Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082120\
Data File : BN011564.D
Acq On : 21 Aug 2020 20:37
Operator : CG/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN082120

Quant Time: Aug 22 05:54:51 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 22 04:52:07 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	110	0.00
2		1,4-Dioxane	0.400	0.398	0.5	109	0.00
3		n-Nitrosodimethylamine	0.400	0.348	13.0	102	0.00
4	S	2-Fluorophenol	0.400	0.372	7.0	117	0.00
5	S	Phenol-d6	0.400	0.331	17.3	119	0.00
6		bis(2-Chloroethyl)ether	0.400	0.343	14.2	123	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	124	0.00
8	S	Nitrobenzene-d5	0.400	0.342	14.5	128	0.00
9		Naphthalene	0.400	0.372	7.0	122	-0.01
10		Hexachlorobutadiene	0.400	0.375	6.3	116	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.370	7.5	129	0.00
12		2-Methylnaphthalene	0.400	0.373	6.8	128	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	134	0.00
14	S	2,4,6-Tribromophenol	0.400	0.326	18.5	128	0.00
15	S	2-Fluorobiphenyl	0.400	0.367	8.3	130	0.00
16		Acenaphthylene	0.400	0.354	11.5	132	0.00
17		Acenaphthene	0.400	0.371	7.3	134	0.00
18		Fluorene	0.400	0.378	5.5	138	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	136	0.00
20		4,6-Dinitro-2-methylphenol	0.400	0.330	17.5	114	-0.01
21		4-Bromophenyl-phenylether	0.400	0.367	8.3	107	0.00
22		Hexachlorobenzene	0.400	0.365	8.8	130	0.00
23		Atrazine	0.400	0.253	36.8#	139	0.00
24		Pentachlorophenol	0.400	0.277	30.8#	119	-0.01
25		Phenanthrene	0.400	0.356	11.0	135	0.00
26		Anthracene	0.400	0.341	14.8	138	0.00
27	SURR	Fluoranthene-d10	0.400	0.336	16.0	139	0.00
28		Fluoranthene	0.400	0.336	16.0	139	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	155	0.00
30		Pyrene	0.400	0.347	13.3	135	0.00
31	S	Terphenyl-d14	0.400	0.345	13.8	141	0.00
32		Benzo(a)anthracene	0.400	0.350	12.5	161	-0.01
33		Chrysene	0.400	0.377	5.8	152	0.00
34		Bis(2-ethylhexyl)phthalate	0.400	0.382	4.5	163	0.00
35		Indeno(1,2,3-cd)pyrene	0.400	0.357	10.8	151	0.00
36	I	Perylene-d12	0.400	0.400	0.0	149	0.00
37		Benzo(b)fluoranthene	0.400	0.361	9.8	178	0.00
38		Benzo(k)fluoranthene	0.400	0.380	5.0	173	0.00
39	C	Benzo(a)pyrene	0.400	0.353	11.8	149	0.00
40		Dibenzo(a,h)anthracene	0.400	0.348	13.0	152	0.00
41		Benzo(g,h,i)perylene	0.400	0.335	16.3	138	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082120\
Data File : BN011564.D
Acq On : 21 Aug 2020 20:37
Operator : CG/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBN082120

Quant Time: Aug 22 05:54:51 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 22 04:52:07 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