Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062320\
Data File : BN011037.D
Acq On : 23 Jun 2020 19:34
Operator : CG/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN062320

Quant Time: Jun 23 20:07:50 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 23 18:53:50 2020
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	100	0.00
2		1,4-Dioxane	0.400	0.431	-7.7	100	0.00
3		n-Nitrosodimethylamine	0.400	0.394	1.5	92	0.00
4	S	2-Fluorophenol	0.400	0.406	-1.5	95	0.00
5	S	Phenol-d6	0.400	0.398	0.5	92	0.00
6		bis(2-Chloroethyl)ether	0.400	0.413	-3.2	95	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	94	0.00
8	S	Nitrobenzene-d5	0.400	0.394	1.5	81	0.01
9		Naphthalene	0.400	0.409	-2.2	98	0.00
10		Hexachlorobutadiene	0.400	0.432	-8.0	98	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.417	-4.2	92	0.00
12		2-Methylnaphthalene	0.400	0.420	-5.0	93	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14	S	2,4,6-Tribromophenol	0.400	0.354	11.5	93	0.01
15	S	2-Fluorobiphenyl	0.400	0.430	-7.5	109	0.00
16		Acenaphthylene	0.400	0.395	1.3	105	0.00
17		Acenaphthene	0.400	0.405	-1.3	105	0.00
18		Fluorene	0.400	0.369	7.8	96	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	95	0.00
20		4,6-Dinitro-2-methylphenol	0.400	0.356	11.0	91	0.01
21		4-Bromophenyl-phenylether	0.400	0.405	-1.3	94	0.01
22		Hexachlorobenzene	0.400	0.422	-5.5	96	0.01
23		Atrazine	0.400	0.394	1.5	96	0.00
24		Pentachlorophenol	0.400	0.371	7.3	88	0.00
25		Phenanthrene	0.400	0.414	-3.5	96	0.00
26		Anthracene	0.400	0.395	1.3	94	0.00
27	SURR	Fluoranthene-d10	0.400	0.396	1.0	94	0.00
28		Fluoranthene	0.400	0.396	1.0	91	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	105	-0.01
30		Pyrene	0.400	0.369	7.8	96	0.00
31	S	Terphenyl-d14	0.400	0.374	6.5	96	0.00
32		Benzo(a)anthracene	0.400	0.393	1.8	101	0.00
33		Chrysene	0.400	0.413	-3.2	108	-0.01
34		Bis(2-ethylhexyl)phthalate	0.400	0.363	9.3	94	0.00
35		Indeno(1,2,3-cd)pyrene	0.400	0.454	-13.5	113	-0.01
36	I	Perylene-d12	0.400	0.400	0.0	88	-0.01
37		Benzo(b)fluoranthene	0.400	0.455	-13.7	113	-0.01
38		Benzo(k)fluoranthene	0.400	0.488	-22.0	125	-0.01
39	C	Benzo(a)pyrene	0.400	0.405	-1.3	88	-0.01
40		Dibenzo(a,h)anthracene	0.400	0.447	-11.7	111	-0.01
41		Benzo(g,h,i)perylene	0.400	0.403	-0.8	103	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062320\
Data File : BN011037.D
Acq On : 23 Jun 2020 19:34
Operator : CG/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
ICVBN062320

Quant Time: Jun 23 20:07:50 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 23 18:53:50 2020
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