



Quantifying Portfolio Credit Risk

– CCruncher Technical Document –

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Version 2.5.1

Abstract

This document provides a comprehensible view of portfolio credit risk modeling using the copula approach (t-Student case). It emphasizes the practical aspects, including short introductions of methods, examples, and a discussion of the implementation details. It takes into account the sensitivity of the risk with respect to the estimated parameters and provides a practical method to incorporate this uncertainty in the portfolio credit risk measure. Finally, a method for optimizing the portfolio credit risk is provided.

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This work is licensed to the public under Creative Commons Attribution-ShareAlike 3.0 Unported License. To view a copy of this license, visit <http://creativecommons.org/licenses/by-sa/3.0/>. We have invested much time and effort in creating CCruncher; please cite it when using it for credit risk analysis.

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Glossary

n	Number of obligors in the portfolio.
s	Number of sectors/factors.
r	Number of ratings.
n_i	Number of obligors in the i -th sector/factor. It is a component of vector \vec{n} that satisfies $\sum_{i=1}^s n_i = n$.
n_{ij}^t	Number of obligors in the i -th sector/factor having the j -th rating at the beginning of the t -th period.
k_{ij}^t	Number of defaults in the i -th sector/factor having the j -th rating at the beginning of the t -th period.
T_i	Default time (aka time-until-default, aka time-to-default, aka survival time) of the i -th obligor.
F	Multivariate distribution of default times or time-until-default. It has n components with marginal distributions F_i .
F_i	Default time distribution of the i -th obligor. Sometimes refers to default time distribution of the i -th rating because obligor's F_i depends only on his rating.
p_i	Probability that i -th obligor/rating defaults within the next 1-year, i.e., $F_i(1)$.
\widehat{R}	Obligors' correlation block matrix of dimension $n \times n$.
R	Factor correlation matrix of dimension $s \times s$.
\vec{w}	Factor loadings, vector of dimension s .
$B_s(\vec{n}, \vec{d}, M)$	Covariance $s \times s$ block matrix in which \vec{n} are the number of individuals per block, \vec{d} are the diagonal values, and M are the block values. See Section A.4.
$t_d(v, \vec{\mu}, \Sigma)$	Multivariate t-Student distribution of dimension d with v degrees of freedom, mean vector $\vec{\mu}$, and dispersion or scatter matrix Σ .
$C_{v, \widehat{R}}^t$	t-Student copula with v degrees of freedom and correlation matrix \widehat{R} .
$C_{v, \vec{w}, R}^t$	Multi-factor t-Student copula with v degrees of freedom, factor loadings \vec{w} , and factor correlation matrix R . It is merely a t-Student copula.

Notes

1. Factors = Sectors. CCruncher considers that dependence between obligors is determined by obligor's industrial sectors causing that sectors and factors be equivalents. This text uses sectors and factors indistinctly to refer to factors.
2. Time unit. We use the 1-year time unit without loss of generality in order to simplify the exposition.

1 — Introduction

CCruncher is a project for quantifying portfolio credit risk. It was developed and is maintained by Gerard Torrent, an enthusiast who lives in Barcelona (Catalonia). The latest version of the project can be downloaded from the CCruncher's site, <http://www.ccruncher.net>.

The CCruncher framework comprises this document and two pieces of software. These elements are released under open-source type licenses. CCruncher is a mature and active project in which users may participate by sending suggestions and contributions to gtorrent@ccruncher.net. Because CCruncher is an evolving project, the version identifier should be mentioned when referring to the project name (e.g., CCruncher-2.5.1).

This document provides a comprehensible view of portfolio credit risk modeling using the copula approach (t-Student case). This document emphasizes the practical aspects, including short introductions of methods, examples, and a discussion of the implementation details.

Portfolio credit risk tries to respond to questions such as: “What is the expected portfolio loss for the next year?” and “How much could I lose in a really bad year?” There are several approaches to answering these questions. All of them attempt to ascertain the portfolio loss distribution and measure the credit risk using a set of selected statistics (e.g., mean, quantile), but they differ in how to obtain this distribution. The following list briefly present the most used models [7, 4, 12][3, chap. 2.4].

Actuarial models consider that the frequency of defaults (Poisson type distributions) is the proper manner in which to analyze observed defaults. Compounding default frequency with the obligors' exposure gives the portfolio loss distribution. The main representative of this approach is the CreditRisk+ model [36].

Intensity models consider that an obligor defaults when occurs the first jump of a Poisson process with random intensity given by a Cox process. Dependency between defaults is induced by assuming that each intensity is a function of a common process and an idiosyncratic process. Actuarial models may be included within this family to the extent that they have a static intensity.

Bernoulli mixture models represent the i -th obligor's default using a Bernoulli variable (with probability of default p_i). These variables are conditionally independent given a vector of default probabilities with a known density. The sum of the defaulted obligors' exposure gives the portfolio loss distribution.

Factor models represent the i -th obligor using a random variable and consider that this obligor defaults when this variable crosses a preset threshold. These variables depend on a low-dimensional random vector of common factors. Conditional on factors, factor models are Bernoulli mixture models. CreditMetricsTM [18] and KMV [6] are models of this type.

Copula models represent the i -th obligor using a random variable and consider that this obligor defaults when this variable satisfies a given condition. The dependence between these variables is explicitly modeled using a copula [10]. The sum of the defaulted obligors' exposure gives the portfolio loss distribution. In the Li copula model [24], random variables are the obligors default times and the default trigger is to cross a preset threshold. The multi-factor Gaussian case of the Li model covers the CreditMetricsTM model.

CCruncher uses the copula approach to formulate the model, then restrict the model into a factor

model in order to reduce the dimensionality, and finally interprets the model as a Bernoulli mixture model to estimate the parameters. The CCruncher model implements the t-Student multi-factor copula even for high-dimensional portfolios, and it is formulated according to the Basel II concepts (PD, EAD, LGD).

CCruncher quantifies the portfolio credit risk by sampling the portfolio loss distribution and computing the appropriate statistics (EL, VaR, ES). The portfolio losses are obtained simulating the time-until-default of obligors and simulating the EADs and LGDs of their assets. The time-until-default of obligors is a multivariate distribution in which the marginals are the probabilities of default given by obligors' ratings and in which the dependence between obligors is determined by obligors' industrial sectors. CCruncher assumes that the dependence structure is a t-Student copula, specifically a t-Student multi-factor copula in which the factors are the industrial sectors.

CCruncher requires a portfolio with a buy-and-hold policy in which each obligor has a rating assigned and belongs to a single economic sector. An example of such a credit portfolio could be an SMEs loan portfolio (buy-and-hold policy, rated obligors, and multiple sectors in which each obligor belongs to a single sector) or a portfolios of bonds. Unlike CreditMetricsTM, CCruncher does not follow a mark-to-market account, but rather follows a two-state model (default/no-default).

The advantages of CCruncher are as follows: 1) the model allows multiple factors; 2) the model allows multiple ratings; 3) obligors can have multiple assets; 4) each asset has its own EAD and LGD; 5) these EADs and LGDs can be stochastic values; 6) EADs and LGDs can vary over time; 7) CCruncher allows risk disaggregation; 8) CCruncher does not presuppose a static transition matrix; 9) factor correlations are appropriately considered; and 10) the uncertainty of parameters is considered.

The disadvantages of CCruncher are as follows: 1) it is computationally intensive; 2) the convergence rate of the Monte Carlo method is rather slow; 3) currently the obligors cannot belong to a weighted sum of factors; 4) dependence only relies on sectors; 5) correlations are constant over time; 6) currently only Gaussian and t-Student copulas are supported; and 7) mark-to-market is not supported.

The content is organized into three chapters. The first chapter details the distribution used to simulate the obligors' default times, and provides methods to estimate the copula parameters. The second chapter describes how to obtain the portfolio loss distribution and quantify the portfolio credit risk. Finally, the third chapter covers four important topics: risk sensitivity, parameters uncertainty, and portfolio optimization. In order to facilitate reading, those issues that are important to understanding the content and require a detailed discussion have been relegated to appendices.

2 — Modeling Default Times

Financial institutions have statistical models that assigns a probability of default to each obligor as well as statistical models for the exposure and recovery of each operation. These models allow us to estimate the Expected Loss of the portfolio, but are insufficient to estimate the measures related to economic capital such as VaR or ES. This is because these measures strongly depend on the dependence between obligors, and models mentioned above do not refer to this concept. This chapter develops the multivariate model using the existing PD, exposure and recovery models.

2.1 The Copula Model

The multivariate time-until-default model described in this section is identical to that described in [24, 13] [3, chap. 2.6]. We recommend perusing the appendices before proceeding. The appendices contain definitions, statements, and notation regarding subjects used in this section.

The copula approach considers that the lifetimes of obligors are the proper manner in which to explain the observed defaults. Obligor's default times (T_1, \dots, T_n) are modeled by a multivariate distribution F such that

$$F(t_1, \dots, t_n) = \Pr\{T_1 \leq t_1, \dots, T_n \leq t_n\}$$

in which n is the number of obligors in the portfolio and T_i is the time-until-default (in years) of the i -th obligor. Applying the Sklar's theorem, we obtain

$$F(t_1, \dots, t_n) = C(F_1(t_1), \dots, F_n(t_n))$$

in which marginal $F_i(t) = \Pr\{T_i \leq t\}$ is the time-until-default distribution of the i -th obligor and C is a copula. CCruncher considers that each obligor has its own time-until-default distribution depending on its rating; therefore, all obligors with identical ratings have identical distributions (see Appendix A.1 to obtain these distributions from existing PD models). Regarding the copula, CCruncher considers that C is a t-Student copula for the following practical reasons:

- it is determined by the correlation matrix (elliptical property).
- it is symmetrical (elliptical property), allowing application of the antithetic reduction variance technique in the Monte Carlo simulation.
- it allows a wide range of dependence structures by parameter ν .
- it exhibits tail dependence when $\nu \neq \infty$.
- it can be expressed as a t-Student multi-factor copula.
- an efficient simulation algorithm exists.
- the Gaussian case ($\nu = \infty$) leads to the CreditMetrics™ model.

The following proposition reformulates F in terms of the multivariate t-Student distribution. This result, combined with the Gaussian multi-factor distribution, provides an efficient simulation algorithm and allows estimating the parameters.

Proposition 2.1 — Multivariate default times distribution. Let $T = (T_1, \dots, T_n)$, the time-until-default of obligors with marginals F_1, \dots, F_n and a t-Student copula $C_{\nu, \widehat{R}}^t$. Then its distribution F is

$$F(t_1, \dots, t_n) = H(t_\nu^{-1}(F_1(t_1)), \dots, t_\nu^{-1}(F_n(t_n)))$$

in which H is the cdf of the multivariate t-Student distribution $t_n(\nu, \vec{0}, \widehat{R})$, and t_ν^{-1} is the inverse of the univariate t-Student distribution $t_1(\nu, 0, 1)$.

Proof. We note that $X = (X_1, \dots, X_n)$ a vector having a multivariate t-Student distribution $t_n(\nu, \vec{0}, \widehat{R})$ and (U_1, \dots, U_n) to represent the components of the copula $C_{\nu, \widehat{R}}^t$,

$$\begin{aligned} F(t_1, \dots, t_n) &= \Pr\{T_1 \leq t_1, \dots, T_n \leq t_n\} = \\ &\quad \text{we use the Corollary A.3: } (T_1, \dots, T_n) = (F_1^{-1}(U_1), \dots, F_n^{-1}(U_n)) \\ &= \Pr\{F_1^{-1}(U_1) \leq t_1, \dots, F_n^{-1}(U_n) \leq t_n\} = \\ &\quad \text{we use the Corollary A.2: } (U_1, \dots, U_n) = (t_\nu(X_1), \dots, t_\nu(X_n)) \\ &= \Pr\{F_1^{-1}(t_\nu(X_1)) \leq t_1, \dots, F_n^{-1}(t_\nu(X_n)) \leq t_n\} = \\ &= \Pr\{t_\nu(X_1) \leq F_1(t_1), \dots, t_\nu(X_n) \leq F_n(t_n)\} = \\ &= \Pr\{X_1 \leq t_\nu^{-1}(F_1(t_1)), \dots, X_n \leq t_\nu^{-1}(F_n(t_n))\} = \\ &= H(t_\nu^{-1}(F_1(t_1)), \dots, t_\nu^{-1}(F_n(t_n))) \end{aligned}$$

□

Corollary 2.1 — Default times simulation (I). Let $T = (T_1, \dots, T_n)$, a random vector with marginals F_1, \dots, F_n and a t-Student copula $C_{\nu, \widehat{R}}^t$. Then we can simulate this distribution using a multivariate t-Student distribution $X \sim t_n(\nu, \vec{0}, \widehat{R})$:

$$(T_1, \dots, T_n) = (F_1^{-1}(t_\nu(X_1)), \dots, F_n^{-1}(t_\nu(X_n)))$$

in which t_ν is the cdf of the univariate t-Student distribution $t_1(\nu, 0, 1)$.

The current situation is not yet satisfactory because the dimension $n \times n$ of the correlation matrix \widehat{R} indicates that any attempt to simulate random values or estimate the parameters cannot be performed in practice when n is large. For these reason we restrict the copula model to the multi-factor formulation.

2.2 The Multi-Factor Formulation

To develop a correlation matrix that is easy to handle, we assume that a reduced number s of industrial sectors exists and that each obligor belongs to a single sector. In addition, we assume that obligors' time-until-default dependence only relies on the involved industrials sectors. With these restrictions and by sorting the obligors by sector, we see that the obligors' time-until-default

correlation is an $n \times n$ correlation block matrix $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$.

$$\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma) = \begin{pmatrix} 1 & \cdots & \sigma_{11} & & \sigma_{1s} & \cdots & \sigma_{1s} \\ \vdots & \ddots & \vdots & \cdots & \vdots & & \vdots \\ \sigma_{11} & \cdots & 1 & & \sigma_{1s} & \cdots & \sigma_{1s} \\ & \vdots & & \ddots & & \vdots & \\ \sigma_{1s} & \cdots & \sigma_{1s} & & 1 & \cdots & \sigma_{ss} \\ \vdots & & \vdots & \cdots & \vdots & \ddots & \vdots \\ \sigma_{1s} & \cdots & \sigma_{1s} & & \sigma_{ss} & \cdots & 1 \end{pmatrix} \quad \Sigma = \begin{pmatrix} \sigma_{11} & \cdots & \sigma_{1s} \\ \vdots & \ddots & \vdots \\ \sigma_{1s} & \cdots & \sigma_{ss} \end{pmatrix}$$

$$\vec{n} = (n_1, \dots, n_s)$$

$$n = \sum_{i=1}^s n_i$$

Consequently, the multivariate time-until-default distribution F can be expressed by its marginals F_i and a t-Student copula $C_{\vec{v}, \widehat{R}}^t$ in which $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$.

As we have observed, Proposition 2.1 establishes the link between the t-Student copula and the multivariate t-Student distribution. Now, we are interested in formulating this multivariate t-Student distribution with correlation matrix $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$ as a t-Student multi-factor distribution.

Proposition 2.2 — t-Student distr. as a multi-factor distribution. Let a random vector $X = (X_1, \dots, X_n)$ be with a multivariate t-Student distribution $t_n(\vec{v}, \vec{0}, \widehat{R})$ in which $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$ is a correlation block matrix and n_i sizes are high enough. Then X can be written as a t-Student multi-factor distribution as follows:

$$X_i^j = \sqrt{\frac{\vec{v}}{V}} \cdot \left(w_i \cdot Z_i + \sqrt{1 - w_i^2} \cdot \varepsilon_i^j \right), \text{ where } \begin{cases} i = 1, \dots, s & s \text{ is the number of factors} \\ j = 1, \dots, n_i & n_i \text{ is } i\text{-th factor size} \\ V \sim \chi_{\vec{v}}^2 & 2 \leq \vec{v} \text{ degrees of freedom} \\ Z \sim N(\vec{0}, R) & R \text{ a } s \times s \text{ correlation matrix} \\ w_i \in (0, 1) & \text{the factor loadings} \\ \varepsilon_i^j \sim N(0, 1) \text{ iid} & \forall i, j \\ V, Z, \varepsilon_i^j \text{ indep.} & \forall i, j \end{cases}$$

The relationship between the multi-factor coefficients \vec{w} , R and the multivariate t-Student correlation matrix $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$ is

$$R_{ij} = \frac{\sigma_{ij}}{\sqrt{\sigma_{ii} \cdot \sigma_{jj}}} \quad \forall i, j$$

$$w_i = \sqrt{\sigma_{ii}} \quad i = 1, \dots, s$$

We call t-Student multi-factor copula to the copula of the t-Student multi-factor distribution. As we have observed, the t-Student multi-factor copula ($C_{\vec{v}, \vec{w}, R}^t$) is merely a t-Student copula ($C_{\vec{v}, \widehat{R}}^t$). The inverse is not true, there exist t-Student copulas with a correlation block matrix that cannot be expressed as a t-Student multi-factor copula (see Appendix A.5 and Proposition A.6).

Note that we expressed the multi-factor distribution in terms of the factor loadings and the factor correlation matrix instead of the more concise form based on the covariance matrix. We do this because the factor loadings have a treatment distinct from the factor correlation in the estimation procedure.

The t-Student multi-factor distribution combined with the Corollary 2.1 provides an efficient algorithm to simulate the obligors' default times disclosed in the next corollary. Later we will

use this algorithm to simulate the portfolio losses.

Corollary 2.2 — Default times simulation (II). Let $T = (T_1, \dots, T_n)$ be a random vector with marginals F_1, \dots, F_r and a t-Student multi-factor copula with ν degrees of freedom, factor loadings \vec{w} , and factor correlation matrix R . Then we can simulate this distribution:

1. Simulate the factors, $\vec{z} \sim N(\vec{0}, R)$.
2. Simulate the degrees of freedom, $\nu \sim \chi_\nu^2$.
3. For each obligor, repeat the following steps:
 - (a) Simulate $\varepsilon \sim N(0, 1)$.
 - (b) Simulate the multivariate t-Student component using the obligor's sector:

$$x = \sqrt{\frac{\nu}{\nu - 1}} \cdot \left(w_i \cdot z_i + \sqrt{1 - w_i^2} \cdot \varepsilon \right)$$
 - (c) Obtain the obligor's default time using the obligor's rating:

$$t = F_j^{-1}(t_\nu(x))$$

We want to estimate the copula parameters of the multivariate default times distribution. These parameters are the degrees of freedom ν , the factor loadings \vec{w} , and the factor correlation matrix R . The historical data available for parameters calibration are panel data type. Calibration with such data leads to the interpretation of the current model as a Bernoulli mixture model.

2.3 The Bernoulli Mixture Model

The available historical data lead us to interpret obligors' defaults as conditionally independent Bernoulli random variables. In this section we reinterpret the multi-factor copula model as a Bernoulli mixture model in order to estimate the copula parameters.

Definition 2.1 — Annual number of defaults. We define the annual number of defaults by sector and rating such as a discrete multivariate random variable $K = (K_{11}, \dots, K_{sr})$ in which

$$K_{ij} = \left\{ \begin{array}{l} \text{number of defaults in the } i\text{-th sector having the} \\ \text{j-th rating at the beginning of the 1-year period} \end{array} \right\}$$

taking in account that

$$N_{ij} = \left\{ \begin{array}{l} \text{number of obligors in the } i\text{-th sector having the} \\ \text{j-th rating at the beginning of the 1-year period} \end{array} \right\}$$

Note that $K_{ij} \leq N_{ij}$. We use the superscript to indicate the observation at time t (e.g., K_{ij}^t).

Example 2.1 — Historical data. Suppose that our portfolio has two sectors, S1 and S2, and four ratings, A, B, C, and D, in which D is the *defaulted* status. We collected the number of defaults by sector and rating from the past 20 years. We organized the collected data as follows:

Year	N (# obligors)						K (# defaults)					
	S1			S2			S1			S2		
	A	B	C	A	B	C	A	B	C	A	B	C
1	529	932	293	685	1249	206	3	12	15	3	25	10
2	522	977	278	638	1315	194	1	18	13	7	28	8
⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮	⋮
20	492	962	310	602	1227	183	6	28	21	4	31	16

Below, we provide an algorithm to simulate the annual number of defaults. The algorithm is a direct application of the multivariate default times distribution using the t-Student multi-factor distribution (see Proposition 2.1 and Proposition 2.2). We will use it to generate samples with which to test the parameters estimation procedures.

Algorithm 2.1 — Annual number of defaults simulation. We have a credit portfolio with n obligors. We identified s distinct sectors and categorized the individual risk profiles in r ratings. The number of obligors in each sector-rating bucket is n_{ij} . Obligor default times are modeled according a multivariate default times distribution with annual probabilities of default p_1, \dots, p_r and a t-Student multi-factor copula $C_{v, \vec{w}, R}^t$, where R is the $s \times s$ factor correlation matrix. Then we can simulate a sample of size M of the annual number of defaults by sector-rating doing the following:

- Compute the Cholesky decomposition of R
 $R = L \cdot L^\top$
- For $m = 1, \dots, M$ repeat the following steps:
 1. Initialize the simulated annual number of defaults
 $k_{ij}^m = 0 \quad i = 1, \dots, s \quad j = 1, \dots, r$
 2. Simulate the correlated factors
 $\vec{z} = L \cdot \vec{y}$ where $\vec{y} \sim N(0, \text{Id})$
 3. Simulate the chi-square factor
 $v \sim \chi_v^2$
 4. For each sector $i = 1, \dots, s$ repeat the following steps:
 - For each rating $j = 1, \dots, r$ repeat the following steps:
 - * For each one of the n_{ij} obligors repeat the following steps:
 - Simulate the t-Student component
 $x = \sqrt{\frac{v}{v-2}} \cdot \left(w_i \cdot z_i + \sqrt{1 - w_i^2} \cdot \varepsilon \right)$ where $\varepsilon \sim N(0, 1)$
 - Check for the default event
If $x < t_v^{-1}(p_j)$ then $k_{ij}^m = k_{ij}^m + 1$
 5. Print the m-th annual number of defaults
 $k_{11}^m, \dots, k_{sr}^m$

Before proceeding to estimate the parameters, we derive the density of the number of observed defaults from the multivariate default times distribution. More information is available in [30, 16, 9].

Proposition 2.3 — Density of the annual number of defaults (I). The density of the annual number of defaults, assuming that multivariate default times have marginals F_1, \dots, F_r

and a t-Student copula $C_{\mathbf{v}, \widehat{R}}^t$ with $\widehat{R} = B_s(\vec{n}, \vec{1}, \Sigma)$, is

$$\begin{aligned} K(k_{11}, \dots, k_{sr}; \mathbf{v}, \widehat{R}) &= \Pr\{K_{11} = k_{11}, \dots, K_{sr} = k_{sr}; \mathbf{v}, \widehat{R}\} = \\ &= \left(\prod_{i=1}^s \prod_{j=1}^r \binom{n_{ij}}{k_{ij}} \right) \cdot \underbrace{\int_{-\infty}^{t_{\mathbf{v}}^{-1}(p_1)}}}_{k_{11} \text{ times}} \underbrace{\int_{t_{\mathbf{v}}^{-1}(p_1)}^{\infty}}_{n_{11}-k_{11} \text{ times}} \cdots \underbrace{\int_{-\infty}^{t_{\mathbf{v}}^{-1}(p_r)}}}_{k_{sr} \text{ times}} \underbrace{\int_{t_{\mathbf{v}}^{-1}(p_r)}^{\infty}}_{n_{sr}-k_{sr} \text{ times}} f(x_1, \dots, x_n) dx_1 \dots dx_n \end{aligned}$$

in which s is the number of sectors, r is the number of ratings, n_{ij} and k_{ij} are respectively the number of obligors and defaults in the i -th sector with the j -th rating, $p_j = F_j(1)$ is the probability of default within 1 year of the j -th rating, $f(\vec{x})$ is the pdf of $t_n(\mathbf{v}, \vec{0}, \widehat{R})$, and $t_{\mathbf{v}}^{-1}$ is the inverse of the cdf of the univariate t-Student. The combinatorial coefficients consider all combinations of exchangeable obligors.

We are not aware of an analytical expression for the density of the annual number of defaults. Furthermore, the numerical solution of the integral is difficult when the number of obligors is larger than just a few. Therefore, we attempt to solve the problem recurring in the multi-factor formulation.

Proposition 2.4 — Conditional density of the annual number of defaults. The conditional density of the annual number of defaults, assuming that multivariate default times have marginals F_1, \dots, F_r and a t-Student multi-factor copula $C_{\mathbf{v}, \vec{w}, R}^t$, is

$$\begin{aligned} K(k_{11}, \dots, k_{sr} | \mathbf{v}, \vec{z}; \mathbf{v}, \vec{w}, R) &= \Pr\{K_{11} = k_{11}, \dots, K_{sr} = k_{sr} | \mathbf{v}, \vec{z}; \mathbf{v}, \vec{w}, R\} = \\ &= \prod_{i=1}^s \prod_{j=1}^r \binom{n_{ij}}{k_{ij}} \cdot H_{ij}(\mathbf{v}, z_i)^{k_{ij}} \cdot (1 - H_{ij}(\mathbf{v}, z_i))^{n_{ij}-k_{ij}} \end{aligned}$$

in which

$$H_{ij}(\mathbf{v}, z_i) = \Phi \left(\frac{\sqrt{\frac{\mathbf{v}}{\mathbf{v}}} \cdot t_{\mathbf{v}}^{-1}(p_j) - w_i \cdot z_i}{\sqrt{1 - w_i^2}} \right)$$

and s is the number of sectors, r is the number of ratings, n_{ij} and k_{ij} are respectively the number of obligors and defaults in the i -th sector with the j -th rating, $p_j = F_j(1)$ is the probability to default within 1 year of the j -th rating, Φ is the cdf of the univariate $N(0, 1)$, and $t_{\mathbf{v}}^{-1}$ is the inverse of the cdf of the univariate t-Student. The combinatorial coefficients consider all combinations of exchangeable obligors.

Proof. We use the pdf of the default times distribution (see Proposition 2.1) to obtain the probability that an obligor k belonging to i -th sector with j -th rating defaults the first year conditioned to z_i and \mathbf{v} :

$$H_{ij}(\mathbf{v}, z_i) = \Pr\{T_k \leq 1 | \mathbf{v}, z_i\} = \Pr\{X_k \leq t_{\mathbf{v}}^{-1}(F_j(1)) | \mathbf{v}, z_i\} = \Pr\{X_k \leq t_{\mathbf{v}}^{-1}(p_j) | \mathbf{v}, z_i\}$$

in which T_k is the k -th component of the multivariate default times and X_k is the k -th component of the multivariate t-Student. Now we express X in the multi-factor form:

$$H_{ij}(\mathbf{v}, z_i) = \Pr \left\{ \sqrt{\frac{\mathbf{v}}{\mathbf{v}}} \cdot \left(w_i \cdot z_i + \sqrt{1 - w_i^2} \cdot \varepsilon_i^j \right) \leq t_{\mathbf{v}}^{-1}(p_j) | \mathbf{v}, z_i \right\}$$

Because \mathbf{v} and z_i are fixed and independent from $\varepsilon_i^j \sim N(0, 1)$, we obtain

$$H_{ij}(\mathbf{v}, z_i) = \Phi \left(\frac{\sqrt{\frac{\mathbf{v}}{v}} \cdot t_v^{-1}(p_j) - w_i \cdot z_i}{\sqrt{1 - w_i^2}} \right)$$

H_{ij} is the probability that an obligor defaults in 1 year. The probability of observing k_{ij} failures in a set of n_{ij} independent obligors follows a binomial distribution:

$$\text{Bin}(k_{ij}; n_{ij}, H_{ij}) = \binom{n_{ij}}{k_{ij}} \cdot H_{ij}^{k_{ij}} \cdot (1 - H_{ij})^{n_{ij} - k_{ij}}$$

The product for all sectors and ratings gives the stated formula. We can do the product because \mathbf{v} and z_i are fixed. □

Corollary 2.3 — Density of the annual number of defaults (II). We can express the density of the annual number of defaults using the conditional density:

$$\begin{aligned} K(k_{11}, \dots, k_{sr}; \mathbf{v}, \vec{w}, R) &= \Pr\{K_{11} = k_{11}, \dots, K_{sr} = k_{sr}; \mathbf{v}, \vec{w}, R\} = \\ &= \int_2^\infty \int_{\mathbb{R}^s} \psi(\mathbf{v}) \cdot \phi(\vec{z}) \cdot K(k_{11}, \dots, k_{sr} | \mathbf{v}, \vec{z}; \mathbf{v}, \vec{w}, R) dz_1 \dots dz_s d\mathbf{v} \end{aligned}$$

in which ψ is the pdf of χ_v^2 and ϕ is the pdf of $N(\vec{0}, R)$.

This expression has a dimension significantly less than before. Despite this, this expression does not have a known analytical form and cannot be approximated numerically with accuracy when the number of factors is high (e.g., 10). The following sections provide effective methods for parameter estimation.

2.4 Frequentist Inference

The cohort approach provides a parametric estimator [5, sec. 3.1] based on estimates p_i (probability that an obligor defaults) and p_{ij} (probability that two obligors default simultaneously). We have adapted to the general case (t-Student) assuming the \mathbf{v} parameter is fixed and known.

Definition 2.2 — Implied covariance estimator. We assume that all obligors in a sector are exchangeable (identical rating per sector) and that we know the degrees of freedom of the t-Student multi-factor copula. We define the estimator of the factors' covariance, $\sigma_{ij} = w_i \cdot w_j \cdot R_{ij}$, as that value $\widetilde{\sigma}_{ij}$ that satisfies

$$p_{ij} = \int_{-\infty}^{t_v^{-1}(p_i)} \int_{-\infty}^{t_v^{-1}(p_j)} f(x, y; \mathbf{v}, \widetilde{\sigma}_{ij}) dx dy \quad \text{where} \quad \begin{cases} p_i = \frac{\sum k_i^t}{\sum n_i^t} \\ p_j = \frac{\sum k_j^t}{\sum n_j^t} \\ p_{ii} = \frac{\sum k_i^t \cdot (k_i^t - 1)}{\sum n_i^t \cdot (n_i^t - 1)} & i = j \\ p_{ij} = \frac{\sum k_i^t \cdot k_j^t}{\sum n_i^t \cdot n_j^t} & i \neq j \end{cases}$$

in which n_i^t and k_i^t are respectively the number of obligors and defaults in the i -th sector and t -th year, $f(x, y; \mathbf{v}, \sigma_{ij})$ is the pdf of the bivariate t-Student, and t_v^{-1} is the inverse of the

cdf of the univariate t-Student. Listing 2.1 implements this estimator in the R programming language.

Listing 2.1: Implied covariance (R script)

```
#install.packages("mvtnorm")
library(mvtnorm)

pri <- function(i, K, N)
{
  if (sum(dim(K) != dim(N))) stop("K and N dim differs")
  if (i <= 0 | ncol(K) < i) stop("i out-of-range")
  ret = sum(K[,i])/sum(N[,i])
  return(ret)
}

pri_j <- function(i, j, K, N)
{
  if (sum(dim(K) != dim(N))) stop("K and N dim differs")
  if (i <= 0 | ncol(K) < i) stop("i out-of-range")
  if (j <= 0 | ncol(K) < j) stop("j out-of-range")
  if (i != j) {
    ret = sum(K[,i]*K[,j])/sum(N[,i]*N[,j])
  }
  else {
    ret = sum(K[,i]*(K[,i]-1))/sum(N[,i]*(N[,i]-1))
  }
  return(ret)
}

f <- function(r, nu, p1, p2, p12)
{
  q1 = qt(p1, df=nu)
  q2 = qt(p2, df=nu)
  R = matrix(nrow=2, ncol=2, 1)
  R[1,2] = R[2,1] = r
  val = pmvt(lower=c(-Inf,-Inf), upper=c(q1,q2), corr=R, df=nu)
  return(p12-val)
}

implied_covariance <- function(nu, K, N)
{
  if (sum(dim(K) != dim(N))) stop("K and N dim differs")
  n = ncol(K)
  ret = matrix(nrow=n, ncol=n, 0)
  for(i in 1:n) {
    for(j in i:n) {
      p1 = pri(i, K, N)
      p2 = pri(j, K, N)
      p12 = pri_j(i, j, K, N)
      ret[i,j] = uniroot(f, nu, p1, p2, p12, interval=c(-0.9,0.9))$root
      ret[j,i] = ret[i,j]
    }
  }
  return(ret)
}
```


Example 2.2 — Implied covariance estimator. Consider the following scenario:

- Two sectors {S1, S2}
- Number of obligors by sector $\vec{n} = (1250, 750)$
- Default probability by sector $p_{S1} = 2.5\%$ and $p_{S2} = 7\%$
- Degrees of freedom $\nu = 15$
- Factor loadings $\vec{w} = (0.3, 0.25)$
- Factors correlation $R_{12} = 0.1$

We simulate the number of defaults of this scenario 1000 times using Algorithm 2.1. Then, we compute the implied covariance estimator and display the results in the table below.

Exact	Estimated
$w_1^2 = 0.09$	$w_1^2 \approx \widetilde{\sigma}_{11} = 0.0882$
$w_2^2 = 0.0625$	$w_2^2 \approx \widetilde{\sigma}_{22} = 0.0662$
$w_1 \cdot w_2 \cdot R_{12} = 0.0075$	$w_1 \cdot w_2 \cdot R_{12} \approx \widetilde{\sigma}_{12} = 0.0072$

The drawbacks of the implied covariance estimator are that: 1) it does not consider parameter uncertainty, 2) it requires knowing in advance the value of ν , and 3) it only supports one rating per sector. This last point is crucial because it implies the absence of an economic cycle. For these reasons, one should not use this estimator. The rest of this chapter is devoted to developing a generic estimation procedure that avoids these problems.

2.5 Bayesian Inference

The sample correlation coefficient of 20 Gaussian values has a high uncertainty [21, pp. 217-221]. This stylized fact reflects our situation because we are trying to estimate factor correlations among other values. Uncertainty is exacerbated because our variables are continuous and latent but we are estimating from observations of type count-data. Thus, considering the estimated parameters as fixed values such as the Maximum Likelihood estimator is a naive and risky proposition. The Bayesian approach appears more appropriate [17, 32] because it considers the uncertainty of the parameters. Later we will see how to transfer this uncertainty to the assessment of credit risk.

We define a numerically feasible likelihood function using the conditional density of the number of defaults, thus implying that the parameters to be inferred are ν, \vec{w}, R and the latent variables $\mathbf{v}^1, \vec{z}^1, \dots, \mathbf{v}^T, \vec{z}^T$.

Definition 2.3 — Likelihood function. Assume a set of observations $\{K^t, N^t\}_{t=1, \dots, T}$ in which N^t and K^t are vectors representing the number of obligors and defaulted obligors in each sector-rating bucket. We define the likelihood of these observations as

$$\begin{aligned}
 L(\nu, \vec{w}, R, \mathbf{v}^1, \vec{z}^1, \dots, \mathbf{v}^T, \vec{z}^T ; K^1, \dots, K^T) &= \\
 &= \prod_{t=1}^T \psi_{\nu}(\mathbf{v}^t) \cdot \phi_R(\vec{z}^t) \cdot K(k_{11}^t, \dots, k_{sr}^t | \mathbf{v}^t, \vec{z}^t ; \nu, \vec{w}, R) = \\
 &= \prod_{t=1}^T \psi_{\nu}(\mathbf{v}^t) \cdot \phi_R(\vec{z}^t) \prod_{i=1}^s \prod_{j=1}^r \binom{n_{ij}^t}{k_{ij}^t} \cdot H_{ij}(\mathbf{v}^t, z_i^t)^{k_{ij}^t} \cdot (1 - H_{ij}(\mathbf{v}^t, z_i^t))^{n_{ij}^t - k_{ij}^t}
 \end{aligned}$$

in which ψ_{ν} is the pdf of χ_{ν}^2 , ϕ_R is the pdf of $N(\vec{0}, R)$, and H_{ij} is the function defined in

Proposition 2.4.

We apply the Metropolis-Hastings algorithm to this conditional likelihood to infer the distribution of variables $v, \vec{w}, R, v^1, \vec{z}^1, \dots, v^T, \vec{z}^T$. Refer to Appendix A.6 for a description of this algorithm adapted to our case.

Proposition 2.5 — Parameters estimation. We can estimate the parameters of the t-Student multi-factor copula from the observed annual number of defaults using the Metropolis-Hastings algorithm detailed in Algorithm A.3, considering

- parameters $\theta \equiv \{v, \vec{w}, R, v^1, \vec{z}^1, \dots, v^T, \vec{z}^T\}$
- observations $y \equiv \{K^t, N^t\}_{t=1, \dots, T}$
- likelihood $f(y | \theta) \equiv L(v, \vec{w}, R, v^1, \vec{z}^1, \dots, v^T, \vec{z}^T ; K^1, \dots, K^T)$ see Definition 2.3
- proposal distributions $q_j(\theta_j | \theta'_j) \equiv N(\theta, \sigma_j)$ all parameters
- prior distribution $v \equiv \text{Unif}(2, 300)$
- prior distribution $w_i \equiv \text{Unif}(0, 1)$ $i = 1, \dots, s$
- prior distribution $r_{ij} \equiv \text{Unif}(-1, +1)$ $i = 1, \dots, s \quad j > i$
- prior distribution $\vec{z}^t \equiv \text{Unif}(-10, +10)$ $t = 1, \dots, T$
- prior distribution $v^t \equiv \text{Unif}(0, 1000)$ $t = 1, \dots, T$

Example 2.3 — Model parameters estimation. Let us consider the following scenario:

- t-Student copula with $v = 20$.
- Three ratings {A, B, D} in which D is the *defaulted* status and three sectors {S1, S2, S3} having the following 1-year probabilities of default, factor loadings, and factor correlation matrix:

$$p_A = 2\% \quad , \quad \vec{w} = (0.4, 0.25, 0.3) \quad , \quad R = \begin{pmatrix} 1 & 0.5 & 0.3 \\ 0.5 & 1 & 0.4 \\ 0.3 & 0.4 & 1 \end{pmatrix}$$

$$p_B = 8\%$$

- The number of obligors by sector-rating is

Obs	N (# obligors)					
	S1		S2		S3	
	A	B	A	B	A	B
1	250	750	500	500	750	250
2	250	750	500	500	750	250
⋮	⋮	⋮	⋮	⋮	⋮	⋮
1000	250	750	500	500	750	250

1000 observations case. We simulate 1000 observations of the number of defaults by sector-rating using Algorithm 2.1. Then we infer the parameters distribution from this sample using Metropolis-Hastings Algorithm A.3. The results are displayed in Figure 2.1. Note that the variance of factor loadings is lower than the factor correlations.

20 observations case. To appreciate the effect of the number of observations in the final result, we repeated the estimation procedure but considered only the first 20 observations instead of the 1000 available. The results are displayed in Figure 2.2. Note that the parameters uncertainty increases dramatically. The scenario depicted here is common; generally, we

have just a few decades of observations. Later we will see how to transfer this parameters uncertainty to the assessment of portfolio risk.

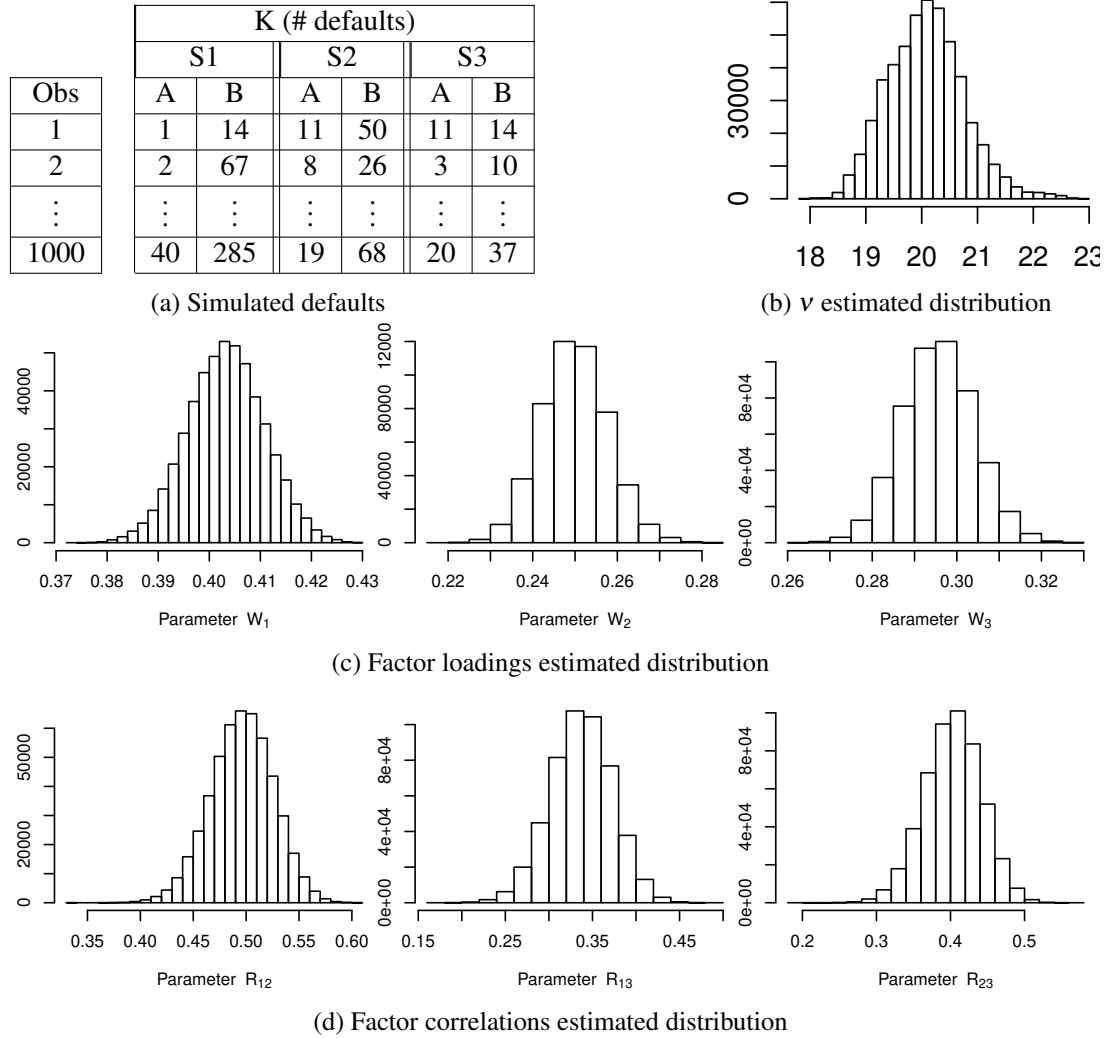


Figure 2.1: Model parameters estimation (1000 obs.)

Averaged PDs. We may be tempted to consider an averaged PD by sector when we lack data disaggregated by rating. The tests we have conducted indicate that in the Gaussian case ($\nu = \infty$), this could be an option because the estimated parameters using the averaged PDs approximately match the expected values. By contrast, when $\nu \neq \infty$, this does not work. In Example 2.3, the averaged 1-year PDs by sectors are

$$\begin{aligned}
 p_{S_1} &= \frac{250}{1000} \cdot p_A + \frac{750}{1000} \cdot p_B = 6.5\% & N_{S_1} &= 250 + 750 = 1000 \\
 p_{S_2} &= \frac{500}{1000} \cdot p_A + \frac{500}{1000} \cdot p_B = 5.0\% & N_{S_2} &= 500 + 500 = 1000 \\
 p_{S_3} &= \frac{750}{1000} \cdot p_A + \frac{250}{1000} \cdot p_B = 3.5\% & N_{S_3} &= 750 + 250 = 1000
 \end{aligned}$$

Constrained model. When we have a significant number of factors (e.g., 10) and a reduced set of observations (e.g., 20), then the uncertainty of the factor correlation matrix can be high. In these cases, one may consider a constrained model [9] that assumes that all factor correlations have the identical value or similar assumptions that reduce the number of correlations from which to estimate $s \times (s - 1) / 2$ to only a few.

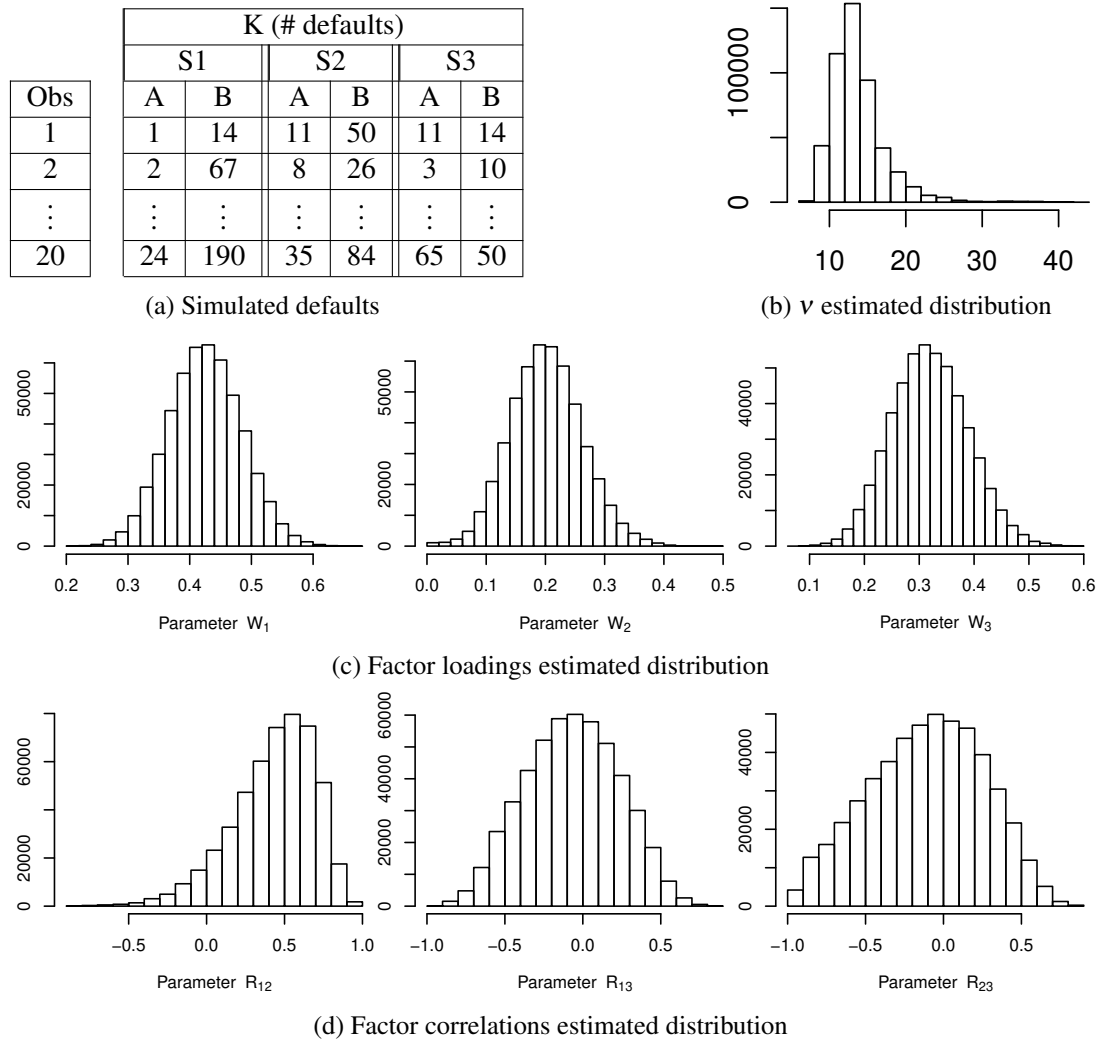


Figure 2.2: Model parameters estimation (20 obs.)

Degrees of freedom. We observed that factor correlations and degrees of freedom are interwoven. When we fix one of them with a value more distinct than generated data and we estimate the other one, the latter reacts to compensate for the modification. This effect is more pronounced when the number of observations is reduced and the amount of information available is scarce, causing the estimations of ν and R to have a high uncertainty. The Gaussian case ($\nu = \infty$) avoids this problem and reduces the number of parameters to estimate (we do not need to estimate ν and ν^t in which $t = 1, \dots, T$). The risk analyst must evaluate the appropriateness of the Gaussian model considering the amount of information available, the number of factors, the renounce to capture the tail dependence, the Deviance Information Criterion (DIC) values, etc.

3 — Modeling Portfolio Loss

3.1 Portfolio Loss Distribution

Figure 3.1 displays the hierarchy of the portfolio components. As we observed in Chapter 2, each obligor belongs to a single sector/factor and has a PD depending only on its rating. CCRuncher considers that obligors' default times distribution has marginals F_i and a t-Student multi-factor copula with ν degrees of freedom, factor loadings \vec{w} , and factor correlation matrix R .

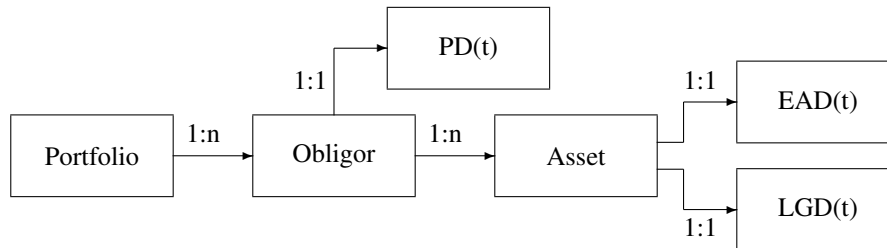


Figure 3.1: Hierarchy of the portfolio components

We define the EAD and LGD concepts used to model the credit risk emphasizing the temporal dimension. These concepts and the PD defined in Appendix A.1 are those of the Basel II accord [28], the de-facto international financial risk regulation framework.

Definition 3.1 — Exposure at Default (EAD). The Exposure at Default of the j -th asset of the i -th obligor at time t indicates the total amount that the bank is exposed if the obligor defaults:

$$EAD_i^j(t) = \left\{ \begin{array}{l} \text{amount in risk of the } j\text{-th asset of the } i\text{-th obligor} \\ \text{if this defaults at time } t, \text{ measured in currency} \end{array} \right\}$$

$EAD_i^j(t)$ is a function when these amounts are fixed and known in advance (e.g., loan, bond) or a distribution when these values are stochastic (e.g., line of credit).

Definition 3.2 — Loss Given Default (LGD). The Loss Given Default of the j -th asset of the i -th obligor at time t indicates the percentage of effective loss over the total exposure. This value depends on endorsements, guarantees, type of asset, default time, etc.

$$LGD_i^j(t) = \left\{ \begin{array}{l} \text{percentage of the } EAD_i^j(t) \text{ that is definitively} \\ \text{lost if the } i\text{-th obligor defaults at time } t \end{array} \right\}$$

$LGD_i^j(t)$ can be a function or a distribution.

Example 3.1 — 10-year Bond. Assume a 10-year, fixed-rate bond with an annual coupon of 4% issued by a company rated BB. The EAD for this asset can be obtained from its expected cash flow because it is known in advance. We suppose that the bank's internal

models indicate that in this case the LGD is a Beta(5, 2) distribution regardless of the default time. Figure 3.2 illustrates the defined concepts applied to this asset.

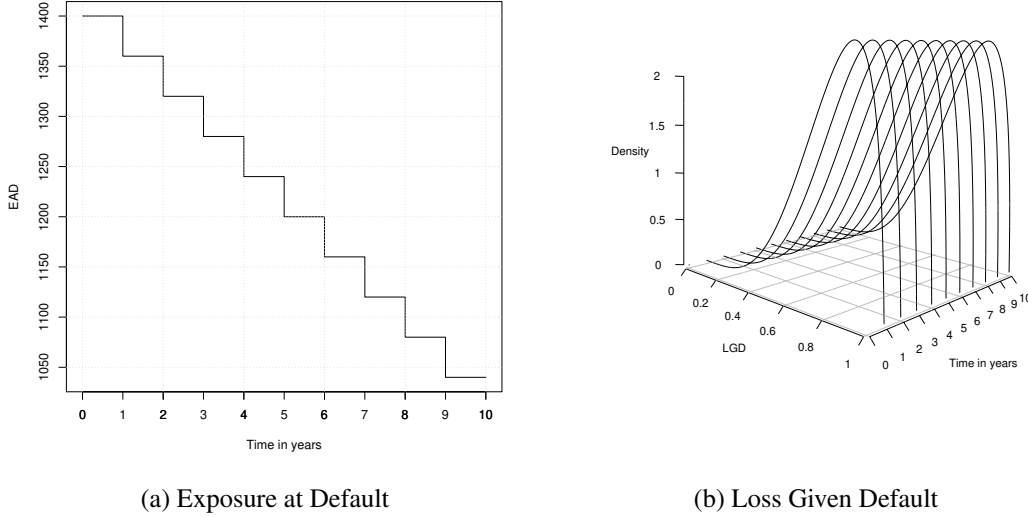


Figure 3.2: 10-year bond example

Definition 3.3 — Portfolio Loss at time T . The Portfolio Loss at time T is the accumulated losses in the time range $[0, T]$ caused by obligors that default (fail to make payments that they are obligated to make).

$$L_T = \{ \text{sum of portfolio losses in time range } [0, T] \}$$

Assume a credit portfolio in which the obligors' default times (T_1, \dots, T_n) is a multivariate distribution. In addition, we consider that each i -th obligor has m_i assets with known EAD and LGD. Thus, the portfolio loss distribution at time horizon t is

$$L_t = \sum_{i=1}^n \mathbb{1}_{[0,t]}(T_i) \cdot \left(\sum_{j=1}^{m_i} \text{EAD}_i^j(T_i) \cdot \text{LGD}_i^j(T_i) \right)$$

CCruncher assumes that the multivariate default times distribution has marginals F_i and a t-Student multi-factor copula with degrees of freedom ν , factor loadings \vec{w} , and factor correlation matrix R . CCruncher generates samples of the portfolio loss distribution simulating the obligors' default times. Portfolio loss distribution is approximated by the empirical distribution of the sample, and credit risk is measured using the appropriate statistics.

The portfolio loss distribution L_t does not have a closed form except in a few cases. One of those cases is the Large Homogeneous Pool (LHP). Vasicek [34] developed an analytical formula in the Gaussian case for the loss distribution of an infinitely large portfolio. Later, Schloegl and O'Kane [29] extended the formula to the t-Student case. The assumptions of these approximations are 1) a one-factor model, 2) an infinitely large portfolio, 3) equal PDs, and 4) equal exposure. These simplified models must be managed with care because they can cause problems [25] when the assumptions are not realized.

Algorithm 3.1 — Portfolio losses sampling. Let a portfolio be categorized in s sectors and r ratings. Assume a multivariate default times distribution with marginals F_i and a t-Student multi-factor copula $C_{v, \vec{w}, R}^t$. We can then simulate a sample of size M of the portfolio loss distribution at time T by performing the following:

- Compute the Cholesky decomposition of $R = L \cdot L^T$
- For $m = 1, \dots, M$ repeat the following steps:
 1. $\text{Loss}^m = 0$
 2. Simulate s independent $N(0, 1)$, \vec{y}
 3. Simulate correlated factors doing $\vec{z} = L \cdot \vec{y}$
 4. Simulate $v \sim \chi_v^2$
 5. For each obligor $i = 1, \dots, n$ repeat the following steps:
 - Set $si = i$ -th obligor sector index
 - Set $ri = i$ -th obligor rating index
 - Simulate $\varepsilon \sim N(0, 1)$
 - Compute $x_i = \sqrt{\frac{v}{v}} \cdot \left(w_{si} \cdot z_{si} + \sqrt{1 - w_{si}^2} \cdot \varepsilon \right)$
 - Compute $t_i = F_{ri}^{-1}(t_v(x_i))$
 - if $t_i \leq T$, then for each asset j belonging to i -th obligor
 - * Evaluate or simulate $\text{EAD}_i^j(t_i)$
 - * Evaluate or simulate $\text{LGD}_i^j(t_i)$
 - * $\text{Loss}^m = \text{Loss}^m + \text{EAD}_i^j(t_i) \cdot \text{LGD}_i^j(t_i)$
 6. print Loss^m

Example 3.2 — Large Homogeneous Pool. Assume a portfolio composed of 1000 exchangeable obligors (with identical ratings and sectors) in which every obligor owes 1€ to 1 year, the 1-year default probability is $p = 10\%$, and the factor loading is 20%. This problem is close to fulfilling the LHP assumptions. In the Gaussian case, the density is [3, chap. 2.5]:

$$f_{p,w}(x) = \frac{\sqrt{1-w^2}}{w} \cdot \exp \left(-\frac{(N^{-1}(x))^2}{2} - \frac{(N^{-1}(p) - \sqrt{1-w^2} \cdot N^{-1}(x))^2}{2 \cdot w^2} \right)$$

in which w is the factor loading, p is the probability to a 1-year horizon, and $x \in (0, 1)$ represents the percentage of loss in the portfolio.

File `CCRUNCHER/samples/test04.xml` contains the CCruncher input file of this example. Figure 3.3 displays the theoretical distribution provided by the LHP approach and the distribution approximated by CCruncher, which, as expected, are quite similar. If we reduce the number of obligors from 1000 to 100, then the LHP assumptions are broken and the distribution provided by the LHP approach fails.

3.2 Credit Risk Measurement

After N simulations (e.g., 20,000, 500,000 or more) of the portfolio loss, we measure the portfolio credit risk computing the corresponding statistics and their standard errors according to

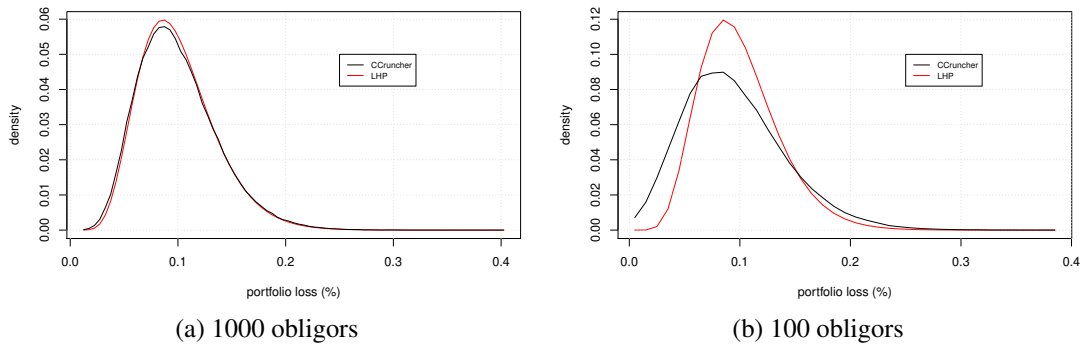
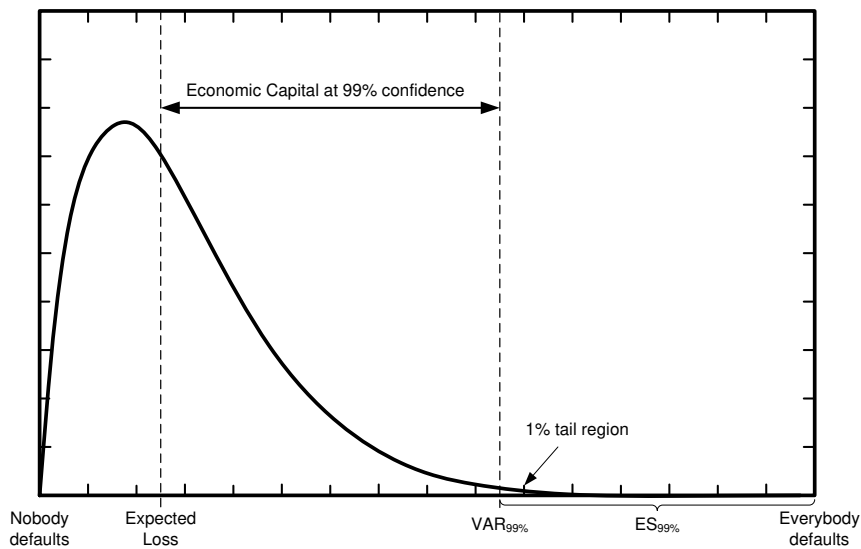


Figure 3.3: Large Homogeneous Pool example

sample size.

3.2.1 Portfolio risk

In this section we provide the definitions and formulas of the most usual credit risk measures. Figure 3.4 illustrates the concepts presented in this section. Any other user-defined risk measure can be estimated from the simulated sample, and its standard error can be obtained using resampling methods.

Figure 3.4: Portfolio Loss Distribution at time T

Expected Loss

The Expected Loss attempts to answer the question, “What is my expected loss for the next year?”

Definition 3.4 — Expected Loss (EL). The Expected Loss of a portfolio at time horizon T

is the mean of the portfolio loss distribution.

$$EL(L_T) = E(L_T)$$

Expected Loss can be approximated using the following statistic:

$$EL = \hat{\mu} \pm \Phi^{-1}\left(\frac{1-\alpha}{2}\right) \cdot \frac{\hat{\sigma}}{\sqrt{N}} \quad \text{being} \quad \begin{cases} \hat{\mu} = \frac{1}{N} \sum_{i=1}^N x_i \\ \hat{\sigma} = \sqrt{\frac{1}{N-1} \sum_{i=1}^N (x_i - \hat{\mu})^2} \end{cases}$$

in which x_i are the simulated portfolio losses, α is the error confidence level, Φ^{-1} is the $N(0, 1)$ inverse cumulative distribution function, and $\hat{\mu}$ and $\hat{\sigma}$ are the mean and standard deviation estimators.

Value at Risk

The Value at Risk [20] is the most commonly used risk measure to respond to the question, ‘‘How much could I lose in a really bad year?’’

Definition 3.5 — Value at Risk (VaR). The Value at Risk of a portfolio at time horizon T and confidence level β is the β -quantile of the portfolio loss distribution at time T . Consider

$$VaR_{\beta}(L_T) = \inf\{x \in \mathbb{R} \mid F(x) \geq \beta\}$$

in which $F(x) = \Pr\{L_T \leq x\}$ is the portfolio loss distribution at time T .

Value at Risk can be approximated using the following statistic:

$$VaR_{\beta} = \widehat{q}_{\beta} \pm \Phi^{-1}\left(\frac{1-\alpha}{2}\right) \cdot \text{stderr}(q_{\beta}) \quad \text{being} \quad \begin{cases} \frac{k}{N} \leq \beta < \frac{k+1}{N} \\ \widehat{q}_{\beta} = x_{k:N} \end{cases}$$

$$\text{stderr}(q_{\beta}) = \sqrt{C_2 - C_1^2} \quad \text{being} \quad \begin{cases} M = [N \cdot \beta + 0.5] \\ a = M - 1 \\ b = N - M \\ W_i = B(a, b, \frac{i+1}{N}) - B(a, b, \frac{i}{N}) \\ C_k = \sum_{i=1}^N W_i \cdot x_i^k \end{cases}$$

in which x_i are the simulated portfolio losses, $x_{k:N}$ is the k -th element of ascendant-sorted values, α is the error confidence level, Φ^{-1} is the $N(0, 1)$ inverse cumulative distribution function, \widehat{q}_{β} is the quantile estimator, $\text{stderr}(q_{\beta})$ is the estimation of the standard error using the Maritz-Jarrett method described in [35, chap. 3.5.3], $[x]$ is the integer part of x , and $B(a, b, x)$ is the regularized incomplete beta function. When the sample size is high, the incomplete beta function can be difficult to ascertain due to convergence problems. Also, a large sample size causes the evaluation of the incomplete beta function multiple being this computation numerically costly.

We note that the variance of the VaR is difficult to evaluate. Fortunately, it is asymptotically normal distributed [4]. This result allows to estimate the VaR value using a numerically stable procedure. This will be done in a future version of this document.

Expected Shortfall

VaR is not a coherent risk measure because it does not fulfill the sub-additive property [19]. Expected Shortfall is a coherent risk measure similar to VaR [1].

Definition 3.6 — Expected Shortfall (ES). The Expected Shortfall (aka Conditional VaR, aka Conditional Tail Expectation) of a portfolio at time horizon T and confidence level β is the average of the worst $\beta\%$ portfolio losses at time T ,

$$ES_{\beta}(L_T) = E(L_T \mid L_T \geq \text{VaR}_{\beta}(L_T))$$

Expected Shortfall can be approximated using the following statistic:

$$ES_{\beta} = \hat{\mu} \pm \Phi^{-1}\left(\frac{1-\alpha}{2}\right) \cdot \frac{\hat{\sigma}}{\sqrt{K}} \quad \text{being} \quad \begin{cases} \hat{\mu} = \frac{1}{K} \sum_{i=1}^K y_i \\ \hat{\sigma} = \sqrt{\frac{1}{K-1} \sum_{i=1}^K (y_i - \hat{\mu})^2} \end{cases}$$

in which $\{y_1, \dots, y_K\} = \{x_i \mid x_i \geq \text{VaR}_{\beta}\}_{i=1, \dots, N}$ are the simulated portfolio losses larger than VaR_{β} , α is the error confidence level, Φ^{-1} is the $N(0, 1)$ inverse cumulative distribution function, and $\hat{\mu}$ and $\hat{\sigma}$ are the mean and standard deviation estimators. This estimator assumes that the estimated VaR_{β} has the exact value. Despite this, if the sample size is high enough, then the estimated value is good enough.

Like VaR, ES is asymptotically normal distributed [4]. This result allows to estimate the value of the ES in a more rigorous way. This will be done in a future version of this document.

Example 3.3 — test05.xml. Assume a portfolio composed of 1000 obligors in which every obligor owes 1€ to 1 year. Then $EAD_i^1(t) = \mathbb{1}_{[0,1]}(t)$, with a $LGD_i^1(t) = 100\%$. There are 4 ratings {A, B, C, D}; D is the *defaulted* status, and the remaining ratings have the following 1-year default probability: $p_A = 5\%$, $p_B = 10\%$, $p_C = 20\%$. Factor loadings are $\vec{w} = (0.4, 0.35)$, and the factors correlation is $\text{Cor}(S_1, S_2) = 0.25$. The t-Student copula has $\nu = 15$ degrees of freedom, and the number of obligors by sector and rating are

S1			S2		
A	B	C	A	B	C
167	166	167	167	167	166

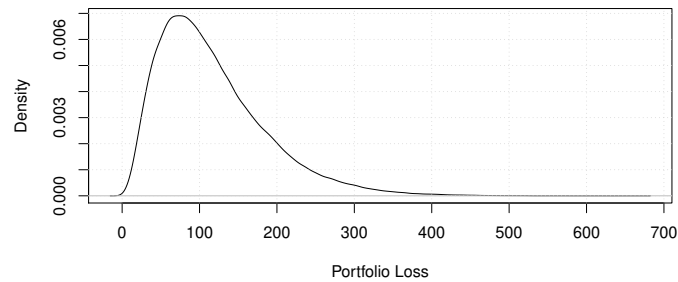
File `$$CCRUNCHER/samples/test05.xml` contains the CCRuncher input file of this example. Figure 3.5 displays the portfolio loss density and risk measures obtained using 200,000 values simulated by CCRuncher.

It is difficult to ascertain the validity of this example. Some authors [14] obtained quasi-analytical expressions for the Gaussian two-factor model with a common PD; however, those expressions are not applicable to this example (t-Student, multiple PDs). One manner in which

to check that the result is correct involves disaggregating the simulated portfolio losses by sector-rating (see Section 3.2.2) and estimating the parameters v , \vec{w} , and R with the first 1000 simulations using the procedure described in Section 2.5. If the estimated parameters match the parameters of the example, we have consistent simulation and estimation algorithms. Observe that in this case, the simulated losses can be equated to the number of defaults because each defaulted obligor is computed as a 1€ loss.

Statistic	Value	Std. Err.
EL	116.58	0.16
VaR _{95%}	252.00	1.09
VaR _{99%}	338.00	2.22
VaR _{99.9%}	442.00	5.44
ES _{95%}	304.63	0.49
ES _{99%}	383.82	0.96
ES _{99.9%}	480.98	2.71

(a) Risk measures



(b) Portfolio loss density

Figure 3.5: 2-factors example

3.2.2 Risk Disaggregation

Once we have measured the credit risk, naturally arises the question “how much of this risk corresponds to the business unit X?” CCruncher allows defining and simulating sub-portfolio losses in order to respond this question.

CCruncher simulates the entire portfolio loss at time T . It also allows (in the same execution) to obtain the sub-portfolios losses. In CCruncher, a sub-portfolio is called a *segment*. The disjoint union of segments covering the entire portfolio is called a *segmentation*, i.e., a segmentation is composed by non-overlapped segments that constitute the entire portfolio. For example, we can define the segmentation of geographic areas where the bank operates composed by four segments {N, S, E, W}. When an obligor defaults, the loss of its assets is added into the entire portfolio loss and into the segments’ loss where each asset resides. CCruncher allows defining and managing multiple segmentations simultaneously.

Expected Loss

Expected Loss is additive respect the portfolio composition, i.e., if L , L_1 , and L_2 are respectively the portfolio, sub-portfolio 1, and sub-portfolio 2 loss distributions in which $L = L_1 + L_2$, then $EL(L_1) + EL(L_2) = EL(L)$. The EL attributable to the k -th segment, in cash, is given by the following statistic:

$$EL_k = \frac{1}{N} \sum_{i=1}^N x_i^k$$

in which x_i^k are the simulated losses of the k -th sub-portfolio. Dividing EL_k by the portfolio EL gives the percentage of EL consumed by this sub-portfolio.

Value at Risk

Value at Risk is not additive, it is not even sub-additive [19]. As mentioned in [3, chap. 5.2.2], calculating VaR risk measure contributions is a natural but difficult attempt. In the CCruncher

case, problems arise when we try to estimate $E(L_k | L = \text{VaR}_\beta(L))$ because the subsample of simulated sub-portfolio losses having a portfolio loss equal to VaR_β has a very small size. For this reason, CCruncher don't provides support for VaR disaggregation.

Expected Shortfall

Expected Shortfall can be disaggregated using the simulated losses of the portfolio and its sub-portfolios. As indicated in [3, chap. 5.2.3], the contribution of k-th sub-portfolio to ES is given by $E(L_k | L \geq \text{VaR}_\beta(L))$. The ES attributable to the k-th segment, in cash, is given by the following statistic:

$$ES_k = \frac{1}{K} \sum_{i=1}^K y_i^k$$

in which x_i are the simulated portfolio losses, x_i^k are the simulated losses of the k-th sub-portfolio, and $\{y_1^k, \dots, y_K^k\} = \{x_i^k | x_i \geq \text{VaR}_\beta\}_{i=1..N}$ are the simulated sub-portfolio losses having a portfolio loss large than VaR_β . Dividing ES_k by the portfolio ES gives the percentage of ES consumed by this sub-portfolio.

Example 3.4 — Risk disaggregation. We use the identical scenario as that described in Example 3.3 and that corresponds to CCruncher input file `test05.xml`. This example defines a segmentation named *sector-rating* with segments S1-A, S1-B, S1-C, S2-A, S2-B, S2-C. Segment X-Y is composed by the obligors located at sector X with the rating Y; each segment has ≈ 166 obligors. Figure 3.6 displays the risk disaggregated respect this segmentation.

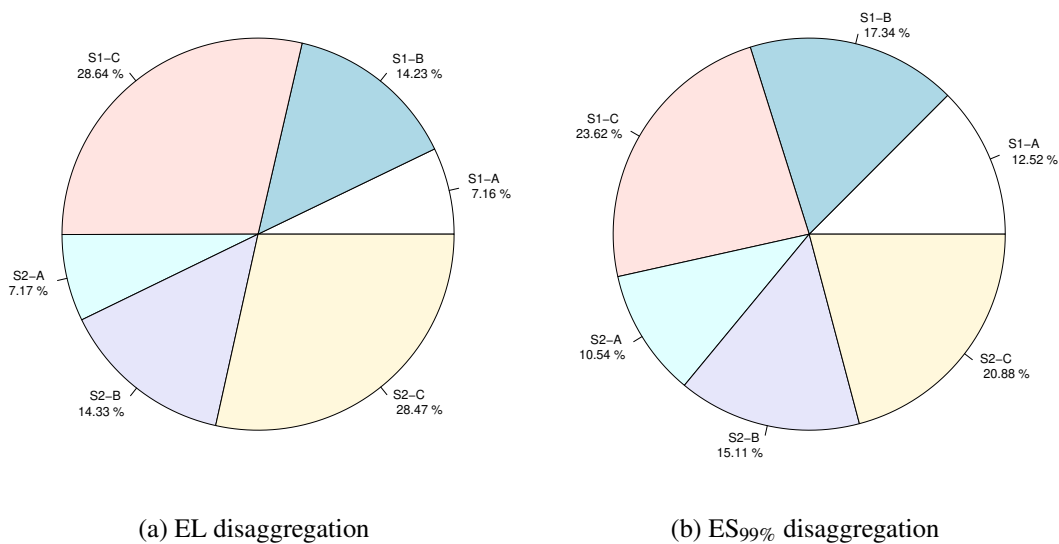


Figure 3.6: Risk disaggregation

3.3 Simulation Details

Antithetic variates. CCruncher generates random samples from the portfolio loss distribution. This sample is used to estimate the risk measures (e.g., VaR) using the statistics described in Section 3.2 or any other required by the analyst. Any one of these measures, considered

individually, can be formulated as a Monte Carlo problem and can apply some variance reduction techniques (e.g., importance sampling [4]), but not simultaneously. CCruncher implements the antithetic variates method (see Appendix A.7) in order to increase the simulation speed instead to reduce the EL variance.

In CCruncher the random sampling of the portfolio loss distribution is performed using $g(X)$ in which X is a multivariate t-Student distribution and $g(x) = F^{-1}(t_v(x))$. We use the fact that the multivariate t-Student distribution X is symmetrical, that is, (x_1, \dots, x_n) is equiprobable to $(-x_1, \dots, -x_n)$, to apply the antithetic method. Expected Loss is the unique statistic that benefits from the variance reduction. However, rather than reducing the variance, we are interested in the reusing of the simulated random numbers to increase the simulation speed. When we use the antithetic method to sample the default times and there are some stochastic EADs or LGDs, we use distinct EADs' and LGDs' random values for each variate.

EAD and LGD interpolation. Given an asset, CCruncher knows its EAD and LGD values at fixed times $\{t_1, \dots, t_n\}$ provided by the user. When we evaluate these functions in a time value distinct from those provided, we use the following piecewise constant interpolation:

$$\left. \begin{aligned} \text{EAD}(t) &= \text{EAD}(t_k) \\ \text{LGD}(t) &= \text{LGD}(t_k) \end{aligned} \right\} \quad \text{where } t_{k-1} < t \leq t_k$$

PD functions approximation. CCruncher allows defining the PD functions using the transition matrix or defining explicitly the PD values at some fixed times. When PDs are defined using the transition matrix, CCruncher uses Proposition A.2 to derive F_i for each month, and intermediate values are interpolated using cubic splines. When PDs are defined explicitly at some fixed times, then intermediate values are interpolated using cubic splines; however, if the resulting function is not strictly increasing, then CCruncher uses linear splines.

Function composition. To speed up the evaluation of the functions $t = F_r^{-1}(t_v(x))$, involved in the simulation of obligors' default times, we interpolate these functions using splines. We only consider the dates where exist some asset event $\{t_1, \dots, t_k\}$ and its values $x_k = t_v^{-1}(F_r(t_k))$. For each rating r , CCruncher try to minimize the number of interpolation points applying the following procedure:

- The initial linear spline $S_r(\cdot)$ is defined at (x_1, t_1) and (x_k, t_k) .
- Repeat until $\max |t_i - S_r(x_i)| < 1$ hour for all $i = 1, \dots, k$
 - Let t_i be the date on which maximum error is achieved
 - Add the interpolation point (x_i, t_i) to the cubic spline function $S_r(\cdot)$.

Spline function $S_r(\cdot)$ use cubic splines; however, if the resulting function is not strictly increasing, then CCruncher uses linear splines.

Interest rate. CCruncher allows considering the Net Present Value of the simulated losses. This is computed using an interest rate curve defined at fixed times and interpolated using linear or cubic splines (selected by user). Allowable methods are simple interest, compound interest and continuous interest. The formulas used to compute the present value are as follows:

$$\begin{aligned} V_0 &= \frac{V_t}{1+r\Delta t} & V_0 &= \frac{V_t}{(1+r)^{\Delta t}} & V_0 &= \frac{V_t}{\exp(r\Delta t)} \\ \text{(Simple)} & & \text{(Compound)} & & \text{(Continuous)} \end{aligned}$$

in which V_t is the value at time t and r is the interest rate given by the interest curve at time t . Don't consider this option if your EADs are the sum of cash flow on distinct dates. In this case, we recommend filling the input file with the EADs' present value and setting to 0 the interest rate curve.

Parallel computing. CCruncher parallelizes the portfolio loss sampling, spawning multiple simulation threads, each with its own Mersenne twister Random Number Generator seeded with consecutive values. When the number of obligors is high, data do not fit in the cache memory L1 and L2. The memory transfer is a time-consuming operation that must be minimized. To achieve this, each thread simulates an obligor *blocksize* times (using the corresponding v and \bar{z}) before moving to the next one.

Parameterization. The CCruncher input file contains the data required to simulate the portfolio losses (obligors, assets, EADs, LGDs, ratings, factor correlations, factor loadings, copula to use, interest rates, segmentations, etc.). This file can be quite large due to portfolio description (e.g., 1GB). When we evaluate the risk sensitivity or the effect of parameters' uncertainty, we must slightly modify the input file and reevaluate the portfolio loss distribution multiple times. To simplify this task, CCruncher provides the macro mechanism, that allows using variables that are replaced by the user-defined values anywhere they are located in the input file, and the automatic evaluation of numerical expressions. Combining these two features, macros and the evaluation of numerical expressions, provides a flexible manner in which to parameterize one's input file (e.g., 2000*\$EUR2USD).

4 — Additional Topics

4.1 Risk Sensitivity

In this section we make a basic sensitivity analysis evaluating the variation of a parameter s , in the risk measure V . The goal is to determine the appearance of the curve $V(s)$ (increasing/decreasing, concave/convex, linear/quadratic/logarithmic, asymptotic behavior) and an estimate of the slope given by $\frac{\partial V}{\partial s}$.

It is difficult to draw general guidelines in the absence of a closed formula for the portfolio credit risk that takes into account all the parameters: number of ratings, PDs over time, number of sectors, number of obligors, EADs and LGDs for each asset, factor loadings, factor correlations, and v . For this reason, we lay out this section analyzing the case `test05.xml` exposed in Example 3.3. In this case, we consider that risk measure is $ES_{99\%}$ and parameters subject to sensibility analysis are: n, \vec{w}, R, v .

We sketch the function $ES_{99\%}(s)$ by calculating $ES_{99\%}$ in various consecutive values of the analyzed parameter. Each of these values is obtained by applying the simulation procedure described in Chapter 3. In order to reduce variations caused by the simulation procedure, we use the same RNG seed in each evaluation.

Example 4.1 — Sensitivity to portfolio size. Common sense tells us that a large and homogeneous portfolio must have the same risk (as percentage of the total exposure) that a portfolio with the same characteristics but with twice obligors. Figure 4.1 shows the $ES_{99\%}$ by varying the number of obligors that compose the portfolio described in `test05.xml`. Note that this value is almost equal in range $[200, \infty)$. In the range $[50, 200]$ we can appreciate the impact of the small number of individuals in the risk measure, not very exaggerated otherwise.

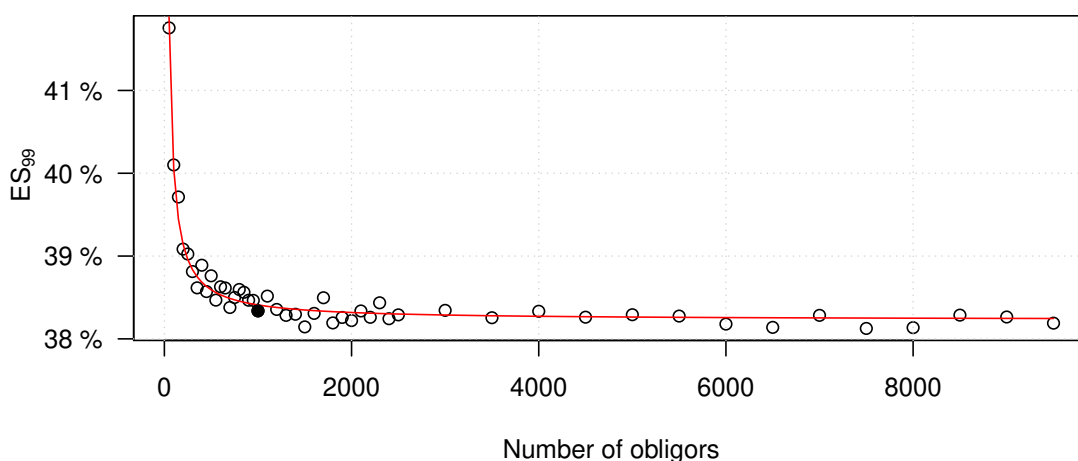


Figure 4.1: $ES_{99\%}$ vs. portfolio size

Two portfolios with distinct number of obligors but the same composition (sector weights, exposures, recoveries) have the same credit risk as long as the number of obligors is high enough. We can use this fact to assess the credit risk of really massive portfolios (e.g. > 0.5 million

obligors). In this cases we can prune the portfolio (preserving the composition) and compute the credit risk of the reduced portfolio.

Example 4.2 — Sensitivity to copula parameters. Figure 4.2 shows the portfolio risk value $ES_{99\%}$ depending on the copula parameters. Each dot is a risk value computed using a simulation, the black dot corresponds to the original value in Example 3.3, and the confidence bounds shows the impact of the parameter value in the error attributable to simulation (this value is magnified in order to appreciate the impact).

If we consider only the range of feasible values (eg. $w_i \in (0, 0.5)$, or $v \in (5, \infty)$), we note that the impact of factor loading loadings and v are similar. However, the impact of factor correlations is much smaller. This is good news, because as we saw in the estimation of the parameters, the uncertainty of factor correlations is much higher than that of factor loadings. We also recall that in the estimation procedure, the values of factor loadings and v are interlaced.

The main drawback of sensitivity analysis is that it measures the impact of the uncertainty of a parameter fixing the other ones. Another drawback is that it does not consider the distribution of the parameters. The section that follows provides a solution to these problems.

4.2 Parameters Uncertainty

The estimation of dependence parameters (degrees of freedom v , factor loadings \vec{w} , and factor correlation matrix R) has an uncertainty that cannot be ignored [32, 17]. Example 2.3 illustrates this phenomenon and its magnitude (see the *20 observations* case).

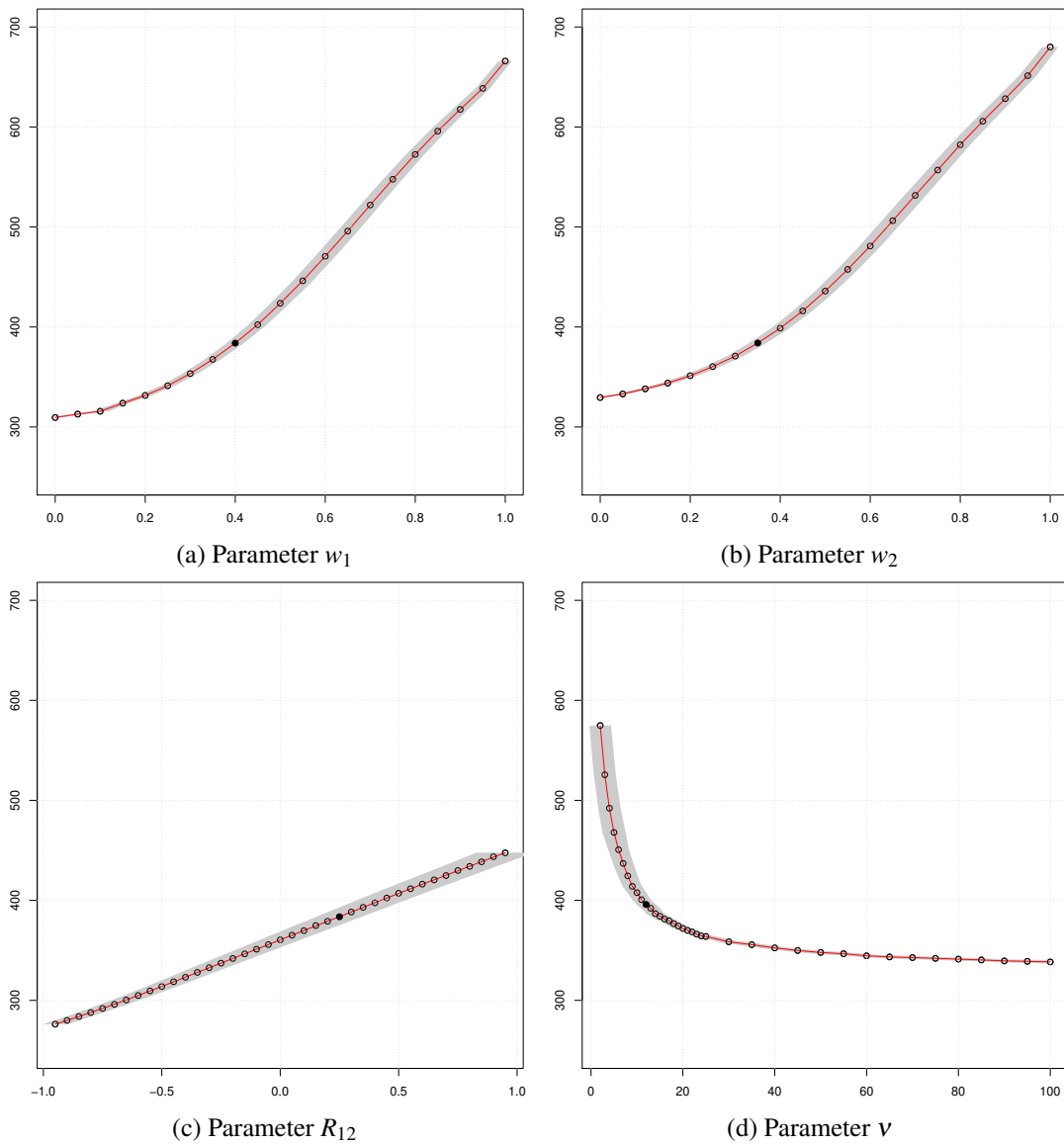
The CCruncher framework encourages incorporating the parameter uncertainty into the credit risk valuation. This indicates that credit risk measures (e.g., $VaR_{99\%}$, $ES_{99\%}$) are regarded as distributions instead of values. For example, in the case of the Value at Risk, we consider the function $VaR_{99\%}(L_T(v, \vec{w}, R))$ that depends on the multivariate random variables v, \vec{w} and R . Note that Expected Loss is not affected by the uncertainty of dependence parameters because the expectation satisfies $E(X + Y) = E(X) + E(Y)$ even if X is not statistically independent of Y .

The solution recommended by CCruncher to determine the credit risk measure distribution is the obvious one and is based on the identical procedure used to determine the distribution of the portfolio loss or the distribution of the parameters: the random sampling of the risk measure distribution.

On one hand, we have a procedure that generates random samples from the multivariate parameters distribution (see Proposition 2.5). These samples can be filtered to obtain a sample of independent parameters values.

On the other hand, we are able to measure the credit risk of a portfolio considering that dependence parameters are fixed values. This is performed sampling the portfolio loss (see Algorithm 3.1) and computing the credit risk measure using the appropriate statistic (see Section 3.2).

We combine both components to define an algorithm that samples the desired credit risk measure. The empirical density given by this sample approximates the risk measure distribution and allows evaluating the effect of the parameter uncertainty. One must be cautious because this procedure unveils the copula parameters uncertainty, but other uncertainties remain hidden such

Figure 4.2: ES_{99%} sensitivity to copula parameters

as: 1) copula selection, 2) factors/sectors selection and assignment, 3) PDs uncertainty, and 4) EAD and LGD models accuracy.

Algorithm 4.1 — Credit risk measure distribution. Assume a portfolio (ratings, sectors, PDs, EADs, LGDs) and the number of defaults observed by sector-rating in the previous years. We can sampling the distribution of a credit risk measure (e.g., VaR_{99%}) by performing the following:

1. Obtain an ergodic sample of the parameters applying the Metropolis-Hastings algorithm (see Proposition 2.5).
2. Remove the first iterations (burn-in period), and determine the minimum common length B at which the ACFs are 0.
3. Let $\{v^t, \vec{w}^t, R^t\}_{t=1, \dots, N}$ be the subsample of the parameters distribution composed by the indexes multiples of B .

4. For each simulated parameter, simulate the portfolio loss distribution $L_T(v^t, \vec{w}^t, R^t)$ using the Algorithm 3.1.
5. For each simulated portfolio loss distribution, compute its risk measure using the statistic described in Section 3.2: $\{\text{VaR}_{99\%}(L_T(v^t, \vec{w}^t, R^t))\}_{t=1, \dots, N}$.

Example 4.3 — Credit risk considering parameters uncertainty. We use the scenario described in Example 3.3. We generate 20 observations of annual numbers of defaults using the Algorithm 2.1 (in the real world, these observations are extracted from the historical records). In the following lines, we apply Algorithm 4.1 to estimate the $\text{ES}_{99\%}$ distribution, considering the parameters' uncertainty.

The first step includes the generation of a sample of the parameters' distribution using the Metropolis-Hastings algorithm. We performed 1,000,000 iterations. Figure 4.3 displays the marginals of this distribution.

The second step is removing the first 20,000 iterations (burn-in period) and determining the thin interval using the ACFs. Figure 4.4 shows the ACF for each parameter. A thin interval of 3000 is sufficient in this example.

The third step is pruning the parameters sample preserving only the simulated values whose index is a multiple of 3000. Below is displayed part of the simulated values:

N	v	w ₁	w ₂	R ₁₂
1	12.54	0.29	0.48	0.21
2	10.77	0.28	0.32	-0.28
⋮	⋮	⋮	⋮	⋮
300	10.06	0.33	0.36	0.02

The fourth step is the simulation of the portfolio loss distribution for each simulated parameter. This is quite an intensive computation. The CCruncher's command line execution and the input file macro expansion can assist in performing this step. Listing 4.1 displays a simple batch script to sequentially execute the required simulations. Those who have a cluster, a cloud or a grid may consider to executing multiple parallel tasks.

Finally, the fifth step is the computation of the risk statistic for each portfolio loss distribution. The density of the $\text{ES}_{99\%}$ can be approximated using the empirical density. Listing 4.2 provides an R script to perform this task.

The exact value of the $\text{ES}_{99\%}$ assuming that parameters are fixed and known is $\approx 383 \text{ €}$ (see Example 3.3). If we do not assume that parameters are known and we infer the parameters from 20 observations of the numbers of defaults, then we obtain the distribution displayed in Figure 4.5. The key point is to dispose a distribution indicating the feasible values that can take the risk measure $\text{ES}_{99\%}$.

Listing 4.1: Execution of multiple CCrunchers (bash script)

```
mkdir data/MH001;
bin/ccruncher -cmd -o data/MH001 -DNU=12.54 -DW1=0.29 \
  -DW2=0.48 -DR12=0.21 samples/test05.xml > data/MH001/ccruncher.out;
mkdir data/MH002;
```

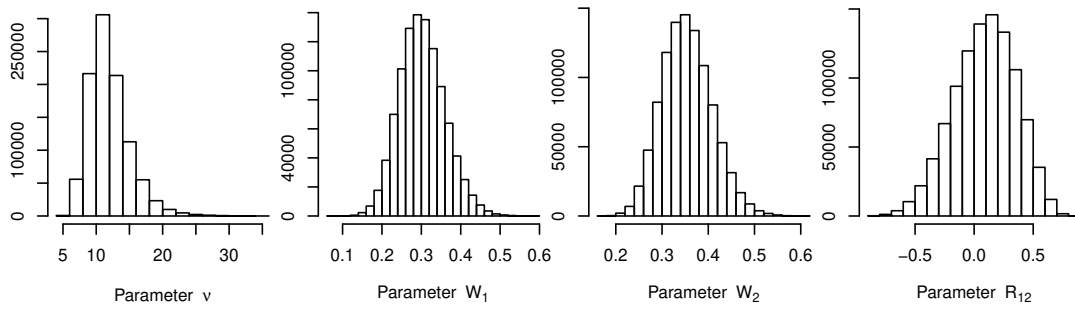


Figure 4.3: Step 1 – Parameters sampling using M-H

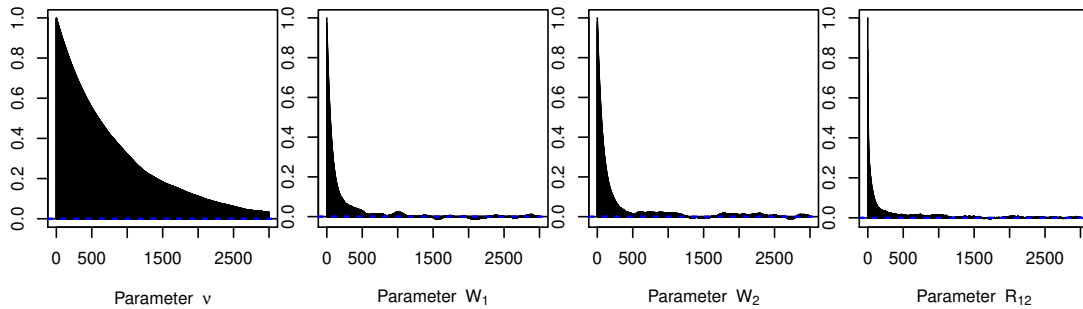


Figure 4.4: Step 2 – ACF of simulated parameters

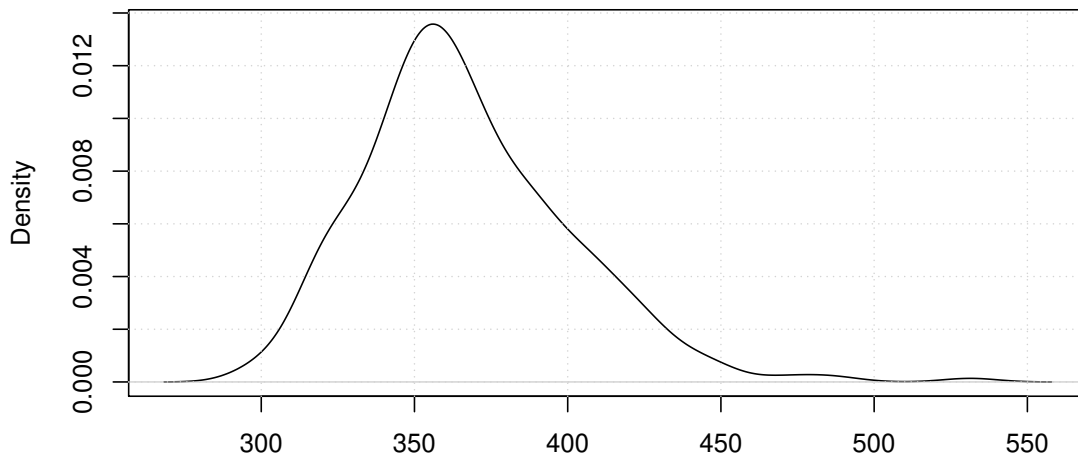
```
bin/ccrunner -cmd -o data/MH002 -DNU=10.77 -DW1=0.28 \
  -DW2=0.32 -DR12=-0.28 samples/test05.xml > data/MH002/ccrunner.out;
...
mkdir data/MH300;
bin/ccrunner -cmd -o data/MH300 -DNU=10.06 -DW1=0.33 \
  -DW2=0.36 -DR12=0.02 samples/test05.xml > data/MH300/ccrunner.out;
```

Listing 4.2: $ES_{99\%}$ distribution (R script)

```
getRisk <- function(dir)
{
  filename = paste(dir, "/portfolio.csv", sep="")
  data <- read.csv(filename, comment.char="#")
  X = sort(data[,1])
  n = length(X)
  Y = X[as.integer(n*0.99):n]
  ES99 = mean(Y)
  sde = sqrt(var(Y)/length(Y))
  return(c(ES99, sde))
}

dirs = dir("data", pattern="MH[[:digit:]]{3}*", full.names=TRUE)
values = matrix(ncol=2, nrow=length(dirs))
colnames(values) = c("ES99", "stderr")
for(i in 1:length(dirs)) {
  values[i,] = getRisk(dirs[i])
}

plot(density(values[,1]))
```

Figure 4.5: ES_{99%} distribution

4.3 Portfolio Optimization

Once we know the portfolio credit risk, naturally arises the question “How can I balance the weights assigned to each business unit so that the total portfolio risk is minimized?” In Section 3.2.2, we have seen that CCruncher allows defining sub-portfolios and it can disaggregate the portfolio credit risk in terms of them. In this section we will see how to modify the weights assigned to each sub-portfolio in order to optimize the portfolio credit risk. First of all, let’s define the problems considered:

Definition 4.1 — Portfolio optimization. Let a portfolio composed of K sub-portfolios. We want to determine the sub-portfolio weights α_k that minimize one of the following problems¹.

EL minimization

$$\min_{\alpha_k} \text{EL}(L)$$

Subjecto to:

$$\left. \begin{array}{l} x^0 = \text{ES}_{99\%}(L) \\ 0 \leq \alpha_k \leq 1 \quad \forall k \\ \sum_{k=1}^K \alpha_k = 1 \end{array} \right\}$$

ES minimization

$$\min_{\alpha_k} \text{ES}_{99\%}(L)$$

Subjecto to:

$$\left. \begin{array}{l} y^0 = \text{EL}(L) \\ 0 \leq \alpha_k \leq 1 \quad \forall k \\ \sum_{k=1}^K \alpha_k = 1 \end{array} \right\}$$

EL-ES minimization

$$\min_{\alpha_k} \sqrt{\text{EL}^2 + \text{ES}_{99\%}^2}$$

Subjecto to:

$$\left. \begin{array}{l} 0 \leq \alpha_k \leq 1 \quad \forall k \\ \sum_{k=1}^K \alpha_k = 1 \end{array} \right\}$$

where $L = L_1 + \dots + L_k$ is the portfolio loss, and x^0 and y^0 are fixed values.

Sub-portfolio weights α_k are the sub-portfolio exposures relative to total portfolio exposure at 1-year. The way to balance the portfolio composition is to add or subtract obligors preserving the characteristics of each sub-portfolio. When we try to model this procedure we find two problems: 1) losses of portfolios with distinct composition are not comparable, 2) it is unclear the procedure

¹We use ES_{99%} but you can use any other risk measure depending on copula parameters, such as ES_{95%} or VaR_{99%}.

to modify the composition of a sub-portfolio (we add half a debtor?, to wich rating and sector we assign the new obligor?, new obligors have an averaged exposure and recovery?).

Fortunately these two problems are solvable. The first one is avoided considering the relative losses on exposure rather than absolute losses. Thus, the relative portfolio loss at time horizon T is

$$L_T = \frac{\sum_{i=1}^n \mathbb{1}_{[0,T]}(t_i) \cdot \left(\sum_{j=1}^{m_i} \text{EAD}_i^j(t_i) \cdot \text{LGD}_i^j(t_i) \right)}{\sum_{i=1}^n \left(\sum_{j=1}^{m_i} \text{avg}_{[0,T]} \left(\text{EAD}_i^j(t) \right) \right)}$$

where T is the time horizon where portfolio loss is computed, n is the number of obligors, (t_1, \dots, t_n) are the obligors' default times, m_i is the number of assets of the i -th obligor, and $\text{EAD}_i^j(t)$ and $\text{LGD}_i^j(t)$ are the EAD and LGD functions of the j -th asset of the i -th obligor evaluated at time t .

The second problem, the portfolio composition, is avoided by equating the number of obligors to exposures. Suppose we have two sub-portfolios A and B with the same exposure, the first one with 500 obligors, and the second one with 500 obligors. We want to balance the portfolio so that ends up with 25% of the exposure in A and 75% in B. The way to implement it is eliminating half of the obligors from A and adding 250 new obligors similar to those of B. That said, if sub-portfolios A and B have a large number of homogeneous individuals, then we can maintain the original number of individuals (500 + 500) and balance the portfolio by multiplying by 0.5 the A's exposures and by 1.5 the B's exposures.

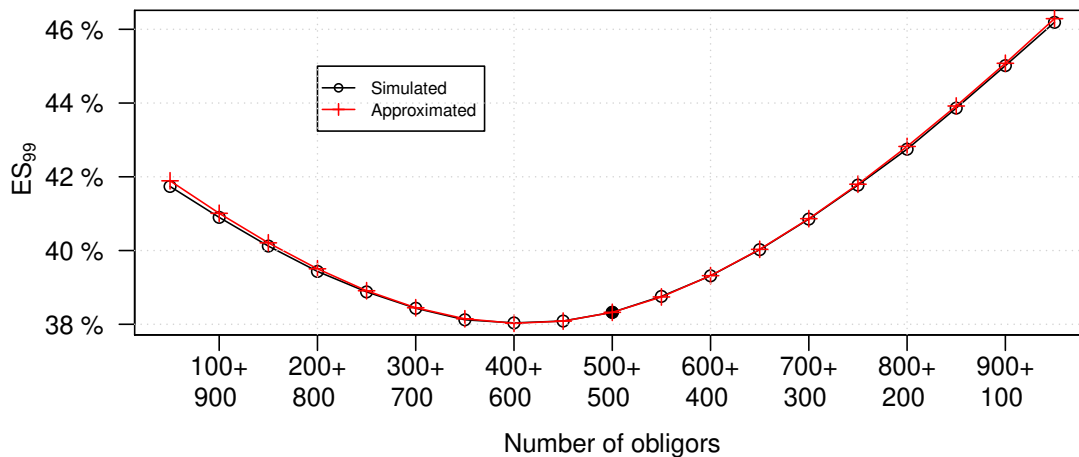


Figure 4.6: $\text{ES}_{99\%}$ for distinct percentages of S1 and S2

Example 4.4 — Sub-portfolio weights. To check the accuracy of the equivalence between the number of individuals and the exposures we analyze the case `test05.xml` (see Example 3.3) considering the sub-portfolios S1 (= S1A + S1B + S1C) and S2 (= S2A + S2B + S2C). In this example we consider only two sub-portfolios, rather than the six defined, so that the values can be clearly represented in the plane. First, we simulate different portfolio compositions (50+950, 100+900, ..., 950+50) and calculate the $\text{ES}_{99\%}$ for each of them. Next, we consider a fixed number of individuals (500+500) and we change the exposure of individuals according to its sector so that we mimic the previous compositions. Figure 4.6

displays the values modifying the composition of the portfolio (simulated) and the values modifying the exposure of the obligors (approximated). In this case, both values are similar, with an error less than 0.15%.

The equivalence between the number of debtors and the exposures has a very important advantage since it allows optimizing without having to do multiple simulations. This is achieved by simulating the losses of the current portfolio and by reusing these values to calculate any desired composition of the portfolio. The following algorithm computes the risk of a weighted portfolio reusing the simulated losses from the current portfolio.

Algorithm 4.2 — Risk of a weighted portfolio. Let a portfolio composed of K sub-portfolios with an average exposure E_k for each of them in the period $[0, T]$. We have simulated the portfolio losses broken down by sub-portfolios. We want compute the EL and $ES_{99\%}$ of this portfolio with weights $\alpha_1, \dots, \alpha_K$.

1. Obtain multiplicative factors: $\lambda_k = \alpha_k \cdot \frac{E}{E_k} = \frac{\alpha_k}{\alpha_k^0}$
2. Create weighted losses: $y_i = \sum_{k=1}^K \lambda_k \cdot x_i^k \quad i = 1, \dots, N$
3. Compute EL (see Section 3.2) using values $\{y_i\}_{i=1, \dots, N}$ and divide this value by E to obtain the EL relative to total exposure.
4. Compute $ES_{99\%}$ (see Section 3.2) using values $\{y_i\}_{i=1, \dots, N}$ and divide this value by E to obtain the $ES_{99\%}$ relative to total exposure.

where $E = \sum_{k=1}^K E_k$ is the total exposure, α_k^0 is the weight of the current k -th sub-portfolio, N is the number of simulated losses, and x_i^k is the i -th simulated loss of the k -th sub-portfolio.

The risk of weighted portfolio is continuous with respect to $\alpha_1, \dots, \alpha_K$ but it is not linear with respect to this parameters because the rank statistic involved in the $VaR_{99\%}$ and $ES_{99\%}$ measures. This is clearly shown in Figure 4.6.

Example 4.5 — Feasible portfolios. We consider the portfolio defined in `test05.xml` (see Example 3.3). We want to know the values that can take the risk measures by changing the composition of the portfolio. To achieve this, we consider the following weights:

$$\left\{ (\alpha_1, \dots, \alpha_6) \mid \alpha_i \in \{0, 0.1, 0.2, \dots, 0.9, 1\} \text{ and } \sum_{i=1}^6 \alpha_i = 1 \right\}$$

This set of weights has 3003 elements. For each element of this set we compute the EL and $ES_{99\%}$ of the weighted portfolio applying the Algorithm 4.2. Result is displayed in Figure 4.7. We have drawn the convex hull of the region and we have identified the relevant points. The original portfolio is marked with a black dot and the corresponding optimal portfolios are indicated with a arrow. The appearance of the EL-ES region is determined by the involved sub-portfolios, in this case: S1A, S1B, S1C, S2A, S2B, S2C. Other segmentations (e.g., geographic areas, types of products) will lead to EL-ES regions with a different shape.

Now we have all the elements to solve the three optimization problems. Instead of presenting the algorithm in pseudo-code, this time we use a basic R script exposed in Listing 4.3 to illustrate the procedure. The optimized portfolio is the case `test05.xml` and optimization results are listed in Table 4.1. Note that $K - 1$ parameters are optimized rather than K . This is because we can obtain the former doing $\alpha_1 = 1 - \sum_{i=2}^K \alpha_i$. Note also that penalty functions are used to ensure the

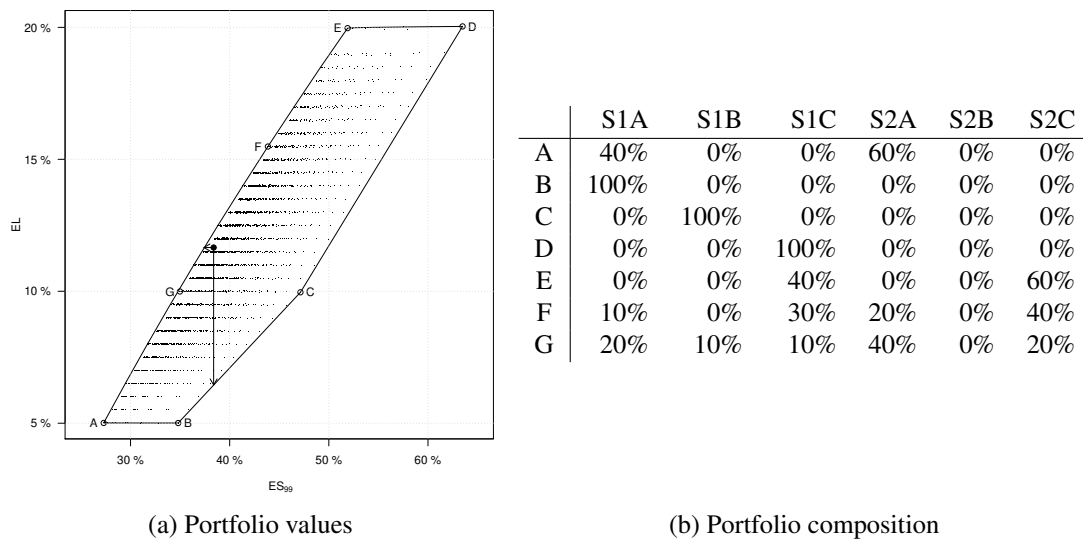


Figure 4.7: test05.xml feasible portfolios

restrictions compliance.

Listing 4.3: Portfolio optimization (R script)

```
evalPortfolio <- function(losses, w0, w1, level=0.99)
{
  if (length(w0) != length(w1)) stop("invalid_parameters")
  if (ncol(losses) != length(w0)) stop("invalid_losses")
  if (level <= 0 || level >= 1) stop("ES_level_out-of-range_(0,1)")
  n = nrow(losses)
  k = ncol(losses)
  lambda = w1/w0
  loss = rep(0, n)
  for(i in 1:k) {
    loss = loss + losses[,i]*lambda[i]
  }
  loss = sort(loss)
  e1 = mean(loss)
  es = mean(loss[as.integer(n*level):n])
  return(c(e1, es))
}

getW <- function(params)
{
  w = c(0, params)
  w[1] = 1-sum(w)
  return(w)
}

f <- function(params, losses, w0, option, val, level=0.99)
{
  if (!(option %in% (1:3))) stop("unrecognized_option")
  w1 = getW(params)
  aux = 0
  if (w1[1] < 0) {
    aux = w1[1]
    w1[1] = 0
  }
}
```

```

eles = evalPortfolio(losses, w0, w1, level)
if (aux != 0) {
  eles = eles + abs(aux)*10
}
if (option == 1) {
  ret = eles[1] + 10*(val-eles[2])^2
  return(ret)
}
else if (option == 2) {
  ret = eles[2] + 10*(val-eles[1])^2
  return(ret)
}
else if (option == 3) {
  ret = sqrt(sum(eles^2))
  return(ret)
}
else return(NA)
}

losses <- read.csv("data/sector-rating.csv", comment.char="#")
exposures = c(167, 166, 167, 167, 167, 166)
w0 = exposures/sum(exposures)
eles0 = evalPortfolio(losses, w0, w0)

# EL minimization
params = w0[2:length(w0)]
opt = nlminb(params, f, losses=losses, w0=w0, option=1, val=eles0[2],
  lower=rep(0,length(params)), upper=rep(1,length(params)))
w1 = getW(opt$par)
evalPortfolio(losses, w0, w1)

# ES minimization
params = w0[2:length(w0)]
opt = nlminb(params, f, losses=losses, w0=w0, option=2, val=eles0[1],
  lower=rep(0,length(params)), upper=rep(1,length(params)))
w2 = getW(opt$par)
evalPortfolio(losses, w0, w2)

# optimization
params = w0[2:length(w0)]
opt = nlminb(params, f, losses=losses, w0=w0, option=3, val=NA,
  lower=rep(0,length(params)), upper=rep(1,length(params)))
w3 = getW(opt$par)
evalPortfolio(losses, w0, w3)

```

	S1A	S1B	S1C	S2A	S2B	S2C	EL	ES
current	16.7%	16.6%	16.7%	16.7%	16.7%	16.6%	11.6%	38.3%
min EL	68.9%	31.1%	0%	0%	0%	0%	6.6%	38.3%
min ES	18.7%	0%	21.3%	36.8%	0%	23.2%	11.6%	37.2%
min EL-ES	41.0%	0%	0%	59.0%	0%	0%	5.0%	27.0%

Table 4.1: test05.xml optimization

A — Appendices

A.1 The Probabilities of Default

Definition A.1 — Probability of Default (PD). T_i is the random variable that represents the default time (aka time-until-default, aka survival time, aka time-to-default) of the i -th obligor. We denote as $F_i(t)$ the cumulative distribution function of T_i :

$$F_i(t) = \Pr\{T_i \leq t\} = \Pr\{\text{i-th obligor defaults before } t \text{ years}\}$$

Note that $F_i(t)$ is the probability of default (PD) of the i -th obligor at time horizon t . PD is a key parameter used in the calculation of credit risk under Basel II. We indicate as p_i the probability that i -th obligor defaults within the time range $[0, 1]$.

$$p_i = \Pr\{T_i \leq 1 \text{ year}\} = F_i(1)$$

Note that the economic cycle is written on the default times distribution shape. Figure A.1 depicts the default time density of an obligor with current rating B , where bumps represent the lean years, and its corresponding cdf (aka PD function).

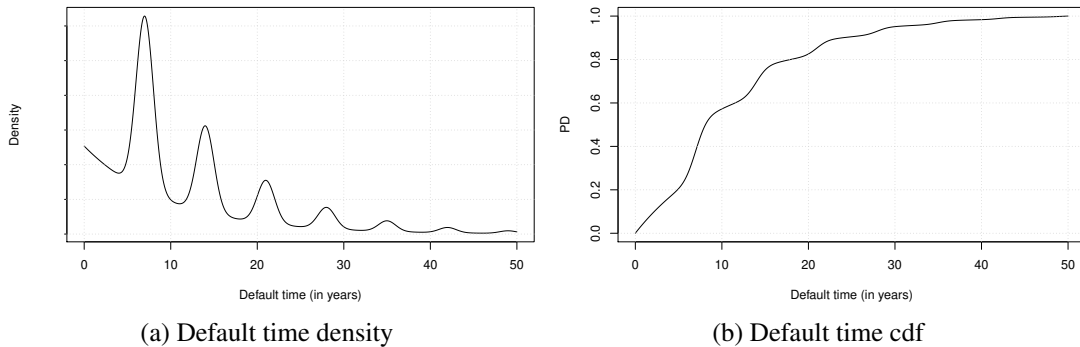


Figure A.1: Probability of Default function

CCruncher allows you to specify the PDs functions declaring the PDs' values at user-defined time points. These values must fulfill the following conditions:

- $F_i(0) = 0$ $i = 1, \dots, r$ $i \neq \text{default}$
- $F_i(t)$ strictly increasing $i = 1, \dots, r$ $i \neq \text{default}$
- $F_i(\infty) = 1$ $i = 1, \dots, r$
- $F_{\text{default}}(t) = 1$ $\forall t \geq 0$

The functions specified in this manner are interpolated using cubic splines when the above conditions are preserved and are interpolated using linear splines when they are not.

In general, what we know of the PD functions is limited. In some cases we have the rating transition matrix, in others we only have the 1-year PD; if obligors are listed at the stock market then we can use the Merton's model to estimate the PD functions. Below, these cases are analyzed. All three manners produce similar results in the short term (e.g., 1 year) but differ in the middle and long term.

A.1.1 Obtaining PDs from the transition matrix

Definition A.2 — Transition matrix. The T -years transition matrix gives the probability of changing from i -th rating to j -th rating in a period of T years:

$$M_T = \begin{pmatrix} m_{11} & \cdots & m_{1r} \\ \vdots & \ddots & \vdots \\ m_{r1} & \cdots & m_{rr} \end{pmatrix}$$

in which r is the number of ratings and m_{ij} is the probability that an obligor with i -th rating ends up having the j -th rating after T years. A valid transition matrix must have its elements in the range $[0, 1]$, the row sums equal to 1, a unique *defaulted* rating, and any rating ends up defaulting.

Example A.1 — 1-year transition matrix. Table A.1 shows a transition matrix, extracted from [18, p. 20], in which the probability that an obligor with rating AA changes to rating B in one year is 0.14%. Note that the last column contains the 1-year default probabilities for each rating and that the last row corresponds to the *defaulted* rating.

	AAA	AA	A	BBB	BB	B	CCC	Default
AAA	90.81	8.33	0.68	0.06	0.12	0.00	0.00	0.00
AA	0.70	90.65	7.79	0.64	0.06	0.14	0.02	0.00
A	0.09	2.27	91.05	5.52	0.74	0.26	0.01	0.06
BBB	0.02	0.33	5.95	86.93	5.30	1.17	0.12	0.18
BB	0.03	0.14	0.67	7.73	80.53	8.84	1.00	1.06
B	0.00	0.11	0.24	0.43	6.48	83.46	4.07	5.21
CCC	0.22	0.00	0.22	1.30	2.38	11.24	64.86	19.78
Default	0.00	0.00	0.00	0.00	0.00	0.00	0.00	100.00

Table A.1: 1-year transition matrix %

Proposition A.1 — Composed and scaled transition matrix. The transition matrix can be composed and scaled in time using the following rules:

- $M_{T_1+T_2} = M_{T_1} \cdot M_{T_2}$
- $M_{k \cdot T} = M_T^k$
- $M_{\frac{T}{k}} = \sqrt[k]{M_T}$

The root of a matrix M can be obtained using the spectral decomposition $M = P \cdot D \cdot P^{-1}$ doing $M^k = P \cdot D^k \cdot P^{-1}$, in which P and D are the eigenvectors and eigenvalue matrices of M .

Sometimes, the scaled transition matrix does not satisfy the Markov conditions (the row sum is equal to one, and all elements are non-negatives). In this case, we must transform this matrix to the relevant Markov matrix. This process is called regularization. Below is the QOM regularization algorithm extracted from [23].

Algorithm A.1 — Transition matrix regularization. The QOM (Quasi-Optimization of the root Matrix) algorithm regularizes an $n \times n$ transition matrix M , row by row. The steps to regularizing the i -th row are

1. Compute the difference between the row sum and one. Divide by n and subtract this value from all non-zero components:

$$m_{ij} \neq 0 \implies m_{ij} = m_{ij} - \frac{1}{n} \left(\sum_{k=1}^n m_{ik} - 1 \right)$$

2. If all the row elements are non-negative and add up to one, then stop: the row is regularized.
3. Adjust any negative row element to zero and go to Step 1.

The algorithm stops after m steps, where $m \leq n$. Apply the previous algorithm to every row. The final matrix is a regularized matrix.

Deriving PD functions from the transition matrix implies that transition probabilities remain constant over time, undermining the economic cycle effect. In practice, this constancy is equivalent to considering an averaged economic cycle. See document [11, sec. 3.2.1] to obtain additional information related with this section and the CreditMetricsTM approach.

Proposition A.2 — PDs derived from transition matrix. If M_T is a transition matrix, then the time-until-default distribution for an obligor with the i -th rating is

$$F_i(t) = \Pr\{T_i \leq t\} = (M_t)_{id}$$

in which T_i is the time-until-default of an obligor with the i -th rating, d is the index of the *defaulted* rating, M_t is the transition matrix scaled to time t , and $(M)_{ij}$ is the matrix element located in the i -th row and the j -th column.

Example A.2 We derive the time-until-default distribution for each rating induced by the transition matrix presented in Table A.1. As a first step, we scale the 1-year transition matrix M_1 to the 1-month transition matrix $M_{\frac{1}{12}}$.

$$M_{\frac{1}{12}} = \begin{pmatrix} 99.1972 & 0.7588 & 0.0320 & 0.0020 & 0.0112 & -0.0011 & -0.0001 & 0.0000 \\ 0.0635 & 99.1744 & 0.7083 & 0.0392 & 0.0015 & 0.0120 & 0.0018 & -0.0007 \\ 0.0074 & 0.2057 & 99.1980 & 0.5102 & 0.0557 & 0.0189 & -0.0001 & 0.0042 \\ 0.0014 & 0.0239 & 0.5507 & 98.8021 & 0.5164 & 0.0871 & 0.0077 & 0.0106 \\ 0.0027 & 0.0113 & 0.0396 & 0.7581 & 98.1557 & 0.8774 & 0.0889 & 0.0665 \\ -0.0007 & 0.0098 & 0.0196 & 0.0111 & 0.6447 & 98.4443 & 0.4463 & 0.4240 \\ 0.0233 & -0.0023 & 0.0170 & 0.1287 & 0.2166 & 1.2298 & 96.4213 & 1.9667 \\ 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 100 \end{pmatrix}$$

We observe that $M_{\frac{1}{12}}$ is not a regular matrix because of negative values. We apply the QOM

algorithm to regularize it.

$$\bar{M}_{12} = \begin{pmatrix} 99.1969 & 0.7586 & 0.0317 & 0.0018 & 0.0110 & 0.0000 & 0.0000 & 0.0000 \\ 0.0634 & 99.1744 & 0.7082 & 0.0391 & 0.0014 & 0.0119 & 0.0017 & 0.0000 \\ 0.0074 & 0.2057 & 99.1980 & 0.5102 & 0.0557 & 0.0189 & 0.0000 & 0.0041 \\ 0.0014 & 0.0239 & 0.5507 & 98.8021 & 0.5164 & 0.0871 & 0.0077 & 0.0106 \\ 0.0027 & 0.0112 & 0.0395 & 0.7581 & 98.1557 & 0.8774 & 0.0889 & 0.0665 \\ 0.0000 & 0.0099 & 0.0197 & 0.0111 & 0.6447 & 98.4444 & 0.4463 & 0.4240 \\ 0.0228 & 0.0000 & 0.0165 & 0.1282 & 0.2161 & 1.2293 & 96.4208 & 1.9663 \\ 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 0.0000 & 100 \end{pmatrix}$$

Finally, we use Proposition A.2 to determine the PD function for each rating and month. The results are displayed in Figure A.2.

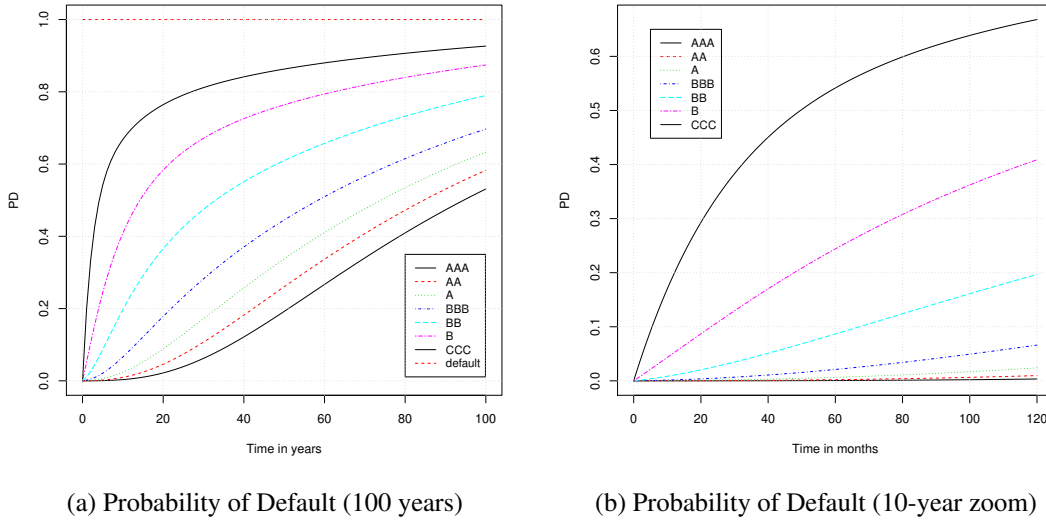


Figure A.2: PDs derived from transition matrix

A.1.2 Obtaining PDs from annual values

In some circumstances, the transition matrix is not available, and we only have the probability of default at 1-year for each rating, p_1, \dots, p_r in which the r -th rating is the *defaulted* status. In the absence of more information, we can consider a diagonal transition matrix such as the following:

$$M_1 = \begin{pmatrix} 1-p_1 & \cdots & 0 & p_1 \\ \vdots & \ddots & \vdots & \vdots \\ 0 & \cdots & 1-p_{r-1} & p_{r-1} \\ 0 & \cdots & 0 & 1 \end{pmatrix}$$

This simplified case has a closed-form solution stated by the following proposition.

Proposition A.3 — PDs derived from annual values. If the time-until-default of an obligor satisfies a transition matrix and only depends on the 1-year probability of default p_i ,

then this time-until-default follows the exponential distribution, i.e., $T_i \sim \text{Exp}(-\ln(1 - p_i))$.

$$F_i(t) = \Pr\{T_i \leq t\} = 1 - e^{-\lambda_i t} \quad \text{where } \lambda = -\ln(1 - p_i) \quad i = 1, \dots, r$$

Proof. We note that ratings are detached from one another. We use Proposition A.2 to obtain the time-until-default distribution of the i -th rating:

$$\begin{pmatrix} 1-p_i & p_i \\ 0 & 1 \end{pmatrix}^t = \begin{pmatrix} 1 & 1 \\ 1 & 0 \end{pmatrix} \cdot \begin{pmatrix} 1 & 0 \\ 0 & 1-p_i \end{pmatrix}^t \cdot \begin{pmatrix} 0 & 1 \\ 1 & -1 \end{pmatrix} = \begin{pmatrix} (1-p_i)^t & 1-(1-p_i)^t \\ 0 & 1 \end{pmatrix}$$

Then,

$$F_i(t) = 1 - (1 - p_i)^t = 1 - e^{-\lambda_i t} \quad \text{where } \lambda_i = -\ln(1 - p_i)$$

□

Example A.3 Suppose that we only know the annual probabilities of default and that these values are those from the 1-year regularized transition matrix displayed in Table A.1. We call the 1-year regularized transition matrix to $\bar{M}_1 = \bar{M}_{\frac{1}{12}}^{12}$ where $\bar{M}_{\frac{1}{12}}$ is computed in Example A.2. In this case, the 1-year diagonal transition matrix is

$$M_1 = \begin{pmatrix} 99.9991 & 0 & 0 & 0 & 0 & 0 & 0 & 0.0009 \\ 0 & 99.9923 & 0 & 0 & 0 & 0 & 0 & 0.0077 \\ 0 & 0 & 99.9400 & 0 & 0 & 0 & 0 & 0.0600 \\ 0 & 0 & 0 & 99.8200 & 0 & 0 & 0 & 0.1800 \\ 0 & 0 & 0 & 0 & 98.9401 & 0 & 0 & 1.0599 \\ 0 & 0 & 0 & 0 & 0 & 94.7997 & 0 & 5.2003 \\ 0 & 0 & 0 & 0 & 0 & 0 & 80.2154 & 19.7846 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 100 \end{pmatrix}$$

We determine the PD's function for each rating using Proposition A.3, i.e., considering that times-until-default are exponential distributions. These results are displayed in Figure A.3. Observe that values around $T = 1$ are similar to those of Example A.2; however, they differ as they move away from this.

A.1.3 Obtaining PDs from Merton's model

In this section we expose the probability of default functions resulting of the Merton's model [3, sec. 3.3]. This method assesses the probability of default of a listed company based on market information. If we assume that the firm's debt amount remains constant over time considering devaluation, then the probability of default under the Merton's framework [6] is

$$F_i(t) = \Phi \left(-\frac{\ln \left(\frac{A_i}{D_i e^{rt}} \right) + \left(\mu_i - \frac{\sigma_i^2}{2} \right) \cdot t}{\sigma_i \sqrt{t}} \right)$$

where Φ is the univariate Gaussian cdf, A_i is the firm's asset value, D_i is the book value of the debt, μ_i is the firm's asset value drift, σ_i is the firm's asset value volatility, and r is the risk free rate. We determine the conditions under which the function F_i is a valid cumulative distribution

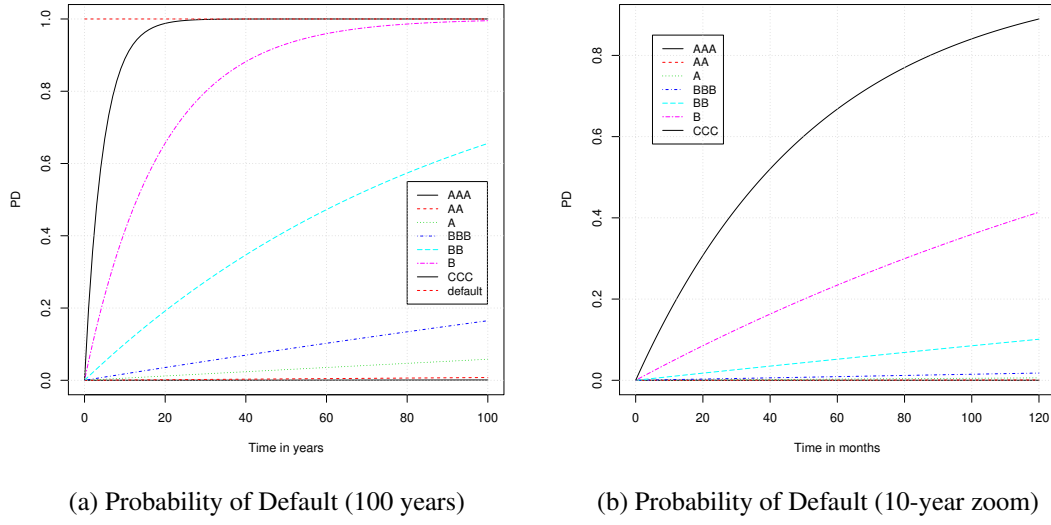


Figure A.3: PDs derived from annual values

function.

$$F_i(0) = 0 \implies \lim_{t \rightarrow 0} -\frac{\ln\left(\frac{A_i}{D_i \cdot e^{rt}}\right) + \left(\mu_i - \frac{\sigma_i^2}{2}\right) \cdot t}{\sigma_i \sqrt{t}} = -\infty \implies \ln\left(\frac{A_i}{D_i}\right) > 0 \implies A_i > D_i$$

$$F_i(\infty) = 1 \implies \lim_{t \rightarrow \infty} -\frac{\ln\left(\frac{A_i}{D_i \cdot e^{rt}}\right) + \left(\mu_i - \frac{\sigma_i^2}{2}\right) \cdot t}{\sigma_i \sqrt{t}} = \infty \implies -r + \mu < \frac{\sigma_i^2}{2}$$

$$F_i'(t) > 0 \implies -2\mu + 2r + \sigma^2 > 0 \implies -r + \mu < \frac{\sigma_i^2}{2}$$

Both conditions are reasonable. The first one indicates that the current market value of the company must exceed the value of its debt. The second one tells us that the expected annual return should be close to the risk-free rate.

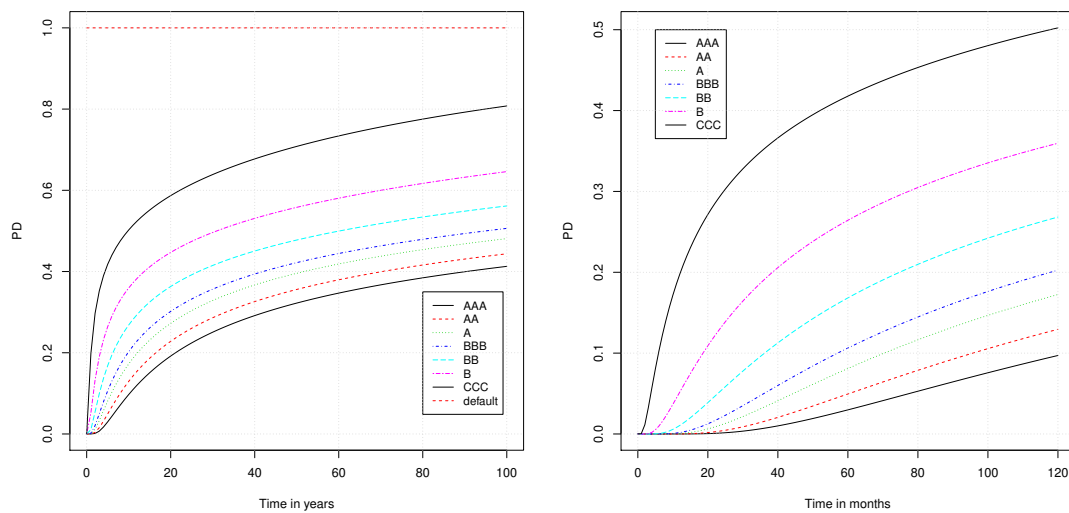
Example A.4 We draw the Merton's probability of default functions so that they are comparable to the functions of the previous examples. For this reason we consider that the annual probabilities of default are those from the 1-year regularized transition matrix displayed in Table A.1. Furthermore we fix the asset value drifts to the risk free rate ($\mu_i = r$) and assumes, arbitrary, that $A_i/D_i = 1.2$. The remaining unknowns σ_i are determined resolving the equation

$$F_i(1) = p_i = \Phi\left(-\frac{\ln\left(\frac{A_i}{D_i}\right) - \frac{\sigma_i^2}{2}}{\sigma_i}\right)$$

The following table exposes the parameters of the probabilities of default functions.

	p_i (in %)	A_i/D_i	μ_i	σ_i
AAA	0.0009	1.2	$r \rightarrow$	0.0422
AA	0.0077	1.2	$r \rightarrow$	0.0479
A	0.0600	1.2	$r \rightarrow$	0.0558
BBB	0.1800	1.2	$r \rightarrow$	0.0620
BB	1.0599	1.2	$r \rightarrow$	0.0778
B	5.2003	1.2	$r \rightarrow$	0.1085
CCC	19.7846	1.2	$r \rightarrow$	0.1928

Finally, we use these values in the $F_i(t)$ Merton's formula to draw the probabilities of default exposed in Figure A.4.



(a) Probability of Default (100 years)

(b) Probability of Default (10-year zoom)

Figure A.4: PDs derived from Merton's model

A.2 Copula Basics

It is common to use correlation as a measure of dependency between random variables. In most cases, this measure does not fully reflect the structure of dependence between the variables. Figure A.5 displays two cases with the identical mean, marginals and correlations but a distinct risk profile because the two cases' underlying dependence structure is distinct (Gaussian copula vs. t-Student(3) copula). The mathematical concept that does reflect the structure of dependence between random variables, however, is the copula, which we define below. The world of copulas is vast and quite interesting. See [26, chap. 5] for a pleasant introduction to copulas.

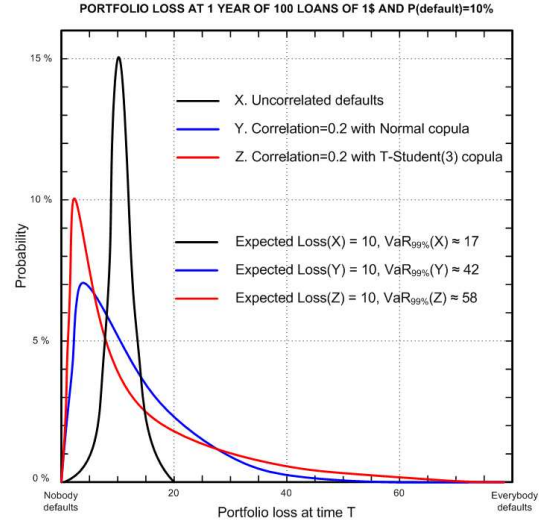


Figure A.5: Dependence structure effect

Definition A.3 — Copula. A copula function C is a multivariate distribution defined on the unit hypercube $[0, 1]^d$ with standard uniform marginals. More precisely,

$$C(u_1, \dots, u_d) = \Pr\{U_1 \leq u_1, \dots, U_d \leq u_d\}$$

in which $U_i \sim \text{Uniform}(0, 1)$ for $i = 1, \dots, d$.

Sklar's theorem [31] states that any multivariate distribution with continuous marginals can be decomposed into the marginals and a copula that reflects the structure of dependence between them. Later, we will use this statement to define and simulate the t-Student copula.

Theorem A.1 — Sklar's theorem. Let F be a d -dimensional distribution function with margins F_1, \dots, F_d . Then there is a d -copula C such that for all $x \in \mathbb{R}^d$,

$$F(x_1, \dots, x_d) = C(F_1(x_1), \dots, F_d(x_d))$$

If F_1, \dots, F_d are all continuous, the C is unique; otherwise, C is uniquely determined on $\text{Ran}F_1 \times \dots \times \text{Ran}F_d$. Conversely, if C is a d -copula and F_1, \dots, F_d are distribution functions, then the function F defined above is a d -dimensional distribution function with margins F_1, \dots, F_d .

Corollary A.1 — Copula of a multivariate distribution. Let $X = (X_1, \dots, X_d)$ be a random vector with a multivariate distribution F and continuous marginals F_1, \dots, F_d . Then its copula is

$$C(u_1, \dots, u_d) = F(F_1^{-1}(u_1), \dots, F_d^{-1}(u_d))$$

Corollary A.2 — Copula simulation. Let $X = (X_1, \dots, X_d)$ be a random vector with a multivariate distribution F and continuous marginals F_1, \dots, F_d . If we have a procedure to

simulate X , then we can simulate its copula C using:

$$(U_1, \dots, U_d) = (F_1(X_1), \dots, F_d(X_d))$$

in which U_i are the copula components.

Corollary A.3 — Multivariate distribution simulation. Let $X = (X_1, \dots, X_d)$ be a random vector with a copula C and continuous marginals F_1, \dots, F_d . If we have a procedure to simulate C , then we can simulate X using

$$(X_1, \dots, X_d) = (F_1^{-1}(U_1), \dots, F_d^{-1}(U_d))$$

in which U_i are the copula components.

Definition A.4 Spearman's rho is a measure of dependence between two variables and it is defined as the Pearson's correlation coefficient between the ranked variables.

It has the important property that its value only relies on the underlying copula C , not on marginals. It can be written as

$$\rho_s = 12 \int_0^1 \int_0^1 C(u, v) du dv - 3$$

A.3 Multivariate t-Student Distribution

The majority of the results presented below have been extracted from [22, 8].

Definition A.5 — Multivariate t-Student distribution. The d -dimensional random vector $X = (X_1, \dots, X_d)$ is said to have a (non-singular) multivariate t-Student distribution with ν degrees of freedom, mean vector $\vec{\mu}$ and positive-definite dispersion or scatter matrix Σ , denoted $t_d(\nu, \vec{\mu}, \Sigma)$, if its density is given by

$$f(\vec{x}) = \frac{\Gamma\left(\frac{\nu+d}{2}\right)}{\Gamma\left(\frac{\nu}{2}\right) \sqrt{(\pi\nu)^d |\Sigma|}} \left(1 + \frac{(\vec{x} - \vec{\mu})^\top \Sigma^{-1} (\vec{x} - \vec{\mu})}{\nu}\right)^{-\frac{\nu+d}{2}}$$

in which Γ is the gamma function and $|\Sigma|$ is the determinant of the matrix.

Proposition A.4 — t-Student covariance. The covariance matrix of the $X \sim t_d(\nu, \vec{\mu}, \Sigma)$ is

$$\text{Cov}(X) = \frac{\nu}{\nu-2} \cdot \Sigma \quad \text{if } \nu > 2$$

Proposition A.5 — Gaussian as limit of the t-Student. The t-Student distribution con-

verges to a Gaussian distribution when ν tends to ∞ .

$$\lim_{\nu \rightarrow \infty} t_d(\nu, \vec{\mu}, \Sigma) = N(\vec{\mu}, \Sigma)$$

Proposition A.6 — Multivariate t-Student characterization. A random vector $T \sim t_d(\nu, \vec{\mu}, \Sigma)$ can be expressed as:

$$T \stackrel{d}{=} \vec{\mu} + \sqrt{\frac{\nu}{V}} \cdot Z \quad \text{where } Z \sim N(\vec{0}, \Sigma) \text{ and } V \sim \chi_\nu^2$$

Proposition A.7 — Multivariate t-Student marginals. Assume $X \sim t_d(\nu, \vec{0}, \Sigma)$. Then its i -th marginal is $X_i \sim t_1(\nu, 0, \sigma_{ii})$.

Definition A.6 — t-Student copula. The t-Student copula, $C_{\nu, \Sigma}^t$, is the copula of the multivariate t-Student distribution $t_d(\nu, \vec{\mu}, \Sigma)$.

Proposition A.8 — t-Student copula equivalence. The copula of $t_d(\nu, \vec{\mu}, \Sigma)$ is identical to that of $t_d(\nu, \vec{0}, R)$ in which R is the correlation matrix implied by the dispersion matrix Σ .

Proposition A.9 — t-Student copula density. The t-Student copula, $C_{\nu, R}^t$, in which R is a correlation matrix, has the following distribution function:

$$C_{\nu, R}^t(u_1, \dots, u_d) = \int_{-\infty}^{t_\nu^{-1}(u_1)} \dots \int_{-\infty}^{t_\nu^{-1}(u_d)} f(x) \, dx$$

in which $f(x)$ is the density function of $t_d(\nu, \vec{0}, R)$ and t_ν^{-1} denotes the quantile function of the univariate distribution $t_1(\nu, 0, 1)$. The copula density is

$$c_{\nu, R}^t(u_1, \dots, u_d) = |R|^{-\frac{1}{2}} \frac{\Gamma(\frac{\nu+d}{2})}{\Gamma(\frac{\nu}{2})} \left[\frac{\Gamma(\frac{\nu}{2})}{\Gamma(\frac{\nu+1}{2})} \right]^d \frac{\left(1 + \frac{\zeta' R^{-1} \zeta}{\nu} \right)^{-\frac{\nu+d}{2}}}{\prod_{i=1}^d \left(1 + \frac{\zeta_i^2}{\nu} \right)^{-\frac{\nu+1}{2}}}$$

in which $\zeta = (t_\nu^{-1}(u_1), \dots, t_\nu^{-1}(u_d))$ is the vector of the t-Student univariate inverse distribution functions.

A.4 Covariance Block Matrix

The content presented in this appendix is detailed and extended in [33].

Definition A.7 — Covariance block matrix. A matrix A is a covariance block matrix $B_k(\vec{n}, \vec{d}, M)$ in which:

- $\vec{n} = (n_1, \dots, n_k)$ with $n_i \in \mathbb{N}$ and $1 \leq n_i$ (number of elements per block),
- $\vec{d} = (d_1, \dots, d_k)$ with $d_i \in \mathbb{R}$ (diagonal block values), and
- M is a $k \times k$ symmetric matrix with values $m_{ij} \in \mathbb{R}$ (block values)

when

- each block B_{ij} is an $n_i \times n_j$ constant matrix with value m_{ij}
- diagonal blocks B_{ii} have diagonal values d_i
- A is definite-positive.

Example A.5 — Correlation block matrix. We have 6 obligors sorted by sector. The first three belong to the banking sector, the next two belong to the energy sector and the last one belongs to the services sector. Obligor default time dependence is determined only by sectors. Thus, the correlation matrix of the underlying elliptical copula is a correlation block matrix.

$$\left. \begin{array}{l} \vec{n} = (3, 2, 1) \\ \vec{d} = (1, 1, 1) \\ M = \begin{pmatrix} 0.5 & 0.2 & 0.1 \\ 0.2 & 0.4 & 0.15 \\ 0.1 & 0.15 & 0.5 \end{pmatrix} \end{array} \right\} \rightarrow B_3(\vec{n}, \vec{d}, M) = \begin{pmatrix} 1 & 0.5 & 0.5 & 0.2 & 0.2 & 0.1 \\ 0.5 & 1 & 0.5 & 0.2 & 0.2 & 0.1 \\ 0.5 & 0.5 & 1 & 0.2 & 0.2 & 0.1 \\ \hline 0.2 & 0.2 & 0.2 & 1 & 0.4 & 0.15 \\ 0.2 & 0.2 & 0.2 & 0.4 & 1 & 0.15 \\ \hline 0.1 & 0.1 & 0.1 & 0.15 & 0.15 & 1 \end{pmatrix}$$

Proposition A.10 — Covariance block matrix eigenvalues. Let $A = B_k(\vec{n}, \vec{d}, M)$ a non-singular matrix, and let G be the $k \times k$ deflated matrix

$$G = \begin{pmatrix} d_1 + (n_1 - 1) \cdot m_{11} & n_2 \cdot m_{12} & \cdots & n_k \cdot m_{1k} \\ n_1 \cdot m_{21} & d_2 + (n_2 - 1) \cdot m_{22} & \cdots & n_k \cdot m_{2k} \\ \vdots & \vdots & \ddots & \vdots \\ n_1 \cdot m_{k1} & n_2 \cdot m_{k2} & \cdots & d_k + (n_k - 1) \cdot m_{kk} \end{pmatrix}$$

Then, the eigenvalues of A are as follows:

- $d_i - m_{ii}$ with multiplicity $n_i - 1$ for $i = 1, \dots, k$.
- λ_i , the eigenvalues of G with multiplicity 1.

Corollary A.4 A covariance block matrix $B_k(\vec{n}, \vec{d}, M)$ is definite-positive if

- $d_i > m_{ii}$ for all $i = 1, \dots, k$, and
- the deflated matrix G is definite-positive.

A.5 Multi-Factor Gaussian Distribution

We only consider the case where the variances are 1, that is $\text{Var}(X_i) = 1$.

Definition A.8 — Gaussian multi-factor distribution. We say that a multivariate distribution X follows a Gaussian multi-factor distribution if it can be expressed in this form:

$$X_i^j = w_i \cdot Z_i + \sqrt{1 - w_i^2} \cdot \varepsilon_i^j \quad \text{where} \quad \begin{cases} i = 1, \dots, k & k \text{ is the number of factors} \\ j = 1, \dots, n_i & n_i \text{ is } i\text{-th factor size} \\ Z \sim N(\vec{0}, R) & R \text{ is a } k \times k \text{ correlation matrix} \\ w_i \in (0, 1) & \text{the factor loadings} \\ \varepsilon_i^j \sim N(0, 1) \text{ iid} & \forall i, j \\ Z, \varepsilon_i^j \text{ independents} & \forall i, j \end{cases}$$

An alternative definition based on the covariance matrix is

$$X_i^j = Z_i + \sqrt{1 - \Sigma_{ii}} \cdot \varepsilon_i^j \quad \text{where} \quad \begin{cases} i = 1, \dots, k & k \text{ is the number of factors} \\ j = 1, \dots, n_i & n_i \text{ is } i\text{-th factor size} \\ Z \sim N(\vec{0}, \Sigma) & \Sigma \text{ is a } k \times k \text{ covariance matrix} \\ \Sigma_{ii} \in (0, 1) & \text{diagonal values of } \Sigma \\ \varepsilon_i^j \sim N(0, 1) \text{ iid} & \forall i, j \\ Z, \varepsilon_i^j \text{ independents} & \forall i, j \end{cases}$$

Proposition A.11 — The Gaussian multi-factor distribution is a Gaussian distribution.

The Gaussian multi-factor distribution defined by factor loadings \vec{w} , the factor correlation matrix R , and $n = \sum_{i=1}^k n_i$ is a multivariate Gaussian distribution with an $n \times n$ block correlation matrix $\hat{R} = B_k(\vec{n}, \vec{1}, \Sigma)$ with values $\Sigma_{ij} = w_i \cdot w_j \cdot R_{ij}$.

Proof. We check that the correlation between components of the Gaussian multi-factor distribution fulfill the equivalence.

Version with the correlation matrix R :

$$\begin{aligned} \text{Var}(X_i^j) &= w_i^2 \cdot \text{Var}(Z_i) + (1 - w_i^2) \cdot \text{Var}(\varepsilon_i^j) + 2 \cdot w_i \cdot \sqrt{1 - w_i^2} \cdot \text{Cov}(Z_i, \varepsilon_i^j) \\ &= w_i^2 + (1 - w_i^2) = 1 \end{aligned}$$

$$\begin{aligned} \text{Cor}(X_{i_1}^{j_1}, X_{i_2}^{j_2}) &= \text{Cov}(X_{i_1}^{j_1}, X_{i_2}^{j_2}) \\ &= w_{i_1} \cdot w_{i_2} \cdot \text{Cov}(Z_{i_1}, Z_{i_2}) + w_{i_1} \cdot \sqrt{1 - w_{i_2}^2} \cdot \text{Cov}(Z_{i_1}, \varepsilon_{i_2}^{j_2}) + \\ &\quad + \sqrt{1 - w_{i_1}^2} \cdot w_{i_2} \cdot \text{Cov}(\varepsilon_{i_1}^{j_1}, Z_{i_2}) + \sqrt{1 - w_{i_1}^2} \cdot \sqrt{1 - w_{i_2}^2} \cdot \text{Cov}(\varepsilon_{i_1}^{j_1}, \varepsilon_{i_2}^{j_2}) \\ &= w_{i_1} \cdot w_{i_2} \cdot \text{Cov}(Z_{i_1}, Z_{i_2}) \\ &= w_{i_1} \cdot w_{i_2} \cdot R_{i_1 i_2} \end{aligned}$$

Version with the covariance matrix Σ :

$$\begin{aligned} \text{Var}(X_i^j) &= \text{Var}(Z_i) + (1 - \Sigma_{ii}) \cdot \text{Var}(\varepsilon_i^j) + 2 \cdot \sqrt{1 - \Sigma_{ii}} \cdot \text{Cov}(Z_i, \varepsilon_i^j) \\ &= \Sigma_{ii} + (1 - \Sigma_{ii}) = 1 \end{aligned}$$

$$\begin{aligned} \text{Cor}(X_{i_1}^{j_1}, X_{i_2}^{j_2}) &= \text{Cov}(X_{i_1}^{j_1}, X_{i_2}^{j_2}) \\ &= \text{Cov}(Z_{i_1}, Z_{i_2}) + \sqrt{1 - \Sigma_{i_2 i_2}} \cdot \text{Cov}(Z_{i_1}, \varepsilon_{i_2}^{j_2}) + \\ &\quad + \sqrt{1 - \Sigma_{i_1 i_1}} \cdot \text{Cov}(\varepsilon_{i_1}^{j_1}, Z_{i_2}) + \sqrt{1 - \Sigma_{i_1 i_1}} \cdot \sqrt{1 - \Sigma_{i_2 i_2}} \cdot \text{Cov}(\varepsilon_{i_1}^{j_1}, \varepsilon_{i_2}^{j_2}) \\ &= \Sigma_{i_1 i_2} \end{aligned}$$

The characterization of the multivariate Gaussian [2, thm. 2.6.2] posits that if every linear combination of the components of a vector X is normally distributed, then X is normally

distributed. It is straightforward to check that the multi-factor Gaussian model fulfills this property and consequently is a multivariate Normal. \square

Proposition A.12 — Gaussian multi-factor distribution \subsetneq multivariate Gaussian block. Any Gaussian multi-factor distribution can be expressed as a multivariate Gaussian block; however, there are multivariate Gaussian blocks that cannot be expressed as a Gaussian multi-factor distribution.

Proof. Proposition A.11 states that any multi-factor distribution is a multivariate Gaussian composed of blocks. Next we see that the reciprocal is false. Let $\widehat{R} = B_k(\vec{n}, \vec{1}, \Sigma)$, the Gaussian distribution correlation matrix. We use the equivalence $\Sigma_{ij} = w_i \cdot w_j \cdot R_{ij}$ stated in Proposition A.11 to determine the cases that cannot be expressed as a multi-factor distribution.

Case 1. Factor loading with imaginary values.

$$\Sigma_{ii} = w_i \cdot w_i \cdot R_{ii} = w_i^2$$

If Σ_{ii} is negative, then w_i has an imaginary value. For example,

$$\widehat{R} = \left(\begin{array}{cc|c} 1 & -0.5 & 0 \\ -0.5 & 1 & 0 \\ \hline 0 & 0 & 1 \end{array} \right) \rightarrow R = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, w = (\sqrt{-0.5}, 1) !!$$

Case 2. Factor correlation with absolute value > 1 .

$$\Sigma_{ij} = w_i \cdot w_j \cdot R_{ij} \rightarrow R_{ij} = \frac{\Sigma_{ij}}{w_i \cdot w_j} = \frac{\Sigma_{ij}}{\sqrt{\Sigma_{ii} \cdot \Sigma_{jj}}}$$

If $\Sigma_{ij}^2 > \Sigma_{ii} \cdot \Sigma_{jj}$, then R_{ij} has an absolute value larger than 1, which is not possible because R_{ij} is a correlation value. For example,

$$\widehat{R} = \left(\begin{array}{cc|cc} 1 & 0.5 & 0.6 & 0.6 \\ 0.5 & 1 & 0.6 & 0.6 \\ \hline 0.6 & 0.6 & 1 & 0.5 \\ 0.6 & 0.6 & 0.5 & 1 \end{array} \right) \rightarrow R = \begin{pmatrix} 1 & \frac{0.6}{0.5} \\ \frac{0.6}{0.5} & 1 \end{pmatrix} !!, w = (\sqrt{0.5}, \sqrt{0.5})$$

\square

It is disturbing that multivariate block Gaussian distributions exist that cannot be expressed as a multi-factor distribution. Fortunately, the cases in which this occurs have no practical application to credit risk modeling in which the number of obligors in each block is higher. The following proposition clarifies this theme.

Proposition A.13 — Gaussian multi-factor distribution \approx multivariate Gaussian block. Assume $X \sim N(\vec{0}, \widehat{R})$ in which $\widehat{R} = B_k(\vec{n}, \vec{1}, \Sigma)$ is a correlation block matrix. Then X can be expressed as a multi-factor Gaussian model if n_i values are large enough.

Proof. We revise the cases in which the multivariate Gaussian distribution cannot be expressed as a Gaussian multi-factor distribution, and we use Proposition A.10 to ascertain that if n_i values are large enough, then that is indeed a valid Gaussian multi-factor distribution.

Case 1. Factor loading with imaginary values.

Factor loading w_i with imaginary value is caused by $\Sigma_{ii} < 0$. We use Proposition A.10 to determine the restrictions regarding Σ_{ii} . We assume a unique factor. Because \hat{R} is a covariance block matrix definite-positive, the deflated matrix has a positive determinant. Then,

$$\begin{aligned} 1 + (n_1 - 1) \cdot \Sigma_{11} &> 0 \\ \Sigma_{11} &> \frac{-1}{n_1 - 1} \\ \text{tending } n_1 &\rightarrow \infty \\ \Sigma_{11} &\geq 0 \end{aligned}$$

Case 2. Factor correlation with absolute value > 1 .

A factor correlation larger than 1 is caused by $R_{12} = \frac{\Sigma_{12}}{\sqrt{\Sigma_{11} \cdot \Sigma_{22}}}$ in which $\Sigma_{12}^2 > \Sigma_{11} \cdot \Sigma_{22}$. We use Proposition A.10 to determine the restrictions regarding Σ_{12} . Because \hat{R} is covariance block matrix definite-positive, the deflated matrix has a positive determinant. Then,

$$\begin{aligned} 1 + (n_1 - 1) \cdot \Sigma_{11} + (n_2 - 1) \cdot \Sigma_{22} + (n_1 - 1) \cdot (n_2 - 1) \cdot \Sigma_{11} \cdot \Sigma_{22} - n_1 \cdot n_2 \cdot \Sigma_{12}^2 &> 0 \\ \Sigma_{12}^2 &< \frac{1}{n_1 \cdot n_2} + \frac{(n_1 - 1)}{n_1 \cdot n_2} \cdot \Sigma_{11} + \frac{(n_2 - 1)}{n_1 \cdot n_2} \cdot \Sigma_{22} + \frac{(n_1 - 1) \cdot (n_2 - 1)}{n_1 \cdot n_2} \cdot \Sigma_{11} \cdot \Sigma_{22} \\ \text{tending } n_1 &\rightarrow \infty \text{ and } n_2 \rightarrow \infty \\ \Sigma_{12}^2 &\leq \Sigma_{11} \cdot \Sigma_{22} \end{aligned}$$

□

A.6 Metropolis-Hastings Algorithm

This section contains a brief introduction to Bayesian inference using the Metropolis-Hastings algorithm and is heavily influenced by the book *Bayesian Modeling Using WinBUGS* [27].

Bayesian statistics differ from classical statistical theory because all unknown parameters are considered as random variables. We want to calculate the distribution $f(\theta | y)$ of the parameters θ given the observed data y . According to the Bayes theorem, this can be written as

$$f(\theta | y) = \frac{f(y | \theta) \cdot f(\theta)}{f(y)} \propto f(y | \theta) \cdot f(\theta)$$

We call *posterior distribution* to $f(\theta | y)$ and *prior distribution* to $f(\theta)$. This prior distribution expresses the information available to the researcher before any data are involved in the statistical analysis. $f(y | \theta)$ is the probability of observing y given θ and is also known as the likelihood.

$$f(y | \theta) = \prod_{i=1}^n f(y_i | \theta)$$

When utilizing Bayesian inference, we can identify a closed form of the posterior distribution in only a few cases. Generally this must be determined by numerical methods. One of the most used is the Metropolis-Hastings (M-H) algorithm, which belongs to the family of Markov Chain Monte Carlo (MCMC) methods. These methods are based on the construction of a Markov chain that converges on the target distribution, which, in our case, is the posterior distribution.

Definition A.9 — Markov chain. A Markov chain is a stochastic process $\{\theta^{(1)}, \dots, \theta^{(T)}\}$

such that

$$f\left(\theta^{(t+1)} \mid \theta^{(t)}, \dots, \theta^{(1)}\right) = f\left(\theta^{(t+1)} \mid \theta^{(t)}\right),$$

that is, the distribution of θ at sequence $t + 1$, given all the preceding θ values, depends only on the previous value. When the Markov chain is irreducible, aperiodic, and positive-recurrent, the distribution of $\theta^{(t)}$ converges on its equilibrium distribution, which is independent of the initial values of the chain $\theta^{(0)}$.

The Markov chain is not an unfamiliar concept to us because the transition matrix (see Definition A.2) is a Markov chain in which its equilibrium state is the *defaulted* rating (i.e., just any obligor ends up defaulting). The Metropolis-Hastings algorithm generates a sample from $f(\theta)$ by constructing a Markov chain that has $f(\theta)$ as the equilibrium distribution.

Algorithm A.2 — Metropolis-Hastings. Let us assume a target distribution $f(x)$ from which we wish to generate a sample $\{x^{(0)}, \dots, x^{(T)}\}$. The Metropolis-Hastings algorithm can be described by the following iterative procedure:

- Set initial values $x^{(0)}$
- For $t = 1, \dots, T$ repeat the following steps:
 1. Set $x = x^{(t-1)}$
 2. Let x' be a random value from the proposal distribution $q(x \rightarrow x') = q(x' \mid x)$
 3. Calculate the acceptance rate $\alpha = \min\left(1, \frac{f(x') \cdot q(x \mid x')}{f(x) \cdot q(x' \mid x)}\right)$
 4. Update $x^{(t)} = \begin{cases} x' & \text{with probability } \alpha \\ x^{(t-1)} & \text{with probability } 1 - \alpha \end{cases}$

Because this is a Markov chain, the chosen starting point $x^{(0)}$ is irrelevant because the algorithm converges on the target distribution no matter the starting point. If the starting point has low probability in respect to the target distribution, then the initial samples do not come from the equilibrium distribution and must be discarded until there is not equilibrium. We call the first B iterations eliminated from the sample the *burn-in period* for this reason.

Bayesian inference. The Metropolis-Hastings algorithm outlined above can be directly implemented to conduct Bayesian inference by substituting x with the unknown parameters θ and the target distribution $f(x)$ with the posterior distribution $f(\theta \mid y)$. Then, using the Bayes theorem, the acceptance rate becomes

$$\alpha = \min\left(1, \frac{f(\theta' \mid y) \cdot q(\theta \mid \theta')}{f(\theta \mid y) \cdot q(\theta' \mid \theta)}\right) = \min\left(1, \frac{f(y \mid \theta') \cdot f(\theta') \cdot q(\theta \mid \theta')}{f(y \mid \theta) \cdot f(\theta) \cdot q(\theta' \mid \theta)}\right)$$

Random-walk Metropolis-Hastings is a special case with symmetrical proposals $q(\theta' \mid \theta) = q(\theta \mid \theta')$ in which $q(\theta' \mid \theta) \equiv N(\theta, \sigma)$. In this case, the acceptance rate depends only on the posterior distribution and in the Bayesian inference case becomes

$$\alpha = \min\left(1, \frac{f(y \mid \theta') \cdot f(\theta') \cdot q(\theta \mid \theta')}{f(y \mid \theta) \cdot f(\theta) \cdot q(\theta' \mid \theta)}\right) = \min\left(1, \frac{f(y \mid \theta') \cdot f(\theta')}{f(y \mid \theta) \cdot f(\theta)}\right)$$

The random walk fixes the shape of the proposal distribution and requires specifying the value σ . This value has a direct non-linear effect on the acceptance ratio (#accepted/#iterations). There is an optimal acceptance ratio that minimizes the auto-correlation and in most cases relies

on the interval $[0.234, 0.44]$. To avoid the manual tuning of the scale parameter σ , we use the Robbins-Monro algorithm [15], which performs an adaptive scaling to achieve the desired acceptance ratio.

Proposition A.14 — The Robbins-Monro process. To achieve an acceptance ratio p in the random-walk Metropolis-Hastings algorithm, consider the following adaptive procedure:

$$\sigma^{(t+1)} = \begin{cases} \sigma^{(t)} \cdot \left(1 + \frac{1}{p \cdot t}\right) & \text{if } \theta' \text{ is accepted} \\ \sigma^{(t)} \cdot \left(1 - \frac{1}{(1-p) \cdot t}\right) & \text{if } \theta' \text{ is rejected} \end{cases}$$

in which the starting value $\sigma^{(1)} > 0$ has an arbitrary value.

Metropolis-Hastings using logarithms. The product of functions involves complicated expressions and accuracy problems. Therefore, it is usual to operate with logarithms. The logarithmic version of the acceptance rate in the Bayesian inference case becomes

$$\begin{aligned} \ln(\alpha) &= \min \left(\ln(1), \ln \left(\frac{f(y|\theta') \cdot f(\theta)}{f(y|\theta) \cdot f(\theta')} \right) \right) = \\ &= \min(0, \ln f(y|\theta') + \ln f(\theta) - \ln f(y|\theta) - \ln f(\theta')) \end{aligned}$$

The single-component Metropolis-Hastings algorithm consists of updates of sequentially univariate components using Metropolis-Hastings steps. One generated observation $\theta^{(t)}$ is obtained after updating all components of the parameter vector. Generally, the sequence of updating the elements of θ does not influence the convergence of the algorithm. Nevertheless, to ensure randomness, a random selection of the updating sequence is recommended. We use the notation $\theta_{\setminus j}$ to indicate the vector θ , excluding its j -th component [$\theta_{\setminus j} = (\theta_1, \dots, \theta_{j-1}, \theta_{j+1}, \dots, \theta_d)$].

Algorithm A.3 — Bayesian inference using Metropolis-Hastings. Let $f(y|\theta)$ be the likelihood function, with priors $f_j(\theta_j)$ and desired acceptance ratios p_j . Then the parameters distribution can be sampled using the following Metropolis-Hastings algorithm:

- Set initial values $\theta^{(0)}$
- Set initial values $\sigma^{(0)}$
- For $t = 1, \dots, T$ repeat the following steps:
 1. Set $\theta = \theta^{(t-1)}$
 2. Shuffle the update sequence
 3. For $j = 1, \dots, d$ repeat the following steps:
 - a. $\theta'_j = N(\theta_j, \sigma_j)$
 - b. Calculate the logarithm of the acceptance rate
 $\ln(\alpha) = \ln f(y|\theta'_j, \theta_{\setminus j}) + \ln f_j(\theta'_j; \theta_{\setminus j}) - \ln f(y|\theta_j, \theta_{\setminus j}) - \ln f_j(\theta_j; \theta_{\setminus j})$
 - c. Simulate $u \sim U(0, 1)$
 - d. Update $\theta_j = \begin{cases} \theta'_j & \text{if } \ln(u) < \ln(\alpha) \\ \theta_j & \text{otherwise} \end{cases}$
 - e. Update $\sigma_j^{(t)} = \begin{cases} \sigma_j^{(t-1)} \cdot \left(1 + \frac{1}{p_j \cdot t}\right) & \text{if } \ln(u) < \ln(\alpha) \\ \sigma_j^{(t-1)} \cdot \left(1 - \frac{1}{(1-p_j) \cdot t}\right) & \text{otherwise} \end{cases}$
 4. Set $\theta^{(t)} = \theta$

The simulated values from the posterior distribution have ergodic properties; however, they are

not independent. Independent values can be obtained by plotting the autocorrelation function (ACF) and estimating the length, B , for which the autocorrelation is canceled. Then only the values whose index is a multiple of B are considered. Notably, the estimated parameters are not independent of one another; in fact, they form a multivariate distribution.

Example A.6 — Beta distribution calibration. We have a sample $\{y_1, \dots, y_{100}\}$ coming from a $\text{Beta}(2, 5)$ distribution. We want to perform Bayesian inference to determine the value of the parameters α and β using the Metropolis-Hastings algorithm presented above. The density of the Beta distribution is

$$f(x; \alpha, \beta) = \frac{\Gamma(\alpha + \beta)}{\Gamma(\alpha) \cdot \Gamma(\beta)} \cdot x^{(\alpha-1)} \cdot (1-x)^{(\beta-1)}$$

The log-likelihood function, then, is

$$\ln f(y | \alpha, \beta) = \sum_{i=1}^{100} \ln \Gamma(\alpha + \beta) - \ln \Gamma(\alpha) - \ln \Gamma(\beta) + (\alpha - 1) \cdot \ln(y_i) + (\beta - 1) \cdot \ln(1 - y_i)$$

We know that $\alpha \in (0, 10)$ and $\beta \in (0, 30)$. Therefore, we consider that prior distributions $f_\alpha()$ and $f_\beta()$ are uniform distributions. Then,

$$\begin{aligned} \ln f_\alpha(\alpha; \beta) &= \ln\left(\frac{1}{10}\right) \\ \ln f_\beta(\beta; \alpha) &= \ln\left(\frac{1}{30}\right) \end{aligned}$$

We apply the Algorithm A.3 with starting point $(\alpha^{(0)}, \beta^{(0)}) = (1, 1)$ and desired acceptance ratios $(p_\alpha, p_\beta) = (0.3, 0.3)$. Figure A.6 displays the results after 25,000 iterations. Notably, the first 500 iterations are not valid because the Markov chain has not achieved the equilibrium distribution. These must be considered in the burn-in period. The histograms of the sampled values approximate the parameters distribution that we are inferring.

A.7 Antithetic variates

The Monte Carlo method consists of sampling a distribution and averaging these values to approximate the mean of this distribution. The main drawback of this method is the low rate of convergence, on the order $1/\sqrt{N}$. Variance reduction methods are procedures used to increase the precision of the estimates of the Monte Carlo method. The main procedures are common random numbers, antithetic variates, control variates, importance sampling and stratified sampling. This section outlines contains the antithetic variates method.

We estimate $E(g(X))$ using the Monte Carlo method. Assume Y_1 and Y_2 as random variables with the identical distribution X . This indicates that

$$E(g(X)) = E(g(Y_1)) = E(g(Y_2)) = E\left(\frac{g(Y_1) + g(Y_2)}{2}\right)$$

and

$$\text{Var}\left(\frac{g(Y_1) + g(Y_2)}{2}\right) = \frac{\text{Var}(g(Y_1)) + \text{Var}(g(Y_2)) + 2 \cdot \text{Cov}(g(Y_1), g(Y_2))}{4} = \frac{\text{Var}(g(X)) + \text{Cov}(g(Y_1), g(Y_2))}{2}$$

If $g(Y_1)$ and $g(Y_2)$ are negatively correlated, then the variance of N simulated values from $Y_1 + Y_2$ is smaller than the variance of $2 \times N$ simulated values from X .

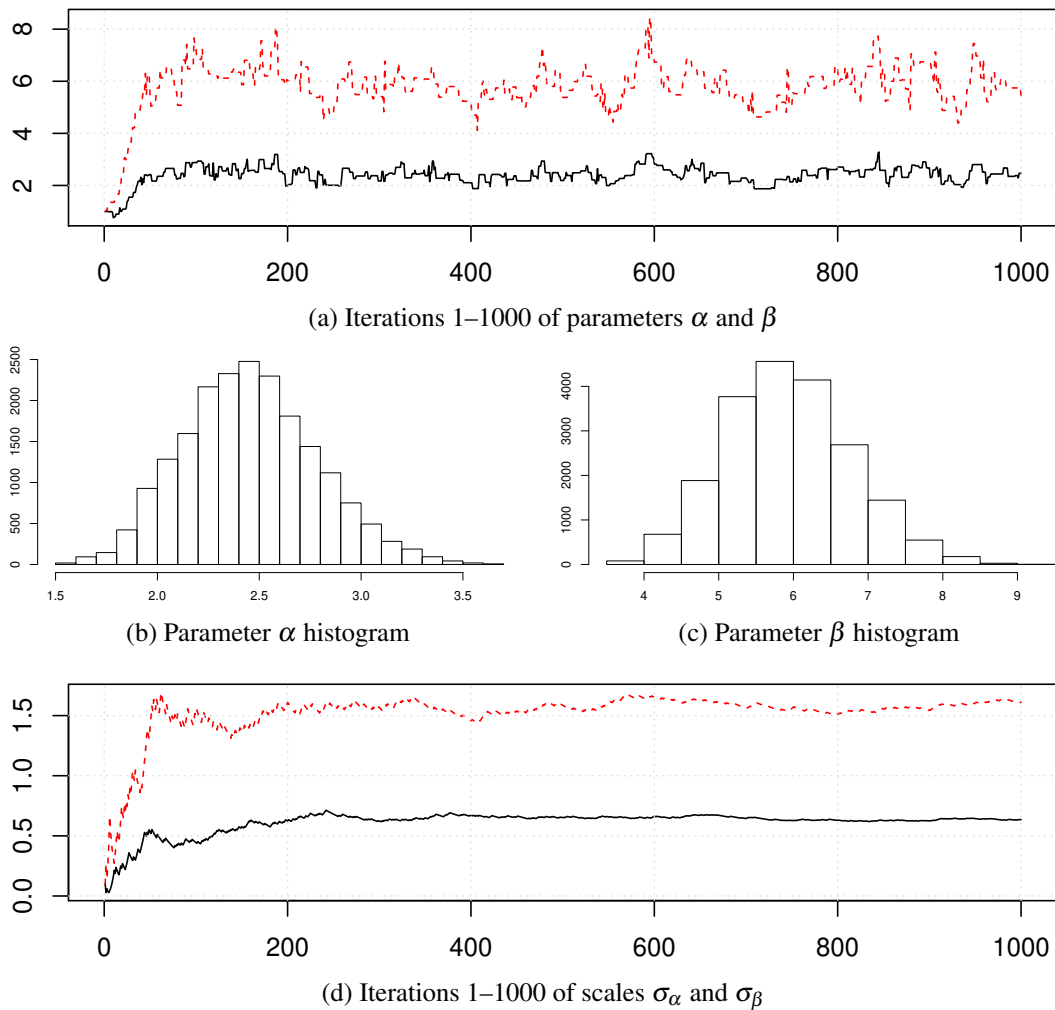


Figure A.6: Metropolis-Hastings algorithm example

Two classic cases are :

$$\begin{aligned} X \sim \text{Uniform}[0, 1] &\implies Y_1 \sim \text{Uniform}[0, 1], & Y_2 \sim 1 - \text{Uniform}[0, 1] \\ X \sim N(0, 1) &\implies Y_1 \sim N(0, 1), & Y_2 \sim -N(0, 1) \end{aligned}$$

Example A.7 In this example extracted from Wikipedia, we show how the antithetical method works and how it is used by CCruncher. We would like to estimate the following integral using the Monte Carlo method:

$$I = \int_0^1 \frac{1}{1+x} dx$$

In this case the random variable is $U \sim \text{Uniform}[0, 1]$, and its variate is $1 - U$. Listing A.1 computes the integral value using plain Monte Carlo, Monte Carlo with antithetic method, and Monte Carlo with antithetic variates but only to reduce the number of simulated values as CCruncher does.

	Sims	N	Mean	Variance
Exact = $\ln(2)$	-	-	0.693147	-
Monte Carlo	10000	10000	0.695685	0.019624
Antithetic	5000	5000	0.693404	0.000593
CCruncher	5000	10000	0.693099	0.019508

Listing A.1: Antithetic example (R script)

```
f <- function(x) { 1/(1+x) }

result = matrix(nrow=4, ncol=4)
colnames(result) = c("Sims", "N", "Mean", "Variance")
rownames(result) = 1:4

rownames(result)[1] = "Exact";
result[1,1] = 0
result[1,2] = 0
result[1,3] = integrate(f, lower=0, upper=1)$value
result[1,4] = 0

u = runif(10000)
rownames(result)[2] = "Monte Carlo";
result[2,1] = length(u)
result[2,2] = length(u)
result[2,3] = mean(f(u))
result[2,4] = var(f(u))

u = runif(5000)
rownames(result)[3] = "Antithetic";
result[3,1] = length(u)
result[3,2] = length(u)
result[3,3] = mean((f(u)+f(1-u))/2)
result[3,4] = var((f(u)+f(1-u))/2)

u = runif(5000)
u = c(u, 1-u)
rownames(result)[4] = "CCruncher";
result[4,1] = length(u)/2
result[4,2] = length(u)
result[4,3] = mean(f(u))
result[4,4] = var(f(u))

result
```


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