Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071020\
Data File : BN011132.D
Acq On : 09 Jul 2020 14:56
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 09 15:34:53 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 09 15:33:27 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	113	0.00
2		1,4-Dioxane	0.400	0.410	-2.5	112	0.00
3		n-Nitrosodimethylamine	0.400	0.410	-2.5	121	0.00
4	S	2-Fluorophenol	0.400	0.466	-16.5	128	0.00
5	S	Phenol-d6	0.400	0.418	-4.5	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.411	-2.7	105	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	113	0.00
8	S	Nitrobenzene-d5	0.400	0.419	-4.7	118	0.00
9		Naphthalene	0.400	0.405	-1.3	115	0.00
10		Hexachlorobutadiene	0.400	0.401	-0.3	104	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.400	0.0	108	0.00
12		2-Methylnaphthalene	0.400	0.406	-1.5	113	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	114	0.00
14	S	2,4,6-Tribromophenol	0.400	0.397	0.8	112	0.00
15	S	2-Fluorobiphenyl	0.400	0.416	-4.0	116	0.00
16		Acenaphthylene	0.400	0.400	0.0	114	0.00
17		Acenaphthene	0.400	0.406	-1.5	115	0.00
18		Fluorene	0.400	0.410	-2.5	115	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	115	0.00
20		4,6-Dinitro-2-methylphenol	0.400	0.477	-19.2	146	0.00
21		4-Bromophenyl-phenylether	0.400	0.386	3.5	110	0.00
22		Hexachlorobenzene	0.400	0.410	-2.5	116	0.00
23		Atrazine	0.400	0.390	2.5	109	0.00
24		Pentachlorophenol	0.400	0.435	-8.7	127	0.00
25		Phenanthrene	0.400	0.403	-0.8	114	0.00
26		Anthracene	0.400	0.393	1.8	111	0.00
27	SURR	Fluoranthene-d10	0.400	0.401	-0.3	114	0.00
28		Fluoranthene	0.400	0.405	-1.3	116	0.00
29	I	Chrysene-d12	0.400	0.400	0.0	103	0.00
30		Pyrene	0.400	0.428	-7.0	115	0.00
31	S	Terphenyl-d14	0.400	0.426	-6.5	113	0.00
32		Benzo(a)anthracene	0.400	0.385	3.8	97	0.00
33		Chrysene	0.400	0.422	-5.5	109	0.00
34		Bis(2-ethylhexyl)phthalate	0.400	0.417	-4.2	108	0.00
35		Indeno(1,2,3-cd)pyrene	0.400	0.526	-31.5#	139	0.00
36	I	Perylene-d12	0.400	0.400	0.0	112	0.00
37		Benzo(b)fluoranthene	0.400	0.377	5.8	102	0.00
38		Benzo(k)fluoranthene	0.400	0.423	-5.7	106	0.00
39	C	Benzo(a)pyrene	0.400	0.401	-0.3	111	0.00
40		Dibenzo(a,h)anthracene	0.400	0.487	-21.7	142	0.00
41		Benzo(g,h,i)perylene	0.400	0.400	0.0	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071020\
Data File : BN011132.D
Acq On : 09 Jul 2020 14:56
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 09 15:34:53 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN063020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 09 15:33:27 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