
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE061719

Quant Time: Jun 17 18:06:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 17 17:23:36 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	88	0.01
2		1,4-Dioxane	0.400	0.375	6.3	83	0.01
3		n-Nitrosodimethylamine	0.400	0.442	-10.5	77	0.02
4	S	2-Fluorophenol	0.400	0.369	7.8	87	0.01
5	S	Phenol-d6	0.400	0.396	1.0	92	0.02
6		bis(2-Chloroethyl)ether	0.400	0.427	-6.7	92	0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	92	0.01
8	S	Nitrobenzene-d5	0.400	0.350	12.5	89	0.03
9		Naphthalene	0.400	0.411	-2.7	94	0.01
10		Hexachlorobutadiene	0.400	0.426	-6.5	97	0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.414	-3.5	98	0.02
12		2-Methylnaphthalene	0.400	0.407	-1.7	97	0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	96	0.02
14	S	2,4,6-Tribromophenol	0.400	0.359	10.3	99	0.01
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	95	0.02
16		Acenaphthylene	0.400	0.398	0.5	97	0.02
17		Acenaphthene	0.400	0.400	0.0	95	0.00
18		Fluorene	0.400	0.385	3.8	93	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	93	0.01
20		4-Bromophenyl-phenylether	0.400	0.391	2.3	93	0.01
21		Hexachlorobenzene	0.400	0.404	-1.0	96	0.01
22		Pentachlorophenol	0.400	0.355	11.3	69	0.03
23		Phenanthrene	0.400	0.387	3.3	93	0.00
24		Anthracene	0.400	0.369	7.8	91	0.00
25	SURR	Fluoranthene-d10	0.400	0.385	3.8	93	0.00
26		Fluoranthene	0.400	0.390	2.5	93	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	91	0.01
28		Pyrene	0.400	0.449	-12.2	92	0.01
29	S	Terphenyl-d14	0.400	0.429	-7.2	90	0.02
30		Benzo(a)anthracene	0.400	0.446	-11.5	95	0.00
31		Chrysene	0.400	0.436	-9.0	88	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.363	9.3	82	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.439	-9.7	90	0.01
34	I	Perylene-d12	0.400	0.400	0.0	90	0.00
35		Benzo(b)fluoranthene	0.400	0.400	0.0	89	0.01
36		Benzo(k)fluoranthene	0.400	0.412	-3.0	91	0.00
37	C	Benzo(a)pyrene	0.400	0.410	-2.5	92	0.00
38		Dibenzo(a,h)anthracene	0.400	0.391	2.3	88	0.02
39		Benzo(g,h,i)perylene	0.400	0.408	-2.0	91	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
Data File : BE100067.D
Acq On : 17 Jun 2019 17:25
Operator : JU/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE061719

Quant Time: Jun 17 18:06:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 17 17:23:36 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