Monte Carlo Simulation Tools for REDD+ Uncertainty Estimates

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Contents

Objective	1
SimVoi Syntax	1
SimVoi Replication	3
Replication Results	8
Distribution analysis	9
Recommendations	14
References	14

Objective

Using an R-based approach, this analysis replicates the Monte Carlo simulations originally performed with SimVoi in Excel. It details the code used in the analysis, compares simulation results between R and Excel, and proposes next steps for enhancement based on statistical tests. Specifically, we identified updates needed to address empirical distributions of the input data. Such updates are recommended to both strengthen compliance with REDD+ requirements and VVB audits, and to provide an effective basis for reducing uncertainty and increasing revenue from emissions credits.

SimVoi Syntax

SimVoi provides seventeen random number generator functions¹ that are operated with the following syntax:

- RandBeta(alpha,beta,,[MinValue],[MaxValue])
- RandBinomial(trials, probability_s)
- RandBiVarNormal(mean1, stdev1, mean2, stdev2, correl12)
- RandCumulative(value_cumulative_table)
- RandDiscrete(value_discrete_table)
- RandExponential(lambda)
- RandInteger(bottom,top)
- RandLogNormal(Mean,StDev)
- RandNormal(mean, standard dev)
- RandPoisson(mean)
- RandSample(population)
- RandTriangular(minimum, most_likely, maximum)
- RandTriBeta(minimum, most_likely, maximum, [shape])
- RandTruncBiVarNormal(mean1, stdev1, mean2, stdev2, correl12, [min1], [max1], [min2], [max2])
- RandTruncLogNormal(Mean, StDev, [MinValue], [MaxValue])
- RandTruncNormal(Mean,StDev,[MinValue],[MaxValue])
- RandUniform(minimum, maximum)

In this workflow, we attempt to replicate the following SimVoi function as identified in Guyana's emissions workbook:

=[1]!randtruncnormal(CarbonStocks.B2,CarbonStocks.B3,0)

¹SimVoi-313-Guide.pdf

According to the documentation, RandTruncNormal() returns a random value from a truncated normal distribution, modeling an uncertain quantity with a bell-shaped density while excluding extreme tail values. If no simulation count is provided, the function repeatedly samples until a value falls between the specified minimum and maximum—or until it reaches 10,000 attempts. In the above example, only a minimum value of 0 is provided, so the default iteration limit is used.

$Import\ data$

```
workbook = "./data/art/GuyanaARTWorkbookMC-thru2022-April2024_values_V2.xlsx"

CarbonStocks = readxl::read_excel(workbook, "CarbonStocks") |>
    janitor::clean_names() |> mutate(across(where(is.numeric), ~ round(.x, 1)))

DegradationEF = readxl::read_excel(workbook, "Degradation EFs") |> janitor::clean_names()|>
    mutate(across(where(is.numeric), ~ round(.x, 1)))

ActivityData = readxl::read_excel(workbook, "Activity Data") |> janitor::clean_names() |>
    mutate(across(where(is.numeric), ~ round(.x, 1)))

flextable(head(CarbonStocks[, 1:8]))|>fontsize(size=8,part="all")

flextable(ActivityData[, 1:8])|>fontsize(size=8,part="all")
```

Table 1: Input values from CarbonStocks tabsheet

Statistic	AG Tree	BG Tree	Saplings	Standing Dead	Lying Dead	Litter	Sum w/o Soil	Soil
mean	205.8	48.3	3.7	2.6	8.6	3.3	272.3	58.7
std_dev	60.4	14.3	2.0	4.0	8.1	1.3	90.0	61.5
minimum	91.6	21.2	0.5	0.0	0.0	1.2	114.4	10.1
maximum	353.7	83.1	18.8	13.7	42.3	8.7	520.3	502.4
90%_CI	9.2	2.2	0.3	0.6	1.2	0.2	N/A	11.0
CI_%_of_mean	0.0	0.0	0.1	0.2	0.1	N/A	N/A	0.2

Table 2: Input values from Degradation EFs tabsheet

Statistic	LDF	Wood Density	LIF	For. Infrastr.	Mining	Mining Infrastr.	Infrast.
Factor_tC	1.05	0.4	46.87	NA	NA	NA	NA
StdDev_tC	0.68	0.03	8.08	NA	NA	NA	NA
CI90_tC	0.08	0	1.6	NA	NA	NA	NA
Factor_tCO2	3.85	1.47	171.84	NA	NA	NA	NA
StdDev_tCO2	2.4900000000000002	0.11	29.63	NA	NA	NA	NA
Cl90_tCO2	0.289999999999998	0.01	5.87	NA	NA	NA	NA
EF_tCO2_AD	NA	NA	NA	8.1	8.1	8.1	8.1
MAD_tCO2	NA	NA	NA	8.1	8.1	8.1	8.1

Table 3: Absolute input values from Activity Data tabsheet

drivers	units	x2011	x2012	x2013	x2014	x2015	x2016	x2017	x2018	x2019	x2020	x2021	x2022	x2023
Forestry infrastructure	ha	186	240	330	204	313	313	227	356	226	195	228	155.6	339
Agriculture	ha	41	440	424	817	379	379	477	512	246	489	216	281.6	475
Mining (medium and large scale) ha	7340	13664	11518	10434	6782	6782	7442	7624	5821	6,452	6,825	5,264.3	5,853
Mining infrastructure	ha													
Infrastructure	ha	118	127	342	141	217	217	195	67	52	102	117	110.6	541
Settlements	ha	-	-	23	71	8	8	7	7	22	60	105	169.4	201

SimVoi Replication

We utilize the replicate function to repeat a simulation assuming a truncated normal distribution.

$CarbonStocks\ data$

```
A_MEAN = CarbonStocks$ag_tree_t_c_ha[1]
         = CarbonStocks$ag_tree_t_c_ha[2]
   B_MEAN = CarbonStocks$bg_tree_t_c_ha[1]
   B SD = CarbonStocks$bg tree t c ha[2]
   C_MEAN = CarbonStocks$saplings_t_c_ha[1]
        = CarbonStocks$saplings t c ha[2]
   D_MEAN = CarbonStocks$standing_dead_wood_t_c_ha[1]
        = CarbonStocks$standing dead wood t c ha[2]
   E MEAN = CarbonStocks$lying dead wood t c ha[1]
        = CarbonStocks$lying_dead_wood_t_c_ha[2]
10
   F_MEAN = CarbonStocks$litter_t_c_ha[1]
   F_SD
         = CarbonStocks$litter_t_c_ha[2]
   G_MEAN = CarbonStocks$sum_pools_w_o_soil[1]
        = CarbonStocks$sum_pools_w_o_soil[2]
14
   H_MEAN = CarbonStocks$soil_t_c_ha[1]
        = CarbonStocks$soil_t_c_ha[2]
16
17
   # 100 simulations sampling 1 observation each iteration.
18
   # NOTE: #no. of simulations reduced to match estimates (see "Observations")
19
   A_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,91.6,353.7,A_MEAN, A_SD)
   B rtruncnormal 100 = truncnorm::rtruncnorm(n=100,21.2,83.1,B MEAN, B SD)
21
   C_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0.5,18.8,C_MEAN, C_SD)
   D_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0.0,13.7,D_MEAN, D_SD)
23
   E_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0.0,42.3,E_MEAN, E_SD)
   F rtruncnormal 100 = truncnorm::rtruncnorm(n=100,1.2,8.7,F MEAN,F SD)
25
   G_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,114.4,520.3,G_MEAN, G_SD)
   H_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,10.1,502.4,H_MEAN, H_SD)
27
   # --- Simulation Estimates
29
   AG_tree_tC_ha
                            = mean(A_rtruncnormal_100)
30
   AG_tree_tCO2_ha
                            = mean(A_rtruncnormal_100)*(44/12)
31
   BG_tree_tC_ha
                            = mean(B_rtruncnormal_100)
   BG_tree_tCO2_ha
                            = mean(B_rtruncnormal_100)*(44/12)
33
   Saplings_tC_ha
                             = mean(C_rtruncnormal_100)
34
   {\tt Saplings\_tCO2\_ha}
                            = mean(C_rtruncnormal_100)*(44/12)
35
   StandDead_tC_ha
                            = mean(D_rtruncnormal_100)
   StandDead tCO2 ha
                             = mean(D rtruncnormal 100)*(44/12)
```

```
LyingDead tC ha
                             = mean(E rtruncnormal 100)
                             = mean(E rtruncnormal 100)*(44/12)
   LyingDead_tCO2_ha
39
                             = mean(F rtruncnormal 100)
   Litter tC ha
40
                             = mean(F rtruncnormal 100)*(44/12)
   Litter_tCO2_ha
   Sum wo Soil tC ha
                             = mean(G rtruncnormal 100)
42
   Sum wo Soil tCO2 ha
                             = mean(G rtruncnormal 100)*(44/12)
   Soil tC ha
                             = mean(H rtruncnormal 100)
44
   Soil tCO2 ha
                             = mean(H rtruncnormal 100)*(44/12)
46
   CarbonStocks_MC_R_df <- data.frame(</pre>
                              = c("tC/ha", "tCO2/ha"),
48
      `AG Tree`
                              = c(AG_tree_tC_ha, AG_tree_tCO2_ha),
49
      `BG Tree`
                              = c(BG_tree_tC_ha, BG_tree_tCO2_ha),
50
                              = c(Saplings_tC_ha, Saplings_tCO2_ha),
     `Saplings`
51
      `Standing Dead`
                              = c(StandDead_tC_ha, StandDead_tCO2_ha),
52
                              = c(LyingDead_tC_ha, LyingDead_tCO2_ha),
      `Lying Dead`
53
     `Litter`
                              = c(Litter_tC_ha, Litter_tCO2_ha),
                              = c(Sum wo Soil tC ha, Sum wo Soil tCO2 ha),
     `Sum w/o Soil`
55
     `Soil`
                              = c(Soil tC ha, Soil tCO2 ha)
56
     )
   Degradation data
   A_MEAN = DegradationEF$ldf[4]
   A SD = DegradationEF$ldf[5]
   B MEAN = DegradationEF$wood density[4]
   B SD = DegradationEF$wood density[5]
   C_MEAN = DegradationEF$lif[4]
        = DegradationEF$lif[5]
   D_MEAN = DegradationEF$forestry_infrastructure[7]
        = DegradationEF$forestry_infrastructure[8]
   E_MEAN = DegradationEF$mining[7]
        = DegradationEF$mining[8]
   E SD
   F_MEAN = DegradationEF$mining_infrastructure[7]
11
         = DegradationEF$mining_infrastructure[8]
12
   G MEAN = DegradationEF$infrastructure[7]
13
        = DegradationEF$infrastructure[8]
14
1.5
   # 100 simulations sampling 1 observation each iteration.
   A rtruncnormal 100 = truncnorm::rtruncnorm(n=100,0,Inf,A MEAN, A SD)
17
   B_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0,Inf,B_MEAN, B_SD)
   C rtruncnormal 100 = truncnorm::rtruncnorm(n=100,0,Inf,C MEAN, C SD)
19
   D rtruncnormal 100 = truncnorm::rtruncnorm(n=100,0,Inf,D MEAN, D SD)
   E rtruncnormal 100 = truncnorm::rtruncnorm(n=100,0,Inf,E MEAN, E SD)
21
   F_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0,Inf,F_MEAN,F_SD)
   G_rtruncnormal_100 = truncnorm::rtruncnorm(n=100,0,Inf,G_MEAN, G_SD)
23
   LDF_EF_tCO2_m2
                             = mean(A rtruncnormal 100)
25
   WD_EF_tCO2_m2
                             = mean(B rtruncnormal 100)
26
                             = mean(C_rtruncnormal_100)
   LIF_EF_tCO2_km
27
                             = mean(D_rtruncnormal_100)
   ForInfr_EF_tCO2_ha
28
   Mining_EF_tCO2_ha
                             = mean(E_rtruncnormal_100)
                             = mean(F_rtruncnormal_100)
   MiningInfr_EF_tCO2_ha
30
   Infrastructure EF tCO2 ha= mean(G rtruncnormal 100)
31
32
```

```
df_logging <- data.frame(</pre>
33
      component = c("LDF", "Wood Density of Harvest", "LIF (Skid Trails)"),
34
                = c("per m3", "per m3", "per km"),
35
                = c(LDF_EF_tCO2_m2, WD_EF_tCO2_m2, LIF_EF_tCO2_km)
     tco2
37
   df degrading <- data.frame(degrading activity = c(</pre>
39
        "Forestry infrastructure", "Mining (medium & large scale)",
40
        "Mining infrastructure", "Infrastructure"),
41
     ef_tco2_ha = c(ForInfr_EF_tCO2_ha, Mining_EF_tCO2_ha,
42
        MiningInfr_EF_tCO2_ha, Infrastructure_EF_tCO2_ha)
43
44
45
   max_rows <- max(nrow(df_logging), nrow(df_degrading))</pre>
46
                         = rep(NA, max_rows - nrow(df_logging))
   logging_nas
47
   degrading_nas
                         = rep(NA, max_rows - nrow(df_degrading))
48
   Degradation_MC_R_df = data.frame(
     Component
                         = c(df logging$component, logging nas),
50
                         = c(df_logging$unit, logging_nas),
     Unit.
51
     tCO2
                         = c(df logging$tco2, logging nas),
52
     Degrading_Activity= c(df_degrading$degrading_activity, degrading_nas),
53
                         = c(df_degrading$ef_tco2_ha, degrading_nas)
     EF tCO2 ha
54
     )
```

Activity data

Please note that for purpose of saving space, columns and cells C1-M8 were implemented with echo-F settings, and can be located in the R-markdown.rmd file used to derive this PDF.

```
A1_MEAN = ActivityData$x2011[1]
          = ActivityData$x2011[17]
   A2_MEAN = ActivityData$x2011[2]
   A2 SD
          = ActivityData$x2011[18]
   A3_MEAN = ActivityData$x2011[3]
           = ActivityData$x2011[19]
   A3 SD
   A4_MEAN = ActivityData$x2011[4]
   A4 SD
           = ActivityData$x2011[20]
   A5 MEAN = ActivityData$x2011[5]
   A5 SD
          = ActivityData$x2011[21]
10
   A6_MEAN = ActivityData$x2011[6]
   A6 SD
           = ActivityData$x2011[22]
12
   A7 MEAN = ActivityData$x2011[7]
13
           = ActivityData$x2011[23]
   A7 SD
14
   A8 MEAN = ActivityData$x2011[8]
15
   A8_SD = ActivityData$x2011[24]
16
17
   B1_MEAN = ActivityData$x2012[1]
   B1_SD
           = ActivityData$x2012[17]
19
   B2_MEAN = ActivityData$x2012[2]
           = ActivityData$x2012[18]
21
   B3_MEAN = ActivityData$x2012[3]
   B3 SD
           = ActivityData$x2012[19]
23
   B4_MEAN = ActivityData$x2012[4]
          = ActivityData$x2012[20]
   B4 SD
   B5_MEAN = ActivityData$x2012[5]
```

```
B5 SD
           = ActivityData$x2012[21]
27
   B6 MEAN = ActivityData$x2012[6]
28
           = ActivityData$x2012[22]
29
   B7_MEAN = ActivityData$x2012[7]
   B7 SD
           = ActivityData$x2012[23]
31
   B8 MEAN = ActivityData$x2012[8]
           = ActivityData$x2012[24]
   B8 SD
33
   A1_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A1_MEAN, A1_SD)
35
   A2_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A2_MEAN, A2_SD)
   A3_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A3_MEAN, A3_SD)
37
   A4_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A4_MEAN, A4_SD)
   A5_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A5_MEAN, A5_SD)
39
   A6_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A6_MEAN, A6_SD)
40
   A7_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A7_MEAN, A7_SD)
   A8_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, A8_MEAN, A8_SD)
42
43
   B1 rtruncnormal 100 = truncnorm::rtruncnorm(n=100, 0, Inf, B1 MEAN, B1 SD)
44
   B2_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, B2_MEAN, B2_SD)
45
   B3 rtruncnormal 100 = truncnorm::rtruncnorm(n=100, 0, Inf, B3 MEAN, B3 SD)
46
   B4_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, B4_MEAN, B4_SD)
   B5 rtruncnormal 100 = truncnorm::rtruncnorm(n=100, 0, Inf, B5 MEAN, B5 SD)
48
   B6_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, B6_MEAN, B6_SD)
   B7_rtruncnormal_100 = truncnorm::rtruncnorm(n=100, 0, Inf, B7_MEAN, B7_SD)
50
   B8 rtruncnormal 100 = truncnorm::rtruncnorm(n=100, 0, Inf, B8 MEAN, B8 SD)
52
   # --- Simulation Estimates ---
   A1 = mean(A1_rtruncnormal_100)
54
   A2 = mean(A2_rtruncnormal_100)
   A3 = mean(A3 rtruncnormal 100)
   A4 = mean(A4_rtruncnormal_100)
   A5 = mean(A5_rtruncnormal_100)
   A6 = mean(A6_rtruncnormal_100)
   A7 = mean(A7_rtruncnormal_100)
   A8 = mean(A8 rtruncnormal 100)
61
62
   B1 = mean(B1 rtruncnormal 100)
63
   B2 = mean(B2_rtruncnormal_100)
   B3 = mean(B3 rtruncnormal 100)
65
   B4 = mean(B4_rtruncnormal_100)
   B5 = mean(B5 rtruncnormal 100)
67
   B6 = mean(B6 rtruncnormal 100)
   B7 = mean(B7 rtruncnormal 100)
69
   B8 = mean(B8_rtruncnormal_100)
```

Below, we presented organize our results into a table and compare with absolute input values (CarbonStocks), and Monte Carlo estimates generated with SimVoi (CarbonStocks (MC)). For external comparisons, we have also saved these results in a new excel tab called "CarbonStocks (MC-R)".

```
1   CarbonStocks_MC_R = flextable(head(CarbonStocks_MC_R_df[, 1:8])) |>
2   fontsize(size = 8, part = "all")
3   CarbonStocks_MC_R
4
5   Degradation_MC_R = flextable(head(Degradation_MC_R_df[, 1:8])) |>
6   fontsize(size = 8, part = "all")
```

7 Degradation_MC_R

Table 4: Results of Monte Carlo simulations of CarbonStocks tabsheet using R

- carbonStocks_MC_R = flextable(CarbonStocks_MC_R_df) |>
- width(width = 1) |> fit_to_width(max_width = 6) |>
- colformat_double(big.mark = ",", digits = 1, na_str = "N/A")
- 4 CarbonStocks_MC_R

Units	AG.Tree	BG.Tree	Saplings	Standing.Dead	Lying.Dead	Litter	Sum.w.o.Soil	Soil
tC/ha	211.1	47.2	3.9	4.2	11.1	3.5	274.5	87.1
tCO2/ha	774.2	173.0	14.5	15.6	40.6	12.7	1,006.6	319.3

Table 5: Results of Monte Carlo simulations of CarbonStocks tabsheet using SimVoi

Units	AG Tree	BG Tree	Saplings	Standing Dead	Lying Dead	Litter	Sum w/o Soil	Soil
tC/ha	181.1	65.0	3.5	7.3	17.1	3.7	277.7	60.6
tCO2/ha	664.2	238.2	12.8	26.9	62.6	13.7	1,018.4	222.3

Table 6: Results of Monte Carlo simulations of DegradationEFs tabsheet using R

Component	Unit	tCO2	Degrading_Activity	EF_tCO2_ha
LDF	per m3	4.2	Forestry infrastructure	10.7
Wood Density of Harvest	per m3	1.5	Mining (medium & large scale)	10.4
LIF (Skid Trails)	per km	176.1	Mining infrastructure	9.8
		N/A	Infrastructure	9.7

Table 7: Results of Monte Carlo simulations of Degradation EFs tabsheet using SimVoi

Component	Unit	t CO2	Degrading Activity	EF (tCO2/ha)
LDF	per m ³	4.54	Forestry infrastructure	7.2
Wood Density of timber harvested	per m³	1.39	Mining (med & large scale)	7.2
LIF (Skid Trails)	per km	185.88	Mining infrastructure	7.2
			Infrastructure	7.2

Table 8: Results of Monte Carlo simulations of Activity Data tabsheet using R

Drivers	units	X2011	X2012	X2013	X2014	X2015
Forestry infrastructure	ha	187	243	337	210	315
Agriculture	ha	40	452	426	815	381
Mining (med & large scale)	ha	7,256	13,578	11,408	10,347	6,816
Mining infrastructure	ha	N/A	N/A	N/A	N/A	N/A
Infrastructure	ha	121	129	341	137	215
Settlements	ha	N/A	N/A	23	73	8

Drivers	units	X2011	X2012	X2013	X2014	X2015
Fire-Biomass burning	ha	47	183	96	258	1,501
Shifting Cultivation	ha	N/A	N/A	N/A	N/A	N/A

Table 8: Results of Monte Carlo simulations of Activity Data tabsheet using SimVoi

Drivers	units	2011	2012	2013	2014	2015	2016	2017	2018	2019	2020	2021	2022
Forestry infrastructure		225	194	270	229	325	288	224	244	228	185	215	83
Agriculture		36	384	463	882	436	401	487	291	261	126	436	305
Mining (med & large scale)		8,835	13,157	7,687	12,583	7,673	7,082	6,619	6,295	5,518	6,980	7,303	6,163
Infrastructure	ha	117	121	331	143	182	236	196	72	33	188	101	117
Settlements		-	-	25	73	7	7	7	9	29	53	110	210
Fire-Biomass burning		53	188	89	301	1,833	1,553	460	681	6,194	3,364	221	365
Shifting Cultivation								451	361	509	478	456	109
Deforestation		9,266	14,044	8,863	14,210	10,457	9,568	7,993	7,592	12,772	11,373	8,842	7,353
Logging - harvest volume	m3	608,730	585,108	624,287	759,684	655,406	500,788	533,106	546,242	521,172	545,355	547,516	547,517
Logging - skid trail length	km	2,302	2,212	2,360	2,872	2,478	1,893	2,016	2,065	1,971	2,062	2,070	2,070
Illegal logging	m3	2776	2,306	2,371	2,836	1,505	2,249	2,706	3,719	2,149	1,281	1,281	1,281
Mining and Infrastructure	ha						36,647	31,919	28,185	23,028	22,795	26,651	26,651

Replication Results

In the following chunk, we compute the final uncertainty estimates for the simulated emission reductions (GHG ER) using a truncated normal distribution. With 10,000 simulation trials, we calculate key statistics, mean, standard deviation, and mean standard error, along with distributional percentiles. By extracting the 5th and 95th percentiles, we form a 90% confidence interval (CI) from which we derive the margin of error (ME) and its corresponding percentage error relative to the mean. These computed metrics replicate the SimVoi univariate summary and provide a statistical summary of the uncertainty associated with the emissions reduction estimates.

```
# Emission Reductions (MC-R)
ER_values <- c(7715885, 10371977, 10040723, 6358705, 7174999, 6977178, 9223423, 7299024)
ER mean emp <- mean(ER values)</pre>
ER_sd_emp
           <- sd(ER_values)
n_sim <- 10000
# Simulate a truncated normal distribution & compute stats
sim_ER <- rtruncnorm(n = n_sim, a = 0, b = Inf, mean = ER_mean_emp, sd = ER_sd_emp)
sim_mean <- mean(sim_ER)</pre>
sim_sd <- sd(sim_ER)
sim se
       <- sim_sd / sqrt(length(sim_ER))
sim_skew <- moments::skewness(sim_ER)</pre>
sim_quant <- quantile(sim_ER, probs = c(0, 0.25, 0.5, 0.75, 1))
lower90 <- quantile(sim ER, probs = 0.05)</pre>
upper90 <- quantile(sim ER, probs = 0.95)
ci90 <- upper90 - lower90
ME \leftarrow ci90 / 2
pct_error <- ME / sim_mean * 100</pre>
ER summary <- data.frame(</pre>
  Metric = c("Mean", "St. Dev.", "Mean St. Error", "Skewness",
             "Minimum", "1st Quartile", "Median", "3rd Quartile", "Maximum",
             "5th Percentile", "95th Percentile", "90% CI", "Margin of Error", "% Error"),
  Value = c(round(sim_mean),
                                   round(sim_sd),
             round(sim_se),
                                   round(sim_skew, 3),
```

```
round(sim_quant[1]), round(sim_quant[2]),
round(sim_quant[3]), round(sim_quant[4]),
round(sim_quant[5]), round(lower90),
round(upper90), round(ci90),
round(ME), round(pct_error, 2))
)
```

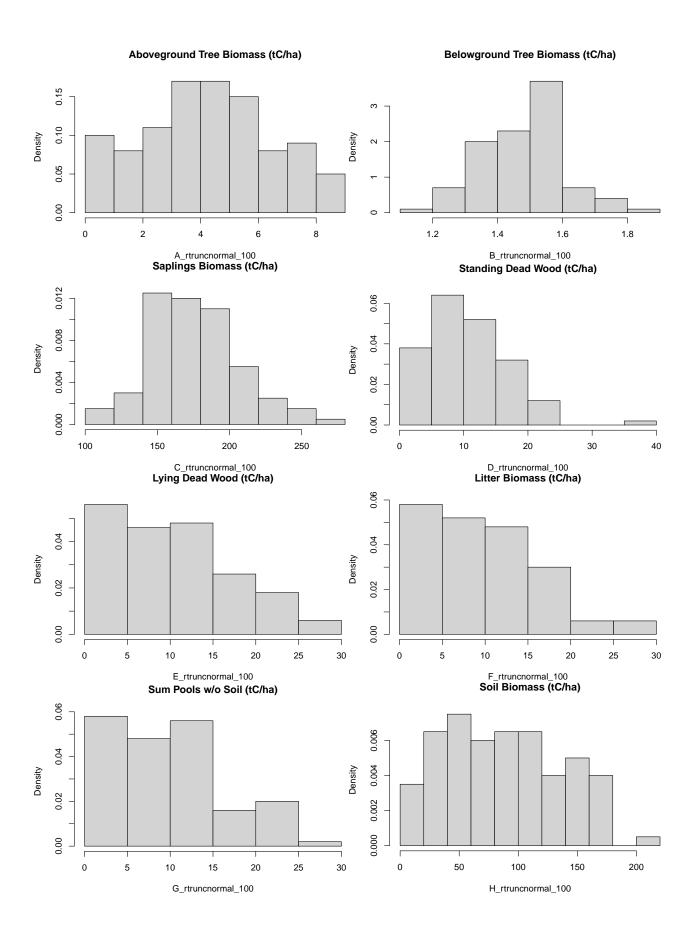
Table 9: Simulated Univariate Summary of Emission Reductions (GHG ER)

Metric	Value
Mean	8,150,375.000
St. Dev.	1,518,508.000
Mean St. Error	15,185.000
Skewness	0.018
Minimum	2,526,130.000
1st Quartile	7,104,729.000
Median	8,143,815.000
3rd Quartile	9,176,853.000
Maximum	13,630,369.000
5th Percentile	5,682,413.000
95th Percentile	10,658,760.000
90% CI	4,976,346.000
Margin of Error	2,488,173.000
% Error	30.530

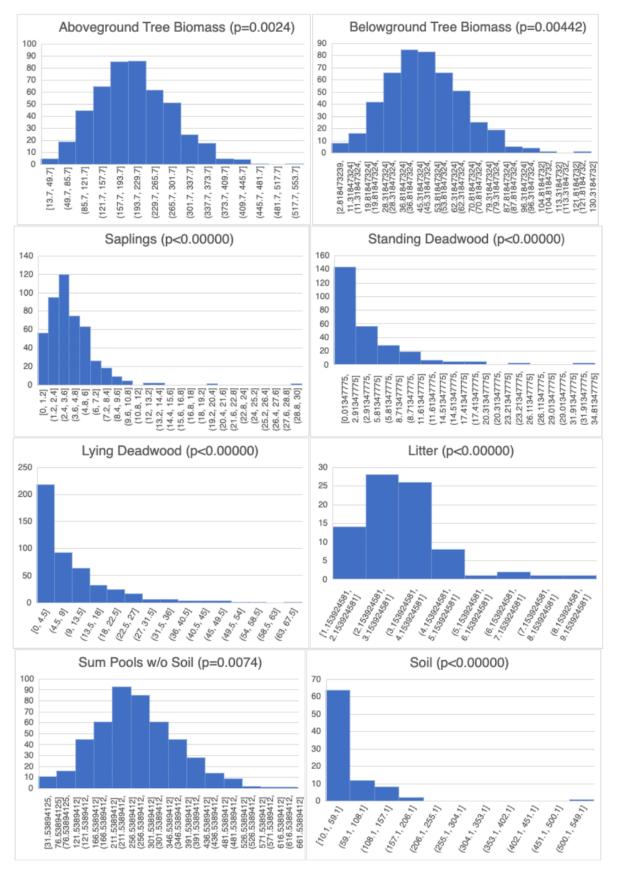
Distribution analysis

This distribution analysis is a critical step in our Monte Carlo simulation process (Bhanti, 2025). It visually and statistically confirms that the simulations accurately reflect the empirical characteristics of the input data. Normality tests (Shapiro, Sandford & Wilk, 1965) and visualizations below (Figures 9-16) guide necessary parameter adjustments and monitor for systematic errors. Failure to perform this step risks masking or exaggerating inherent biases, which could undermine the reliability of our estimates. Figures 1–8 display the simulated distributions, while Figures 9–16 show the original data distributions. Discrepancies between them highlight opportunities for effectively reducing uncertainty. We recommend to refine the current simulation approach by incorporating data-specific models fitted with log-normal or beta distributions.

```
# Distribution of Monte Carlo derived estimates
hist(A_rtruncnormal_100, freq=F, main="Aboveground Tree Biomass (tC/ha)")
hist(B_rtruncnormal_100, freq=F, main="Belowground Tree Biomass (tC/ha)")
hist(C_rtruncnormal_100, freq=F, main="Saplings Biomass (tC/ha)")
hist(D_rtruncnormal_100, freq=F, main="Standing Dead Wood (tC/ha)")
hist(E_rtruncnormal_100, freq=F, main="Lying Dead Wood (tC/ha)")
hist(F_rtruncnormal_100, freq=F, main="Litter Biomass (tC/ha)")
hist(G_rtruncnormal_100, freq=F, main="Sum Pools w/o Soil (tC/ha)")
hist(H_rtruncnormal_100, freq=F, main="Soil Biomass (tC/ha)")
```



Figures 1-8: Distribution analysis of simulated estimates of carbon stock variables



Figures 9-16: Distribution analysis of carbon stock input variables (Bhanti, 2025-03-17)

When preparing for Monte Carlo simulations, it is best practice to start by examining descriptive statistics to characterize the empirical distributions of input variables. This preliminary analysis typically includes statistical tests for normality and visualizations of univariate distributions, such as histograms, kernel density plots, and Q-Q plots. Together, these tools provide critical insights into the shape, spread, symmetry, skewness, and presence of potential outliers in the data. Although this preliminary step may seem minor, it substantially influences uncertainty estimates, which can directly translate into increased financial returns, particularly within forest project landscapes exhibiting non-normal data distributions.

Accurately characterizing data distributions also helps in identifying and addressing biases, thereby ensuring high data quality and increasing confidence in subsequent estimations of biomass and carbon emissions. Selecting appropriate statistical distributions, informed by exploratory analyses, significantly enhances the reliability and precision of Monte Carlo simulations. Consequently, such careful statistical characterizations reduce overall uncertainty in forest biomass and emissions estimates. In turn, this strengthens the credibility of jurisdictional claims made under REDD+ programs and maximizes potential financial returns for Guyana from carbon financing initiatives.

Univariate distribution visualizations additionally provide auditors with useful diagnostic resources, enabling rapid identification and characterization of biases commonly encountered in biomass data. These diagrams help auditors efficiently assess the technical rigor and statistical approaches implemented by the project to monitor and manage uncertainty (ART, 2021: 8). Winrock strongly recommends incorporating distribution analyses early in a project's quantitative planning and throughout its technical standard operating procedures (SOPs). Such early integration represents a low hanging fruit with cost-effective strategy and significant potential in reducing audit findings, lowering uncertainty, and enhancing financial outcomes for Guyana's REDD+ activities. Specifically, early attention to data distributions directly informs appropriate simulation selection from the available options in SimVoi.

To effectively guide practitioners and stakeholders in selecting appropriate statistical distributions for Monte Carlo methods within forestry and REDD+ contexts, the following two tables present findings from a rapid review of relevant literature. The review identified and summarized statistical distributions frequently encountered in forestry, biomass estimation, and emissions analysis, which are dis-aggregated below between discrete and continuous types and according to their inherent statistical characteristics.

Table 9: Continuous data distributions, example cases $\mathfrak E$ equations used in Monte Carlo simulations.

Distribution	Statistical Use Cases	PDF
Normal	Symmetric, bell-shaped distribution used for modeling continuous variables: biomass/ha	$f(x) = \frac{1}{\sigma\sqrt{2\pi}} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right)$
Lognormal	Right-skewed distribution suitable for variables constrained to positive values (e.g., emission rates).	$f(x) = \frac{1}{x\sigma\sqrt{2\pi}}\exp\left(-\frac{(\ln x - \mu)^2}{2\sigma^2}\right)$
Exponential	Models waiting times between independent events, such as forest fire occurrences or logging events.	$f(x) = \lambda e^{-\lambda x}, \\ x \ge 0$
Cont. Uniform	Assumes all values in an interval [a, b] are equally likely; useful for random spatial sampling in forests.	$f(x) = \frac{1}{b-a},$ $a < x < b$
Chi-Square	Often used in goodness-of-fit tests to evaluate model accuracy in biomass estimation.	$f(x) = \frac{1}{2^{k/2} \Gamma(k/2)} x^{\frac{k}{2} - 1} e^{-x/2}, x > 0$
t-Distribution	Suitable for small sample sizes with unknown population stdev (e.g., limited forest carbon data).	$f(x) = \frac{\Gamma\left(\frac{v+1}{2}\right)}{\sqrt{\nu\pi}\Gamma\left(\frac{v}{2}\right)} \left(1 + \frac{x^2}{v}\right)^{\frac{v+1}{2}}$
Gamma	Models positively skewed data, such as biomass growth rates or carbon accumulation over time.	$f(x) = \frac{x^{k-1}e^{-x/\theta}}{\theta^k \Gamma(k)}$
Weibull	Flexible distribution used in reliability analysis, e.g., modeling tree mortality.	$f(x) = \frac{k}{\lambda} \left(\frac{x}{\lambda}\right)^{k-1} e^{-(x/\lambda)^k}$

Table 10: Discrete data distributions, example cases $\mathfrak E$ equations used in Monte Carlo simulations.

Distribution	Statistical Use Cases	PMF
Bernoulli	Binary outcome probability, e.g., presence/absence of deforestation in an area.	$P(X = x) = p^{x}(1-p)^{1-x},$ $x \in \{0,1\}$
Binomial	Probability of fixed #no. of successes over n Bernoulli trials, e.g., no. of heads in ten coin flips.	$P(X = k) = \binom{n}{k} p^k (1 - p)^{n-k},$ k = 0, 1,, n
Poisson	Models counts of independent events within an interval, e.g., number of wildfire incidents per year.	$P(X = k) = \frac{\lambda^k e^{-\lambda}}{k!},$ k = 0,1,2,
Geometric	Models #no. of trials until the first success, e.g., number of inspections until detecting deforestation.	$P(X = k) = (1 - p)^{k-1}p,$ k = 1,2,
Negative Binomial	Counts #no. failures until r successes occur, treats overdispersed or repeated deforestation detections.	$P(X = k) = {k + r - 1 \choose k} (1 - p)^r p^k,$ k = 0,1,2,
Discrete Uniform	Assumes outcome in a finite set is equally likely, e.g., random sampling of inventory across a forest.	$P(X = x) = \frac{1}{n},$ $x = 1,2,,n$

Discrete distributions describe forestry monitoring scenarios where data outcomes are countable and finite. Common examples include the number of deforestation events, occurrences of wildfires, or counts of logged trees within a defined monitoring interval. Accurate representation of discrete events using appropriate distributions such as Binomial, Poisson, or Negative Binomial significantly enhances the accuracy of model predictions and uncertainty assessments. For instance, employing a Poisson distribution to model occurrences of illegal logging events can improve the precision of estimated deforestation emissions and reduce uncertainty around compliance risks.

In contrast, continuous distributions capture variables capable of taking any value within a specified range and are particularly relevant in forestry when modeling measurements such as tree heights, carbon stock densities, or biomass values. Continuous distributions like the Normal (Gaussian), Lognormal, Weibull, and

Gamma distributions frequently arise in ecological modeling and biomass estimations due to their ability to realistically represent ecological variability and complex environmental factors. For example, using a Lognormal distribution for tree biomass data often provides more reliable estimates, particularly when the dataset is right-skewed due to natural variability in tree growth and forest conditions.

Central to these distributions are two mathematical concepts: Probability Mass Functions (PMFs) for discrete data and Probability Density Functions (PDFs) for continuous data. PMFs allocate specific probabilities to discrete outcomes, essential for accurately simulating events such as species occurrences or forest disturbances. PDFs describe the relative likelihood of continuous data points, enabling the robust estimation of variables like forest carbon content or annual biomass increment.

In Monte Carlo simulations, precise definition and utilization of PMFs and PDFs are crucial. These functions underpin random sampling processes that directly influence the reliability, precision, and credibility of uncertainty estimates. Given that forestry data is known to exhibit non-normal distributions due to inherent ecological heterogeneity that, informed selection and rigorous application of these functions are vital. Accurate modeling of the underlying data distribution enhances biomass and emissions estimates, significantly reduces uncertainty, and bolsters the financial and ecological credibility of REDD+ reporting initiatives (Morgan & Henrion, 1990; IPCC, 2019; ART, 2021).

Practitioners are encouraged to conduct exploratory data analysis early in their project planning stages, integrating statistical tests of normality and visual assessments (histograms, kernel density plots, Q-Q plots). Such preliminary analyses assist in diagnosing data distributions accurately, improving model selection, reducing potential auditor findings, and ultimately enhancing the financial and environmental outcomes of national REDD+ monitoring programs.

Recommendations

- Regarding simulation parameters, "RandTruncNormal(Mean,StDev,MinValue,MaxValue)) uses values of RandNormal until a value is found between MinValue and MaxValue or until it has made 10,000 attempts." The formula applied in Guyana's verified workbook uses a minimum value of 0. Considering the levels of quantitative variance between input values in the "CarbonStocks" tabsheet and SimVoisimulated estimates in "CarbonStocks (MC)" tabsheet, we may assume that approximately less than ~100 simulations were completed before values were reached "between MinValue and MaxValue". It is useful to consider that with increasing numbers of simulations the closer the mean values generated should be to absolute input values. Therefore, in order to test this assumption we simply ran 10,000 simulations, which generated mean estimates of within 2 tC ha-1 of input values (205 tC ha-1) 10 times of 10 reruns. By reducing the number of iterations per simulation, we confirm our assumptions that the current SimVoi setup provided estimates from ~<100 simulations.
- We recommend increasing the number of simulation iterations from 100 to a higher number, in order to reduce variance in and between simulated estimates, which thereby provide more robust uncertainty quantification.
- It is advised to consider performing sensitivity analyses in order to identify which input parameters contribute most to overall uncertainty, thereby informing targeted data collection and model refinement efforts.
- Furthermore, distribution analysis confirms that our simulation outputs approximate the expected statistical distributions of the underlying data. In the next phase, we recommend incorporating the observed distributional characteristics into the Monte Carlo simulation framework. This adjustment—using distribution-specific models (e.g., log-normal, beta, etc.)—is essential for reducing uncertainty and ensuring that our estimates carry statistical significance.

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Annex I: SimVoi replication in base-R

Using the following base functions from R avoids need for any package installations and limits risk of bugs or common problems with R versioning arising in the workflow. However, the base functions below are a little clunkier and may require additional guidance. Regarding R and package versions, please consult the Annex II which presents record of the current runtime used in this analysis.

```
randtruncnormal_sim_10000 <- rnorm(n=10,mean=MEAN,sd=SD)
hist(randtruncnormal_sim_10000, freq=F)
AG_Tree_tC_ha
               = mean(randtruncnormal_sim_10000)
AG_Tree_tCO2_ha = AG_Tree_tC_ha*(44/12)
AG_Tree_tC_ha
AG Tree tCO2 ha
#curve(dnorm(x, mean=MEAN, sd=SD), from=0, to=450, add=T, col="red")
# 10,000 simulations sampling 10 observations
randtruncnormal_sim_10000_10 = replicate(n=10000, rnorm(n=10,mean=MEAN,sd=SD))
hist(apply(X = randtruncnormal_sim_10000_10, MARGIN=2, FUN=mean))
sd(apply(X = randtruncnormal sim 10000 10, MARGIN=2, FUN=mean))
mean(apply(X = randtruncnormal sim 10000 10, MARGIN=2, FUN=mean))
(mean(apply(X = randtruncnormal sim 10000 10, MARGIN=2, FUN=mean)))*(44/12)
# 10,000 simulations sampling 100 observations
randtruncnormal_sim_10000_100 = replicate(n=10000,rnorm(n=100,mean=MEAN,sd=SD))
hist(apply(X = randtruncnormal_sim_10000_100, MARGIN=2, FUN=mean))
sd(apply(X = randtruncnormal_sim_10000_100, MARGIN=2, FUN=mean))
mean(apply(X = randtruncnormal_sim_10000_100, MARGIN=2, FUN=mean))
(mean(apply(X = randtruncnormal_sim_10000_100, MARGIN=2, FUN=mean)))*(44/12)
# 10,000 simulations sampling 1,000 observations
randtruncnormal sim 10000 1000 = replicate(n=10000, rnorm(n=1000, mean=MEAN, sd=SD))
hist(apply(X = randtruncnormal sim 10000 1000, MARGIN=2, FUN=mean))
sd(apply(X = randtruncnormal_sim_10000_1000, MARGIN=2, FUN=mean))
mean(apply(X = randtruncnormal sim 10000 1000, MARGIN=2, FUN=mean))
(mean(apply(X = randtruncnormal_sim_10000_1000, MARGIN=2, FUN=mean)))*(44/12)
# 10,000 simulations sampling 10,000 observations
randtruncnormal_sim_10000_10000 = replicate(n=10000,rnorm(n=10000,mean=MEAN,sd=SD))
hist(apply(X = randtruncnormal_sim_10000_10000, MARGIN=2, FUN=mean))
sd(apply(X = randtruncnormal_sim_10000_10000, MARGIN=2, FUN=mean))
mean(apply(X = randtruncnormal_sim_10000_10000, MARGIN=2, FUN=mean))
(mean(apply(X = randtruncnormal_sim_10000_10000, MARGIN=2, FUN=mean)))*(44/12)
```

Annex II: Rapid literature review of Monte Carlo methods in REDD+

Table A.2: Search parameters, resource scope, and objectives informing search

$\overline{\mathbf{REDD}}$ +2	MC Application	Region	Key Findings
ADD	Uncertainty of SAAB	Rondônia, Brazil	Estimated \pm 20%
	estimate		measurement error in
			SAAB using Monte
			Carlo simulations;
			emphasized large trees'
			role in biomass.

²1. ADD: Avoided deforestation degradation, IFM: Improved forest management, JNR: Jurisdictional nested REDD+

ADD	AGB Uncertainty	Kenya, Mozambique	Assessed mixed-effects models in estimating
ADD	Blanket uncertainty propagation	Ghana	mangrove biomass. AGB prediction error >20%; addressed error propagation from trees to pixels in remote sensing.
ADD	Plot-based uncertainty	New Zealand	Cross-plot variance greatest magnitude of uncertainty
JNR	Multi-scale AGB uncertainty modeling	Minnesota, USA	Cross-scale tests showing effects of spatial resolution on AGB uncertainty.
N/A	Allometric uncertainty modeling	Panama	Allometric models identified as largest source of biomass estimation error.
ADD	Sampling and allometric uncertainty	Tapajos Nat Forest, Brazil	Significance of allometric models on uncertainty of root biomass, 95% CI, 21 plots.
ADD	Uncertainty of volume estimates	Santa Catarina, Brazil	Negligible effects of residual uncertainty on large-area estimates
N/A	Uncertainty metrics in model selection	Oregon, USA	Uncertainty estimates call for local validation or new local model development
ADD	AGB model uncertainty	French Guiana	AGB sub-model errors dominate uncertainty; height and wood-specific gravity errors are minor but can cause bias.
IFM	Emission factor uncertainty	Central Africa	Model selection is the largest error source (40%); weighting models reduces uncertainty in emission factors.
NA	Uncertainty in ecosystem nutrient estimate	New Hampshire, USA	Identified 8% uncertainty in nitrogen budgets, mainly from plot variability (6%) and allometric errors (5%).

${\bf Annex~III:~Runtime~snapshot}$

devtools::session_info()

⁻ Session info -----setting value
version R version 4.3.0 (2023-04-21)

os macOS 15.3.2 system aarch64, darwin20

ui X11 language (EN)

collate en_US.UTF-8
ctype en_US.UTF-8
tz America/Vancouver

date 2025-03-25

pandoc 3.6.1 @ /usr/local/bin/ (via rmarkdown)

_	- Packages									
	package	*	version	date	(UTC)	lib	sourc	:e		
	animation		2.7	2021-	10-07	[1]	CRAN	(R	4.3.3)	
	askpass		1.2.1						4.3.3)	
	assertthat		0.2.1	2019-	03-21	[1]	CRAN	(R	4.3.0)	
	backports		1.5.0	2024-	05-23	[1]	CRAN	(R	4.3.3)	
	BIOMASS	*	2.2.3	2025-	02-24	[1]	CRAN	(R	4.3.3)	
	boot		1.3-31	2024-	-08-28	[1]	CRAN	(R	4.3.3)	
	broom	*	1.0.7	2024-	09-26	[1]	CRAN	(R	4.3.3)	
	c2z	*	0.2.0	2023-	-08-10	[1]	CRAN	(R	4.3.0)	
	cachem		1.1.0	2024-	05-16	[1]	CRAN	(R	4.3.3)	
	caret	*	7.0-1	2024-	12-10	[1]	CRAN	(R	4.3.3)	
	cellranger		1.1.0	2016-	07-27	[1]	CRAN	(R	4.3.0)	
	chromote		0.4.0	2025-	01-25	[1]	CRAN	(R	4.3.3)	
	class		7.3-23	2025-	01-01	[1]	CRAN	(R	4.3.3)	
	classInt		0.4-11	2025-	-01-08	[1]	CRAN	(R	4.3.3)	
	cli		3.6.3	2024-	06-21	[1]	CRAN	(R	4.3.3)	
	codetools		0.2-20	2024-	-03-31	[1]	CRAN	(R	4.3.1)	
	colorspace		2.1-1	2024-	07-26	[1]	CRAN	(R	4.3.3)	
	data.table		1.16.4	2024-	12-06	[1]	CRAN	(R	4.3.3)	
	dataMaid	*	1.4.1	2021-	10-08	[1]	CRAN	(R	4.3.0)	
	DBI		1.2.3	2024-	-06-02	[1]	CRAN	(R	4.3.3)	
	DEoptimR		1.1-3-1	2024-	11-23	[1]	CRAN	(R	4.3.3)	
	DescTools	*	0.99.59	2025-	01-26	[1]	CRAN	(R	4.3.3)	
	devtools		2.4.5	2022-	10-11	[1]	CRAN	(R	4.3.0)	
	dials	*	1.3.0	2024-	07-30	[1]	CRAN	(R	4.3.3)	
	DiceDesign		1.10	2023-	12-07	[1]	CRAN	(R	4.3.1)	
	digest		0.6.37	2024-	08-19	[1]	CRAN	(R	4.3.3)	
	dplyr	*	1.1.4	2023-	11-17	[1]	CRAN	(R	4.3.1)	
	e1071		1.7-16	2024-	-09-16	[1]	CRAN	(R	4.3.3)	
	easypackages		0.1.0	2016-	12-05	[1]	CRAN	(R	4.3.0)	
	ellipsis		0.3.2	2021-	04-29	[1]	CRAN	(R	4.3.0)	
	evaluate		1.0.3	2025-	01-10	[1]	CRAN	(R	4.3.3)	
	Exact		3.3	2024-	07-21	[1]	CRAN	(R	4.3.3)	
	expm		1.0-0	2024-	08-19	[1]	CRAN	(R	4.3.3)	
	extrafont	*	0.19	2023-	01-18	[1]	CRAN	(R	4.3.3)	
	extrafontdb		1.0	2012-	06-11	[1]	CRAN	(R	4.3.3)	
	fastmap		1.2.0	2024-	05-15	[1]	CRAN	(R	4.3.3)	
	flextable	*	0.9.7	2024-	10-27	[1]	CRAN	(R	4.3.3)	
	${\tt fontBitstreamVera}$		0.1.1	2017-	02-01	[1]	CRAN	(R	4.3.3)	
	fontLiberation		0.1.0	2016-	10-15	[1]	CRAN	(R	4.3.3)	
	fontquiver		0.2.1	2017-	02-01	[1]	CRAN	(R	4.3.3)	
	forcats	*	1.0.0	2023-	01-29	[1]	CRAN	(R	4.3.0)	
	foreach		1.5.2	2022-	02-02	[1]	CRAN	(R	4.3.0)	

formatR	*	1.14	2023-01-17	[1]	CRAN	(R 4.3.3)
fs		1.6.5	2024-10-30	[1]	CRAN	(R 4.3.3)
furrr		0.3.1	2022-08-15	[1]	CRAN	(R 4.3.0)
future		1.34.0	2024-07-29	[1]	CRAN	(R 4.3.3)
future.apply		1.11.3	2024-10-27	[1]	CRAN	(R 4.3.3)
gdtools		0.4.1	2024-11-04	[1]	CRAN	(R 4.3.3)
generics		0.1.3	2022-07-05	[1]	CRAN	(R 4.3.0)
-	Ψ.	3.5.1	2024-04-23	[1]	CRAN	(R 4.3.0)
ggplot2 gld	т	2.6.7	2025-01-17	[1]	CRAN	(R 4.3.1)
0		0.16.3	2023-01-17	[1]	CRAN	(R 4.3.3)
globals						(R 4.3.1)
glue		1.8.0	2024-09-30 2024-12-17	[1] [1]	CRAN	
gower		1.0.2			CRAN	(R 4.3.3)
GPfit		1.0-8	2019-02-08	[1]	CRAN	(R 4.3.0)
gridExtra		2.3	2017-09-09	[1]	CRAN	(R 4.3.0)
gtable		0.3.6	2024-10-25	[1]	CRAN	(R 4.3.3)
hardhat		1.4.0	2024-06-02	[1]	CRAN	(R 4.3.3)
haven		2.5.4	2023-11-30	[1]	CRAN	(R 4.3.1)
hms		1.1.3	2023-03-21	[1]	CRAN	(R 4.3.0)
htmltools	*	0.5.8.1	2024-04-04	[1]	CRAN	(R 4.3.1)
htmlwidgets		1.6.4	2023-12-06	[1]	CRAN	(R 4.3.1)
httpuv		1.6.15	2024-03-26	[1]	CRAN	(R 4.3.1)
httr		1.4.7	2023-08-15	[1]	CRAN	(R 4.3.0)
infer	*	1.0.7	2024-03-25	[1]	CRAN	(R 4.3.1)
ipred		0.9-15	2024-07-18	[1]	CRAN	(R 4.3.3)
iterators		1.0.14	2022-02-05	[1]	CRAN	(R 4.3.0)
janitor	*	2.2.1	2024-12-22	[1]	CRAN	(R 4.3.3)
jsonlite	*	1.8.9	2024-09-20	[1]	CRAN	(R 4.3.3)
kableExtra	*	1.4.0	2024-01-24	[1]	CRAN	(R 4.3.1)
kernlab	*	0.9-33	2024-08-13	[1]	CRAN	(R 4.3.3)
KernSmooth		2.23-26	2025-01-01	[1]	CRAN	(R 4.3.3)
knitr	*	1.49	2024-11-08	[1]	CRAN	(R 4.3.3)
later		1.4.1	2024-11-27	[1]	CRAN	(R 4.3.3)
latex2exp	*	0.9.6	2022-11-28	[1]	CRAN	(R 4.3.0)
latexpdf	*	0.1.8	2023-12-19	[1]	CRAN	(R 4.3.3)
lattice	*	0.22-6	2024-03-20	[1]	CRAN	(R 4.3.1)
lava		1.8.1	2025-01-12	[1]	CRAN	(R 4.3.3)
lhs		1.2.0	2024-06-30	[1]	CRAN	(R 4.3.3)
lifecycle		1.0.4	2023-11-07			(R 4.3.1)
listenv		0.9.1	2024-01-29	[1]		
lmom		3.2		[1]		
lubridate	*	1.9.4	2024-12-08			
magrittr	*	2.0.3		[1]		
MASS		7.3-58.4	2023-03-07	[2]		
Matrix		1.6-5		[1]		
memoise		2.0.1		[1]		(R 4.3.0)
mime		0.12	2021-09-28			(R 4.3.0)
miniUI		0.12		[1]		(R 4.3.0)
minpack.lm		1.2-4	2018-05-18	[1]		(R 4.3.0)
minpack.im modeldata	ı.					
	*	1.4.0	2024-06-19	[1]		(R 4.3.3)
ModelMetrics		1.2.2.2	2020-03-17	[1]		(R 4.3.0)
moments	*	0.14.1	2022-05-02	[1]		(R 4.3.3)
munsell		0.5.1		[1]		
mvtnorm		1.3-3		[1]		
nlme		3.1-166	2024-08-14	[1]	CRAN	(R 4.3.3)

nnet		7.3-20	2025-01-01	[1]	CRAN	(R 4.3.3)
officer	*	0.6.7	2024-10-09	[1]	CRAN	(R 4.3.3)
openssl		2.3.1	2025-01-09	[1]	CRAN	(R 4.3.3)
pander		0.6.6	2025-03-01	[1]	CRAN	(R 4.3.3)
parallelly		1.41.0	2024-12-18	[1]	CRAN	(R 4.3.3)
parsnip	*	1.2.1	2024-03-22	[1]	CRAN	(R 4.3.1)
pillar		1.10.1	2025-01-07	[1]	CRAN	(R 4.3.3)
pkgbuild		1.4.6	2025-01-16	[1]	CRAN	(R 4.3.3)
pkgconfig		2.0.3	2019-09-22	[1]	CRAN	(R 4.3.0)
pkgload		1.4.0	2024-06-28	[1]	CRAN	(R 4.3.3)
plyr		1.8.9	2023-10-02	[1]	CRAN	(R 4.3.1)
pROC		1.18.5	2023-11-01	[1]	CRAN	(R 4.3.1)
processx		3.8.5	2025-01-08	[1]	CRAN	(R 4.3.3)
prodlim		2024.06.25		[1]	CRAN	(R 4.3.3)
profvis		0.4.0	2024-09-20	[1]	CRAN	(R 4.3.3)
promises		1.3.2	2024-11-28	[1]	CRAN	(R 4.3.3)
proxy		0.4-27	2022-06-09	[1]	CRAN	(R 4.3.0)
ps		1.8.1	2024-10-28	[1]	CRAN	(R 4.3.3)
purrr	*	1.0.2	2023-08-10	[1]	CRAN	(R 4.3.0)
R6	Ċ	2.5.1	2021-08-19	[1]	CRAN	(R 4.3.0)
ragg		1.3.3	2024-09-11	[1]	CRAN	(R 4.3.3)
rappdirs		0.3.3	2021-01-31	[1]	CRAN	(R 4.3.0)
Rcpp		1.0.14	2025-01-12	[1]	CRAN	(R 4.3.3)
readr	*	2.1.5	2024-01-10	[1]	CRAN	(R 4.3.1)
readxl	*	1.4.3	2023-07-06	[1]	CRAN	(R 4.3.1)
recipes	*	1.1.0	2024-07-04	[1]	CRAN	(R 4.3.3)
remotes	т	2.5.0	2024-07-04	[1]	CRAN	(R 4.3.3)
reshape2	*	1.4.4	2020-04-09	[1]	CRAN	(R 4.3.1)
rlang	-1-	1.1.4	2024-06-04	[1]	CRAN	(R 4.3.3)
rmarkdown	*	2.29	2024 00 04	[1]	CRAN	(R 4.3.3)
robustbase	-1-	0.99-4-1	2024-11-04	[1]	CRAN	(R 4.3.3)
rootSolve		1.8.2.4	2023-09-21	[1]	CRAN	(R 4.3.3)
rpart		4.1.24	2025-01-07	[1]	CRAN	(R 4.3.3)
	*	1.2.1	2024-03-25	[1]	CRAN	(R 4.3.1)
rsample rstudioapi	т	0.17.1	2024 03 23	[1]	CRAN	(R 4.3.1)
Rttf2pt1		1.3.12	2024 10 22 2023-01-22	[1]	CRAN	(R 4.3.3)
rvest		1.0.4	2024-02-12	[1]	CRAN	(R 4.3.1)
scales	*	1.3.0	2024 02 12		CRAN	(R 4.3.1)
sessioninfo	-1-	1.2.2	2023 11 26	[1]		(R 4.3.1)
sf		1.0-19	2024-11-05	[1]		(R 4.3.3)
shiny		1.10.0	2024 11 03	[1]		(R 4.3.3)
snakecase		0.11.1	2023-08-27	[1]		(R 4.3.0)
stringi		1.8.4	2023 06 27	[1]		(R 4.3.0)
stringr	*	1.5.1	2023-11-14	[1]	CRAN	(R 4.3.1)
survival	-1-	3.8-3	2024-12-17	[1]	CRAN	(R 4.3.1)
survival		2.1.3	2023-12-08	[1]	CRAN	(R 4.3.1)
systemfonts		1.1.0	2023-12-08	[1]	CRAN	(R 4.3.1)
terra		1.8-29	2024-03-13	[1]	CRAN	(R 4.3.3)
textshaping		0.4.1	2023-02-20	[1]	CRAN	(R 4.3.3)
tibble	*	3.2.1	2024-12-00	[1]	CRAN	(R 4.3.3)
tidymodels	*	1.2.0	2023-03-20	[1]	CRAN	(R 4.3.0)
tidyr	*	1.3.1	2024-03-25	[1]	CRAN	(R 4.3.1) $(R 4.3.1)$
tidyselect	т	1.2.1	2024-01-24	[1]	CRAN	(R 4.3.1)
tidyverse	*	2.0.0	2024-03-11	[1]	CRAN	(R 4.3.1)
ciayverse	*	2.0.0	2023-02 - 22	ГΤ]	OUVIN	(IL 4.3.U)

timechange		0.3.0	2024-01-18	[1]	CRAN	(R 4.3.1)
timeDate		4041.110	2024-09-22	[1]	CRAN	(R 4.3.3)
tinytex	*	0.54	2024-11-01	[1]	CRAN	(R 4.3.3)
truncnorm	*	1.0-9	2023-03-20	[1]	CRAN	(R 4.3.3)
tune	*	1.2.1	2024-04-18	[1]	CRAN	(R 4.3.1)
tzdb		0.4.0	2023-05-12	[1]	CRAN	(R 4.3.0)
units		0.8-5	2023-11-28	[1]	CRAN	(R 4.3.1)
urlchecker		1.0.1	2021-11-30	[1]	CRAN	(R 4.3.0)
useful	*	1.2.6.1	2023-10-24	[1]	CRAN	(R 4.3.1)
usethis		3.1.0	2024-11-26	[1]	CRAN	(R 4.3.3)
uuid		1.2-1	2024-07-29	[1]	CRAN	(R 4.3.3)
vctrs		0.6.5	2023-12-01	[1]	CRAN	(R 4.3.1)
viridisLite		0.4.2	2023-05-02	[1]	CRAN	(R 4.3.0)
webshot	*	0.5.5	2023-06-26	[1]	CRAN	(R 4.3.0)
webshot2	*	0.1.1	2023-08-11	[1]	CRAN	(R 4.3.0)
websocket		1.4.2	2024-07-22	[1]	CRAN	(R 4.3.3)
withr		3.0.2	2024-10-28	[1]	CRAN	(R 4.3.3)
workflows	*	1.1.4	2024-02-19	[1]	CRAN	(R 4.3.1)
workflowsets	*	1.1.0	2024-03-21	[1]	CRAN	(R 4.3.1)
xfun		0.50	2025-01-07	[1]	CRAN	(R 4.3.3)
xml2		1.3.6	2023-12-04	[1]	CRAN	(R 4.3.1)
xtable		1.8-4	2019-04-21	[1]	CRAN	(R 4.3.0)
yaml		2.3.10	2024-07-26	[1]	CRAN	(R 4.3.3)
yardstick	*	1.3.1	2024-03-21	[1]	CRAN	(R 4.3.1)
zip		2.3.1	2024-01-27	[1]	CRAN	(R 4.3.1)

- [1] /Library/Frameworks/R.framework/Versions/4.1-arm64/Resources/library [2] /Library/Frameworks/R.framework/Versions/4.3-arm64/Resources/library

[#]Sys.getenv()

^{#.}libPaths()