# Uncertainty in Risk Modelling 

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# Uncertainty in Risk Modelling 

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#### Abstract

In this thesis, we explore the uncertainty issues in risk modelling arising from the different approaches proposed in the literature and currently being used in the industry.

The first type of methods that we discuss assume that the returns of the stocks follows a generalized hyperbolic distribution. Data is calibrated by the Expectation-Maximization (EM) algorithm in order to estimate the parameters in the underlying distribution. Once we have the parameters, we estimate the Value at Risk (VaR) and Expected Shortfall (ES) by using Monte Carlo simulations.

Furthermore, we calibrate data to different copulas, including the Gauss Copula, the t Copula and the Gumbel Copula for estimation of VaR and ES using these copula structures.

The results from both methods are then compared. It can be concluded that uncertainty issues in risk modelling are very significant and can be troublesome as the values of the same risk measure computed using different methods demonstrate great oscillations.


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Harjas Singh

## Dedication

In every age
There exist quiet heroes
Who, in ther selfless devotion,
lay aside their own needs and comforts
and consecrate their energies and talents
to healing the wounds of a troubled humanity.
To these rare and often unhonored souls
this thesis is humbly dedicated.

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## Chapter 1

## Introduction

According to the Cambridge dictionary, risk is "the possibility of something bad happening." In financial terms, risk is defined as "the potential for financial loss and uncertainty about its extent" [15]. The crucial part for companies and regulators is precisely the extent of risk. Let us first begin with some definitions in order to quantify this risk.

### 1.1 Risk Measure

### 1.1.1 Basic Properties

Mathematically speaking, a risk measure is defined as a mapping from a set of random variables to the real numbers. In practice, the random variables are generally representative of portfolio returns. We represent a risk measure associated with a random variable $X$ as $\rho(X)$. A risk measure $\rho: \mathcal{L} \rightarrow \mathbb{R} \cup\{+\infty\}$ has the following properties each with its own financial interpretation [3]:

- Normalized, i.e.

$$
\rho(0)=0 .
$$

From a financial perspective, normalization is equivalent to saying that a portfolio
with no holding has zero risk.

- Translation invariance, i.e.

$$
\text { If } a \in \mathbb{R} \text { and } X \in \mathcal{L}, \text { then } \rho(X+a)=\rho(X)-a
$$

From a financial perspective, translation invariance says that adding a fixed amount of capital to the portfolio will decrease the risk of the portfolio by the same amount.

- Monotonicity, i.e.

$$
\text { If } X_{1}, X_{2} \in \mathcal{L} \text { and } X_{1} \leq X_{2} \text {, then } \rho\left(X_{2}\right) \leq \rho\left(X_{1}\right)
$$

Viewed from the lens of finance, monotonicity has the interpretation that if a portfolio performs better than another one almost surely, then the first portfolio has less risk.

- Law invariance under $\mathbb{P}$, i.e.

If $X_{1}, X_{2} \in \mathcal{L}$ have the same distribution with respect to P , then $\rho\left(X_{1}\right)=\rho\left(X_{2}\right)$.

Law-invariant risk measures allot the same level of risk to financial portfolios that have an identical distribution with respect to $\mathbb{P}$ a priori. In reality, it is the most pervasive type of risk measures that are used. Principally, any risk measure whose computation relies on statistical methods and thus depends on distributions of random variables must be law-invariant.

- Sub-additivity, i.e.

$$
\text { If } Z_{1}, Z_{2} \in \mathcal{L} \text {, then } \rho\left(Z_{1}+Z_{2}\right) \leq \rho\left(Z_{1}\right)+\rho\left(Z_{2}\right)
$$

This alludes to the principle commonly known as the diversification effect in finance, i.e. a portfolio that are well diversified carry lower risk than their counterparts. The idea is that the bad events for different assets in the portfolio may not happen simultaneously.

- Positive homogeneity, i.e.

$$
\text { If } \alpha \geq 0 \text { and } Z \in \mathcal{L}, \text { then } \rho(\alpha Z)=\alpha \rho(Z) .
$$

This implies that the risk scaling is linear, i.e. doubling your financial position doubles risk.

- Convexity, i.e.

$$
\text { If } Z_{1}, Z_{2} \in \mathcal{L} \text { and } \lambda \in[0,1] \text { then } \rho\left(\lambda Z_{1}+(1-\lambda) Z_{2}\right) \leq \lambda \rho\left(Z_{1}\right)+(1-\lambda) \rho\left(Z_{2}\right)
$$

Convexity incorporates both sub-additivity and positive homogeneity and states that a linear combination (with non-negative weights) of two financial positions has less risk than the corresponding linear combination of risks associated with the financial positions.

Not all properties are satisfied by all risk measures. In particular, a risk measure $\rho$ that satisfies monotonicity, sub-additivity, homogeneity, and translation invariance is called a coherent risk measure [3].

### 1.1.2 Examples of Risk Measures

A commonly used example of a risk measure is Value-at-Risk or VaR. Widely believed to be developed by JP Morgan during the late 1980's [2], VaR estimates the loss of a set of investments, given a certain probability and time interval. Mathematically speaking,


Figure 1.1: Graph highlighting Value-at-Risk.
if $\alpha \in(0,1)$ represents the level of confidence that we seek, then

$$
\begin{align*}
\operatorname{VaR}_{\alpha}(X) & =-\inf \left\{x \in \mathbb{R}: F_{X}(x)>\alpha\right\}  \tag{1.1}\\
& =F_{-X}^{\leftarrow}(1-\alpha)  \tag{1.2}\\
& =\inf \{m \in \mathbb{R}: \mathbb{P}(X+m \leq 0) \leq \alpha\} \tag{1.3}
\end{align*}
$$

where $F_{X}$ is the cumulative distribution function (cdf) associated with the random variable X and $F_{-X}^{\leftarrow}$ is the generalized inverse of the cumulative distribution function associated with -X. Hence, $\operatorname{VaR}_{\alpha}(X)$ is the minimal amount of cash that needs to be raised and added to the position $X$ to make it "acceptable".

VaR has remained the industry standard for Basel regulators till recently, but it is

2. Alternative situation to figure 1:VAR is the same, but the potential risk is higher


Figure 1.2: Graph highlighting shortcomings of VaR (does not capture tail-risk).
gradually being phased out in favour of Expected Shortfall (ES), which is another risk measure. Over the next few paragraphs, we discuss shortcomings of VaR and how ES elegantly takes care of those limitations.

We begin by observing that VaR is not a coherent risk measure. Indeed, it does not satisfy the sub-additivity property in some scenarios. In addition to this, it does not capture tail-risk, i.e. even if the distribution is heavy-tailed with a lot of large losses beyond the quantile level, the VaR does not change.

To account for these shortcomings, Expected Shortfall was introduced. Expected shortfall at $\alpha \%$ level is the expected return on the portfolio in the worst $\alpha \%$ of cases. Mathematically speaking, if $X$ is the payoff of a portfolio at some future time and $0<\alpha<1$ then we define the expected shortfall as

$$
\begin{equation*}
E S_{\alpha}(X)=-\frac{1}{\alpha} \int_{0}^{\alpha} \operatorname{VaR}_{\gamma}(X) \mathrm{d} \gamma \tag{1.4}
\end{equation*}
$$

Finally, we define the entropic risk measure, which uses the exponential utility func-


Figure 1.3: Graph highlighting Expected Shortfall.
tion in order to quantify risk. The entropic risk measure with the risk aversion parameter $\beta>0$ is defined as

$$
\begin{equation*}
\rho^{e n t}(X)=\frac{1}{\beta} \log \left(\mathbb{E}\left[e^{-\beta X}\right]\right) . \tag{1.5}
\end{equation*}
$$

Different from Expected Shortfall, the entropic risk measure is convex but not coherent.
We refer to [6] to a more comprehensive treatment of risk measures.

### 1.2 Calculating Risk Measures for Portfolios

There are several approaches to calculate risk measures. We highlight a couple of them in this section. In addition we discuss the data set that we will use to model returns and
subsequently estimate corresponding risk of the portfolio.

### 1.2.1 Methods to calculate Risk Measures

Of the few approaches that can be used to estimate risk measures, two are prominent in existing literature and practice. The first one is a direct approach, involving knowledge of the underlying joint distribution. The other one is an indirect approach, and models dependence using copulas. We discuss them one by one.

## Direct approach: Modelling Joint Distribution

Generally speaking, the marginal distributions (probability density functions) of the constituents of the portfolio are known (or can be safely assumed) a priori and can be calibrated using well known univariate techniques. A simple example is the return distribution underlying the Black-Scholes option pricing model. In the Black-Scholes framework, an asset is assumed to have log returns that are normally distributed [4], thereby allowing pricing of options on the asset. In such a framework, once the return distribution is established beforehand, it is simply a question of calibrating the empirical data set for the unknown parameters, i.e. the mean and the variance. Following this, we can also check if our assumptions about the distribution were valid using, for example, tests of normality [8]. Finally, once we are convinced that the return distribution is indeed as expected, we may plug in the calibrated parameters to obtain the required answer which, in this case, is the option price.

On the other hand, the joint distribution of the portfolio may be sufficiently complex to preclude such an exercise. However, under certain assumptions, we can calibrate the joint distribution of the portfolio and subsequently estimate risk measures. This is the approach that we follow in Chapter 2, where we assume that the joint distribution is a Generalized Hyperbolic Distribution and is subsequently calibrated using the Expecta-
tion Maximization (EM) algorithm.

Once we know the underlying joint distribution, and have calibrated it, we may then begin calculating risk measures. For this, we simulate a return from the distribution $n$ times, where n is a sufficiently large number (in practice, $n>1,000,000$ ). These numbers are then stored in a vector, following which Value-at-Risk and Expected Shortfall are calculated using quantiles [7].

## Indirect approach: Modelling Dependence via Copulas

Sometimes, it may be difficult to directly model the joint distribution of the portfolio. In this case, we rely on an implicit measure of dependence, i.e. the copula associated with that portfolio. Copulas capture the dependence structure of the joint distribution, and using this information we can evaluate risk measures. This is explained in greater detail in Chapter 3.

### 1.2.2 Data

The data that we use to perform analyses in this thesis consists of weekly returns of four stocks that trade on NASDAQ. These four stocks are Apple Inc. (ticker: AAPL), Ford Motor Company (ticker: F), Suncor Energy Inc. (ticker: SU) and Bank of America (ticker: BAC). The reason these particular stocks were chosen is that they represent four different sectors of the stock market, and thus their returns should, in general, be different over time. A portfolio consisting of these stocks will also be fairly representative of the entire stock market since the sectors are different. The reason weekly returns are chosen is to avoid the issue of volatility clustering that would be present in noisy daily return data.

We use the last 20 years of data (from May 1998 to May 2018), in order to have a sufficient number of weeks to conduct our analyses (1045 data points per stock). Instead
of prices, we employ log returns of closing prices to quantify movements allowing us to compare stocks at different price levels. Adjusted prices are used to account for stock splits and dividends.

We also assume an equal investment of $\$ 10,000$ in each of the stocks at inception and track the movement of the portfolio with this as the reference point.

## Chapter 2

## Risk modelling via joint distribution

Generalized hyperbolic distributions are a large class of normal mixture distributions, containing as special and limiting cases many well-known classes of distributions, such as Student's $t$-distributions, variance-gamma distributions, normal-inverse Gaussian distributions, and hyperbolic distributions. More importantly, compared to normal distributions, generalized hyperbolic distributions generally have semi-heavy tails and can accommodate skewness. Thus it has been very appealing to use them to model various real-life phenomena. Their use in financial modelling started in the 1990s. See, for example, [11, 16]. As for the case of finite Gaussian mixture models, it is conventional to use the Expectation-Maximization Algorithm or its variants to estimate the parameters in generalized hyperbolic distributions. See, for example, [9, 15]. In this chapter, we fit a generalized hyperbolic distribution to return data of four stocks to model their joint distribution. After estimating the parameters using our modified MCECM algorithm, we use the Monte Carlo methods to evaluate various risk measures of a portfolio of these four stocks.

### 2.1 Algorithmic Setup

### 2.1.1 Generalized Hyperbolic Distributions

Earlier risk models were often based on multivariate normal distributions. In recent decades, more advanced distributions are demanded, among which normal mixture models have become increasingly popular. Pioneering mixture models introduce mixture to variance only, resulting in, e.g., a already very large classes of spherical and elliptical distributions. More generally, one can introduce mixtures to both mean and variance, which produces in particular generalized hyperbolic distributions.

We refer to [15] for basic properties of generalized hyperbolic distributions. A ddimensional random vector $X$ is said to follow a generalized hyperbolic distribution if

$$
\begin{equation*}
X \stackrel{\text { def }}{=}(\mu+W \gamma)+\sqrt{W} \Sigma^{\frac{1}{2}} Z \tag{2.1}
\end{equation*}
$$

where $\mu, \gamma \in \mathbb{R}^{d}, \Sigma$ is a positive-definite $d \times d$ matrix, $Z \sim \mathrm{~N}_{d}\left(0, I_{d}\right)$ follows the standard $d$-dimensional multivariate normal distribution, and

$$
W \sim \mathrm{~N}^{-}(\lambda, \chi, \psi)
$$

follows a generalized inverse Gaussian (GIG) distribution and is independent of $Z$. Following the convention, we write

$$
X \sim \mathrm{GH}_{d}(\lambda, \chi, \psi, \mu, \Sigma, \gamma)
$$

To be specific, $W$ has the following density:

$$
\begin{equation*}
f_{W}(w)=\frac{\chi^{-\frac{\lambda}{2}} \psi^{\frac{\lambda}{2}}}{2 K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)} w^{\lambda-1} \exp \left(-\frac{\chi}{2} \frac{1}{w}-\frac{\psi}{2} w\right), \quad w>0 \tag{2.2}
\end{equation*}
$$

where $K_{\lambda}$ is a modified Bessel function of the second kind with index $\lambda$. In some literature, $K_{\lambda}$ is also called modified Bessel function of the third kind. It can be explicitly written for example as follows. For any $\lambda \in \mathbb{R}$ and $x>0$,

$$
\begin{align*}
K_{\lambda}(x) & =\frac{1}{2}\left(\frac{x}{2}\right)^{\lambda} \int_{0}^{\infty} \exp \left(-t-\frac{x^{2}}{4 t}\right) \frac{\mathrm{d} t}{t^{\lambda+1}} \\
& =\frac{1}{2} \int_{0}^{\infty} t^{\lambda-1} \exp \left(-\frac{x}{2}\left(t+t^{-1}\right)\right) \mathrm{d} t . \tag{2.3}
\end{align*}
$$

We refer to [1, 12] for basic facts on Bessel functions.
The appearance of $K_{\lambda}$ in $f_{W}$ is due to normalization: $\int_{0}^{\infty} f_{W}(w) \mathrm{d} w=1$. In other words, it holds that

$$
\begin{equation*}
\int_{0}^{\infty} w^{\lambda-1} \exp \left(-\frac{\chi}{2} \frac{1}{w}-\frac{\psi}{2} w\right) \mathrm{d} w=2\left(\frac{\chi}{\psi}\right)^{\frac{\lambda}{2}} K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right) \tag{2.4}
\end{equation*}
$$

which follows easily from (2.3) by a change of variable. Thus, by a simple computation using (2.2) and (2.4), one easily sees that

$$
\begin{align*}
\mathbb{E}\left[W^{\alpha}\right] & =\frac{\chi^{-\frac{\lambda}{2}} \psi^{\frac{\lambda}{2}}}{2 K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)} \int_{0}^{\infty} w^{\alpha+\lambda-1} \exp \left(-\frac{\chi}{2} \frac{1}{w}-\frac{\psi}{2} w\right) \mathrm{d} w  \tag{2.5}\\
& =\left(\frac{\chi}{\psi}\right)^{\frac{\alpha}{2}} \frac{K_{\lambda+\alpha}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)}{K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)} \tag{2.6}
\end{align*}
$$

By the definition of $X$, it is clear that $X \mid W$ is normally distributed with the following density:

$$
\begin{align*}
& f_{X \mid W}(x \mid w) \\
= & \frac{1}{(2 \pi)^{\frac{d}{2}} \operatorname{det}(w \Sigma)^{\frac{1}{2}}} \exp \left(-\frac{(x-\mu-w \gamma)^{\prime}(w \Sigma)^{-1}(x-\mu-w \gamma)}{2}\right)  \tag{2.7}\\
= & \frac{1}{(2 \pi)^{\frac{d}{2}} \operatorname{det}(\Sigma)^{\frac{1}{2}} w^{\frac{d}{2}}} \exp \left(-\frac{(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)}{2} \frac{1}{w}-\frac{\gamma^{\prime} \Sigma^{-1} \gamma}{2} w+(x-\mu)^{\prime} \Sigma^{-1} \gamma\right) .
\end{align*}
$$

Therefore, by (2.4), $X$ has the following density:

$$
\begin{align*}
f_{X}(x)= & \int_{0}^{\infty} f_{X \mid W}(x \mid w) f_{W}(w) \mathrm{d} w \\
= & \frac{\exp \left((x-\mu)^{\prime} \Sigma^{-1} \gamma\right)}{(2 \pi)^{\frac{d}{2}} \operatorname{det}(\Sigma)^{\frac{1}{2}}} \cdot \frac{\chi^{-\frac{\lambda}{2}} \psi^{\frac{\lambda}{2}}}{2 K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)}  \tag{2.8}\\
& \quad \times \frac{2 K_{\lambda-\frac{d}{2}}\left(\left(\chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)\right)^{\frac{1}{2}}\left(\psi+\gamma^{\prime} \Sigma^{-1} \gamma\right)^{\frac{1}{2}}\right)}{\left(\chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)\right)^{-\frac{\lambda-\frac{d}{2}}{2}}\left(\psi+\gamma^{\prime} \Sigma^{-1} \gamma\right)^{\frac{\lambda-\frac{d}{2}}{2}}}
\end{align*}
$$

Consequently, $W \mid X$ has the following density:

$$
\begin{align*}
& f_{W \mid X}(w \mid x)=\frac{f_{X \mid W}(x \mid w) f_{W}(w)}{f_{X}(x)} \\
= & \frac{\left(\chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)\right)^{-\frac{\lambda-\frac{d}{2}}{2}}\left(\psi+\gamma^{\prime} \Sigma^{-1} \gamma\right)^{\frac{\lambda-\frac{d}{2}}{2}}}{2 K_{\lambda-\frac{d}{2}}\left(\left(\chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)\right)^{\frac{1}{2}}\left(\psi+\gamma^{\prime} \Sigma^{-1} \gamma\right)^{\frac{1}{2}}\right)}  \tag{2.9}\\
& \quad \times w^{\lambda-\frac{d}{2}-1} \exp \left(-\frac{\chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu)}{2} \frac{1}{w}-\frac{\psi+\gamma^{\prime} \Sigma^{-1} \gamma}{2} w\right),
\end{align*}
$$

and therefore,

$$
\begin{equation*}
\left.W \left\lvert\, X=x \sim \mathrm{~N}^{-}\left(\lambda-\frac{d}{2}, \chi+(x-\mu)^{\prime} \Sigma^{-1}(x-\mu), \psi+\gamma^{\prime} \Sigma^{-1} \gamma\right)\right.\right) \tag{2.10}
\end{equation*}
$$

### 2.1.2 The Expectation-Maximization Algorithm

A comprehensive treatment in the EM algorithm can be found in [9, 14]. We now use it to fit a multivariate generalized hyperbolic distribution $\operatorname{GH}_{d}(\lambda, \chi, \psi, \mu, \Sigma, \gamma)$ to $n$ observations

$$
\mathbf{x}=\left(x_{1}, \ldots, x_{n}\right)
$$

that are independently produced from an iid experiment. For this purpose, let $X_{i}$ be the random outcome at the $i$-th observation, and write

$$
\mathbf{X}=\left(X_{1}, \ldots, X_{n}\right) .
$$

Let

$$
\mathbf{W}=\left(W_{1}, \ldots, W_{n}\right)
$$

be the corresponding mixtures. Then

$$
X_{i} \sim \operatorname{GH}_{d}(\lambda, \chi, \psi, \mu, \Sigma, \gamma), \quad W_{i} \sim \mathrm{~N}^{-}(\lambda, \chi, \psi), \quad i=1, \ldots, n .
$$

Moreover, $\left(X_{i}, W_{i}\right), i=1, \ldots, n$, are independent random pairs.
In the EM algorithm, we regard $(\mathbf{X}, \mathbf{W})$ as the complete information, $\mathbf{X}$ as the observable information, and $\mathbf{W}$ as the missing information. Put $\theta=(\lambda, \chi, \psi, \mu, \Sigma, \gamma)$, the parameter collection in the model. Given an estimate $\theta^{(m)}$ of $\theta$ at the $m$-th step, the estimate of $\theta$ at the $(m+1)$-th step is given by the argument in maximizing the following $Q$-function:

$$
\begin{align*}
& Q\left(\theta \mid \theta^{(m)}\right) \\
= & \mathbb{E}_{\mathbf{X}, \mathbf{W} \mid \mathbf{x}, \theta^{(m)}}\left[\ln f_{\mathbf{X}, \mathbf{W}}(\mathbf{X}, \mathbf{W} \mid \theta)\right] \\
= & \sum_{i=1}^{n} \mathbb{E}_{X_{i}, W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{X_{i}, W_{i}}\left(X_{i}, W_{i} \mid \theta\right)\right]  \tag{2.11}\\
= & \sum_{i=1}^{n} \mathbb{E}_{X_{i}, W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{X_{i} \mid W_{i}}\left(X_{i} \mid W_{i}, \mu, \Sigma, \gamma\right)\right]+\sum_{i=1}^{n} \mathbb{E}_{X_{i}, W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{W_{i}}\left(W_{i} \mid \lambda, \chi, \psi\right)\right] \\
= & \sum_{i=1}^{n} \mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{X_{i} \mid W_{i}}\left(x_{i} \mid W_{i}, \mu, \Sigma, \gamma\right)\right]+\sum_{i=1}^{n} \mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{W_{i}}\left(W_{i} \mid \lambda, \chi, \psi\right)\right] \\
:= & Q_{1}\left(\mu, \Sigma, \gamma \mid \theta^{(m)}\right)+Q_{2}\left(\lambda, \chi, \psi \mid \theta^{(m)}\right),
\end{align*}
$$

where the second equality is due to independence of the pairs $\left(X_{i}, W_{i}\right)$ 's and [9, Proposition 1.1].

The E-step. Here $E$ stands for expectation. This step deals with computing $Q$.
Recall from (2.10) that

$$
W_{i} \mid x_{i}, \theta^{(m)} \sim \mathrm{N}^{-}\left(\lambda^{(m)}-\frac{d}{2}, \chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right), \psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right) .
$$

Thus by (2.6), we have the following

$$
\begin{aligned}
a_{i}^{(m)} & :=\mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[W_{i}\right] \\
& =\left(\frac{\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)}{\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}}\right)^{\frac{1}{2}}
\end{aligned}
$$

$$
\times \frac{K_{\lambda^{(m)}-\frac{d}{2}+1}\left(\sqrt{\left(\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)\right)\left(\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)}\right)}{K_{\lambda(m)-\frac{d}{2}}\left(\sqrt{\left(\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)\right)\left(\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)}\right)}
$$

$$
b_{i}^{(m)}:=\mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[W_{i}^{-1}\right]
$$

$$
=\left(\frac{\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)}{\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}}\right)^{-\frac{1}{2}}
$$

$$
\times \frac{K_{\lambda^{(m)}-\frac{d}{2}-1}\left(\sqrt{\left(\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)\right)\left(\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)}\right)}{K_{\lambda^{(m)}-\frac{d}{2}}\left(\sqrt{\left(\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)\right)\left(\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)}\right)}
$$

Also, put

$$
c_{i}^{(m)}=\mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln W_{i}\right] .
$$

The computation of $c_{i}^{(m)}$ is more sophisticated, and we will deal with it in the next
section. By (2.7), we have

$$
\begin{aligned}
& Q_{1}\left(\mu, \Sigma, \gamma \mid \theta^{(m)}\right) \\
= & \sum_{i=1}^{n} \mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{X_{i} \mid W_{i}}\left(x_{i} \mid W_{i}, \mu, \Sigma, \gamma\right)\right] \\
= & \sum_{i=1}^{n}\left(x_{i}-\mu\right)^{\prime} \Sigma^{-1} \gamma-\frac{\gamma^{\prime} \Sigma^{-1} \gamma}{2} \sum_{i=1}^{n} a_{i}^{(m)}-\sum_{i=1}^{n} \frac{\left(x_{i}-\mu\right)^{\prime} \Sigma^{-1}\left(x_{i}-\mu\right)}{2} b_{i}^{(m)}-\frac{n}{2} \ln \operatorname{det}(\Sigma)+C,
\end{aligned}
$$

where $C$ is a constant independent of $\mu, \Sigma, \gamma$.
By (2.2), we have

$$
\begin{aligned}
& Q_{2}\left(\lambda, \chi, \psi \mid \theta^{(m)}\right) \\
= & \sum_{i=1}^{n} \mathbb{E}_{W_{i} \mid x_{i}, \theta^{(m)}}\left[\ln f_{W_{i}}\left(W_{i} \mid \lambda, \chi, \psi\right)\right] \\
= & -\frac{n}{2} \lambda \ln \chi+\frac{n}{2} \lambda \ln \psi-n \ln \left(2 K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)\right)-\frac{\psi}{2} \sum_{i=1}^{n} a_{i}^{(m)}-\frac{\chi}{2} \sum_{i=1}^{n} b_{i}^{(m)}+(\lambda-1) \sum_{i=1}^{n} c_{i}^{(m)} .
\end{aligned}
$$

The M-step. This step deals with finding the argument when maximizing $Q$. Namely,

$$
\begin{aligned}
& \left(\mu^{(m+1)}, \Sigma^{(m+1)}, \gamma^{(m+1)}\right)=\arg \max _{\mu, \Sigma, \gamma} Q_{1}\left(\mu, \Sigma, \gamma \mid \theta^{(m)}\right) \\
& \left(\lambda^{(m+1)}, \chi^{(m+1)}, \psi^{(m+1)}\right)=\arg \max _{\lambda, \chi, \psi} Q_{2}\left(\lambda, \chi, \psi \mid \theta^{(m)}\right) .
\end{aligned}
$$

Monotonicity of Likelihood. Define the likelihood function as follows:

$$
l(\theta)=\log f_{\mathbf{X}}(\mathbf{x} \mid \theta)
$$

[9, Theorem 2.1] assets that each step in the EM algorithm increases the likelihood:

$$
\begin{equation*}
l(\theta) \geq l\left(\theta^{(m)}\right) \quad \text { whenever } \quad Q\left(\theta \mid \theta^{(m)}\right) \geq Q\left(\theta^{(m)} \mid \theta^{(m)}\right) \tag{2.14}
\end{equation*}
$$

We refer to [10] and [14, Chapters $3 \& 4]$ for convergence properties of the EM algorithm.

### 2.2 A Modified EM Algorithm

### 2.2.1 Modified MCECM Algorithm

Many variants of the EM algorithm have been introduced. See, e.g., [14, Chapters 5 \& 6]. Since the maximization of $Q$ is split into two components $Q_{1}$ and $Q_{2}$ which separate $\theta$ into two independent components $(\mu, \Sigma, \gamma)$ and $(\lambda, \chi, \psi)$, one sees that the EM algorithm coincides with the Expectation-Conditional Maximization algorithm (ECM), in which following each E-step the M-step (now called the CM-step) consists of a few consecutive maximizations subject to updated constraints on prescribed subsets of the parameter collection. If one also updates the E-step whenever a conditional maximization in the CM-step is conducted, we come to the MCECM algorithm (It should be alerted that MCECM also refers to Monte Carlo ECM in some literature). In what follows, we describe a modified version of the MCECM algorithm.

Suppose we have finished the $m$-th step and obtain $\theta^{(m)}$. After updating the E-step of computing $Q\left(\theta \mid \theta^{(m)}\right)$, we have the option to first maximize $Q_{1}$ to obtain

$$
\left(\mu^{(m+1,1)}, \Sigma^{(m+1,1)}, \gamma^{(m+1,1)}\right)
$$

, and then partially update the known parameters from $\theta^{(m)}$ to

$$
\theta^{(m, 0,1)}=\left(\lambda^{(m)}, \chi^{(m)}, \psi^{(m)}, \mu^{(m+1,1)}, \Sigma^{(m+1,1)}, \gamma^{(m+1,1)}\right)
$$

Then we continue to update the E-step, namely, calculate $Q\left(\theta \mid \theta^{(m, 0,1)}\right)$, and then maxi-
mize $Q_{1}$ again to partially update the known parameters from $\theta^{(m, 0,1)}$ to

$$
\theta^{(m, 0,2)}=\left(\lambda^{(m)}, \chi^{(m)}, \psi^{(m)}, \mu^{(m+1,2)}, \Sigma^{(m+1,2)}, \gamma^{(m+1,2)}\right)
$$

Repeating this process, one obtains

$$
\theta^{(m, 0, k)}=\left(\lambda^{(m)}, \chi^{(m)}, \psi^{(m)}, \mu^{(m+1, k)}, \Sigma^{(m+1, k)}, \gamma^{(m+1, k)}\right), \quad k \in \mathbb{N} .
$$

Intuitively,

$$
\theta^{(m, 0, \infty)}:=\lim _{k \rightarrow \infty} \theta^{(m, 0, k)}=\left(\lambda^{(m)}, \chi^{(m)}, \psi^{(m)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

gives the best fit of the model to the $n$ observations, when $\left(\lambda^{(m)}, \chi^{(m)}, \psi^{(m)}\right.$ is fixed, subject to the initial value $\left(\mu^{(m)}, \Sigma^{(m)}, \gamma^{(m)}\right)$ of $(\mu, \Sigma, \gamma)$.

Next, we update the E-step, calculating $Q\left(\theta \mid \theta^{(m, 0, \infty)}\right)$, and maximize $Q_{2}\left(\lambda, \chi, \psi \mid \theta^{(m, 0, \infty)}\right)$ to update the parameters from $\theta^{(m, 0, \infty)}$ to

$$
\theta^{(m, 1, \infty)}=\left(\lambda^{(m+1,1)}, \chi^{(m+1,1)}, \psi^{(m+1,1)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

We continue to update the E-step, calculating $Q\left(\theta \mid \theta^{(m, 1, \infty)}\right)$, and maximize $Q_{2}\left(\lambda, \chi, \psi \mid \theta^{(m, 1, \infty)}\right)$ to update the parameters from $\theta^{(m, 1, \infty)}$ to

$$
\theta^{(m, 2, \infty)}=\left(\lambda^{(m+1,2)}, \chi^{(m+1,2)}, \psi^{(m+1,2)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

Repeating the process, we obtain

$$
\theta^{(m, k, \infty)}=\left(\lambda^{(m+1, k)}, \chi^{(m+1, k)}, \psi^{(m+1, k)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right), \quad k \in \mathbb{N}
$$

and

$$
\theta^{(m, \infty, \infty)}:=\lim _{k \rightarrow \infty} \theta^{(m, k, \infty)}=\left(\lambda^{(m+1, \infty)}, \chi^{(m+1, \infty)}, \psi^{(m+1, \infty)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

Finally, we complete the $(m+1)$-th step by setting

$$
\theta^{(m+1)}:=\theta^{(m, \infty, \infty)}
$$

In reality, it may be unfeasible or unnecessary to obtain $\left.\theta^{(m, 0, \infty)}\right)$ and then $\theta^{(m, \infty, \infty)}$. Instead, one may stop first at $\theta^{\left(m, 0, k_{1}\right)}$ and turn to work on $Q_{2}$ and stop at $\theta^{\left(m, k_{2}, k_{1}\right)}$. In practice, $k_{1}=k_{2}=1$ is used. We will stick to $k_{1}=k_{2}=3$. Repeatedly using (2.14), one sees that

$$
l\left(\theta^{(m+1)}\right) \geq l\left(\theta^{(m)}\right)
$$

It deserves mentioning that, alternatively, one can first work on $Q_{2}$ and then on the partially updated $Q_{1}$. However, we will not take this alternative.

### 2.2.2 A further CM split of $\lambda, \chi, \psi$

We will use the MATLAB tool "fmincon" to optimize $Q_{2}$. In Section 2.1.2, one sees that $Q_{2}$ is highly nonlinear and has three variables $\lambda, \chi, \psi$. These two factors add a significant level of difficulties, instability and inaccuracy into the optimization task. Indeed, if one assumes the true values of $\mu, \Sigma, \gamma$ are known (and thus never update them in the EM algorithm), then simulation studies often reveal that the estimates of $\lambda, \chi, \psi$ produced by the EM algorithms may be wrong, in particular, when the true values of $\lambda, \chi, \psi$ are somewhat extreme. We include the MATLAB codes for these simulation studies on EM estimates of $W$, i.e., estimates of $\lambda, \chi, \psi$ in Appendix A.2. One reason for such failures is that "fmincon" can only find local minimums and has reduced efficiency over highdimensional regions.

On the other hand, if one assumes the true values of all parameters but $\lambda$ (respectively, $\chi, \psi)$ are known (and thus never update them in the EM algorithm), then simulation studies show that the EM algorithm produces expected estimates of $\lambda$ (respectively, $\chi, \psi)$. We include the MATLAB codes for these simulation studies on EM estimates of $W$, i.e., estimates of $\lambda, \chi, \psi$ in Appendix A.3.

Due to the above reasons, we modify the CM-step in the maximization of $Q_{2}$. Namely, we optimize $Q_{2}\left(\lambda, \theta_{*}^{(m, k)} \mid\left(\lambda^{(m+1, k)}, \theta_{*}^{(m, k)}\right)\right.$, where

$$
\theta_{*}^{(m, k)}=\left(\chi^{(m+1, k)}, \psi^{(m+1, k)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right),
$$

over $\lambda$ to update the parameters to the following:

$$
\left(\lambda^{(m+1, k+1)}, \chi^{(m+1, k)}, \psi^{(m+1, k)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

Then we optimize $Q_{2}\left(\chi, \theta_{* *}^{(m, k)} \mid\left(\chi^{(m+1, k)}, \theta_{* *}^{(m, k)}\right)\right.$, where

$$
\theta_{* *}^{(m, k)}=\left(\lambda^{(m+1, k+1)}, \psi^{(m+1, k)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right),
$$

over $\chi$ to update the parameters to the following:

$$
\left(\lambda^{(m+1, k+1)}, \chi^{(m+1, k+1)}, \psi^{(m+1, k)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

Finally, we optimize $Q_{2}\left(\psi, \theta_{* * *}^{(m, k)} \mid\left(\psi^{(m+1, k)}, \theta_{* * *}^{(m, k)}\right)\right.$, where

$$
\theta_{* * *}^{(m, k)}=\left(\lambda^{(m+1, k+1)}, \chi^{(m+1, k+1)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

over $\psi$ to update the parameters to the following:

$$
\left(\lambda^{(m+1, k+1)}, \chi^{(m+1, k+1)}, \psi^{(m+1, k+1)}, \mu^{(m+1, \infty)}, \Sigma^{(m+1, \infty)}, \gamma^{(m+1, \infty)}\right)
$$

### 2.3 Computational considerations

### 2.3.1 Initial values

For $X \sim \operatorname{GH}_{d}(\lambda, \chi, \psi, \mu, \Sigma, \gamma)$ as defined in (2.1), we have

$$
\begin{align*}
\mathbb{E}[X] & =\mu+\mathbb{E}[W] \gamma,  \tag{2.15}\\
\mathrm{V}[X] & =\mathrm{V}[W] \gamma \gamma^{\prime}+\mathbb{E}[W] \Sigma \tag{2.16}
\end{align*}
$$

In practice, one usually starts with

$$
\gamma^{(0)}=0,
$$

which corresponds to the special class of symmetric generalized hyperbolic distributions. Thus by (2.15), we set

$$
\mu^{(0)}=\frac{1}{n} \sum_{i=1}^{n} x_{i},
$$

and by (2.16), we set

$$
\Sigma^{(0)}=S,
$$

where $S$ is the sample covariance matrix. Now a second application of (2.16) inspires to restrict

$$
\mathbb{E}[W]=\left(\frac{\chi}{\psi}\right)^{\frac{1}{2}} \frac{K_{\lambda+1}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)}{K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)}=1
$$

We set

$$
\chi^{(0)}=\psi^{(0)}=1,
$$

and set $\lambda^{(0)}$ by solving

$$
K_{\lambda^{(0)}+1}(1)=K_{\lambda^{(0)}}(1),
$$

which by inspection has the solution $\lambda^{(0)}=-0.5$.

### 2.3.2 Computing $c_{i}^{(m)}$

The computation of $c_{i}^{(m)}$ involves evaluating $\mathbb{E}[\ln W]$, where $W \sim \mathrm{~N}^{-}(\lambda, \chi, \psi)$. We have at least three methods to do this. The first one is to use Monte Carlo simulation, namely, simulating iid $W_{i} \sim \mathrm{~N}^{-}(\lambda, \chi, \psi), i=1, \ldots, N$, and estimating

$$
\mathbb{E}[\ln W] \approx \frac{1}{N} \sum_{i=1}^{N} W_{i}
$$

A drawback is that the convergence is quite slow, with rate $O\left(N^{-\frac{1}{2}}\right)$.
Alternatively, by (2.5),

$$
\begin{align*}
\mathbb{E}[\ln W] & =\frac{\chi^{-\frac{\lambda}{2}} \psi^{\frac{\lambda}{2}}}{2 K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)} \int_{0}^{\infty}(\ln w) w^{\lambda-1} \exp \left(-\frac{\chi}{2} \frac{1}{w}-\frac{\psi}{2} w\right) \mathrm{d} w  \tag{2.17}\\
& =\frac{\partial \mathbb{E}\left[W^{\alpha}\right]}{\partial \alpha}(0) \tag{2.18}
\end{align*}
$$

Thus one may compute $\mathbb{E}[\ln W]$ via differentiation in (2.18) or via integration via improper integration in 2.17). To avoid numerical instability of differentiation, we will use integration, which can be easily done by implementing the integral command in MATLAB.

Therefore, since
$W_{i} \mid x_{i}, \theta^{(m)} \sim \mathrm{N}^{-}\left(\lambda^{(m)}-\frac{d}{2}, \chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right), \psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)$,
we have

$$
\begin{aligned}
c_{i}^{(m)}= & \left(\frac{\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}}{\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)}\right)^{\frac{\lambda^{(m)}-\frac{d}{2}}{2}} \\
& \times \frac{1}{2 K_{\lambda(m)-\frac{d}{2}}\left(\sqrt{\left(\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)\right)\left(\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}\right)}\right)} \\
& \times \int_{0}^{\infty}(\ln w) w^{\lambda^{(m)}-\frac{d}{2}-1} \exp \left(-\frac{\chi^{(m)}+\left(x_{i}-\mu^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1}\left(x_{i}-\mu^{(m)}\right)}{2} \frac{1}{w}\right. \\
& \left.-\frac{\psi^{(m)}+\left(\gamma^{(m)}\right)^{\prime}\left(\Sigma^{(m)}\right)^{-1} \gamma^{(m)}}{2} w\right) \mathrm{d} w
\end{aligned}
$$

### 2.3.3 Maximizing $Q_{1}$

We refer to [19] for basics on vector and matrix differentiation. Differentiating $Q_{1}$ we have

$$
\begin{aligned}
\frac{\partial Q_{1}}{\partial \mu} & =-n \Sigma^{-1} \gamma+\sum_{i=1}^{n} \Sigma^{-1}\left(x_{i}-\mu\right) b_{i}^{(m)} \\
\frac{\partial Q_{1}}{\partial \gamma} & =\sum_{i=1}^{n} \Sigma^{-1}\left(x_{i}-\mu\right)-\Sigma^{-1} \gamma \sum_{i=1}^{n} a_{i}^{(m)} \\
\frac{\partial Q_{1}}{\partial \Sigma^{-1}} & =\sum_{i=1}^{n}\left(x_{i}-\mu\right) \gamma^{\prime}-\frac{1}{2} \gamma \gamma^{\prime} \sum_{i=1}^{n} a_{i}^{(m)}-\frac{1}{2} \sum_{i=1}^{n}\left(x_{i}-\mu\right)\left(x_{i}-\mu\right)^{\prime} b_{i}^{(m)}+\frac{n}{2} \Sigma .
\end{aligned}
$$

Setting the partial derivatives to 0 and solving the equations, we obtain

$$
\begin{aligned}
\gamma^{(m+1)} & =\frac{\frac{1}{n} \sum_{i=1}^{n} b_{i}^{(m)}\left(x_{i}-\bar{x}\right)}{1-\frac{1}{n} \sum_{i=1}^{n} b_{i}^{(m)} \cdot \frac{1}{n} \sum_{i=1}^{n} a_{i}^{(m)}}, \\
\mu^{(m+1)} & =\frac{\sum_{i=1}^{n} x_{i}}{n}-\frac{\sum_{i=1}^{n} a_{i}}{n} \gamma, \\
\Sigma & =\frac{1}{n} \sum_{i=1}^{n}\left(x_{i}-\mu^{(m+1)}\right)\left(x_{i}-\mu^{(m+1)}\right)^{\prime} b_{i}^{(m)}-\frac{1}{n} \sum_{i=1}^{n} a_{i}^{(m)} \cdot \gamma^{(m+1)}\left(\gamma^{(m+1)}\right)^{\prime}
\end{aligned}
$$

### 2.3.4 Maximizing $Q_{2}$

In practice, one shall maximize $2 Q_{2} / n$ to remove the magnitude $n$ in case the data size is very large and unnecessary constants. Namely, we need to find the argument in minimizng the following function:

$$
\begin{aligned}
& Q_{2}^{\prime}\left(\lambda, \chi, \psi \mid \theta^{(m)}\right) \\
= & \lambda \ln \chi-\lambda \ln \psi+2 \ln \left(K_{\lambda}\left(\chi^{\frac{1}{2}} \psi^{\frac{1}{2}}\right)\right)+\psi \frac{\sum_{i=1}^{n} a_{i}^{(m)}}{n}+\chi \frac{\sum_{i=1}^{n} b_{i}^{(m)}}{n}-2 \lambda \frac{\sum_{i=1}^{n} c_{i}^{(m)}}{n} .
\end{aligned}
$$

We will use the MATLAB tool "fmincon" to solve this nonlinear optimization problem with constraints: $\chi \geq 0$ and $\psi \geq 0$. Note that the precise constraints of the parameters $\lambda, \chi, \psi$ are as follows: $\chi \geq 0, \psi>0$ if $\lambda>0 ; \chi>0, \psi>0$ if $\lambda=0 ; \chi>0, \psi \geq 0$ if $\lambda<0$.

### 2.4 Numerical Results

In this section, we discuss the results from calibrating and simulating Generalized Hyberbolic Distributions with our sample data.

### 2.4.1 EM Estimates of Parameters

Running the code in Appendix A. 1 enables us to calculate the best estimates of the parameters involved in the Generalized Hyperbolic Distribution. We list them here.

$$
\begin{aligned}
& \lambda=-1.8552 \\
& \chi=2.2326 \\
& \psi=5.1476 * 10^{-5}
\end{aligned}
$$

$$
\begin{aligned}
& \gamma^{\prime}=\left[\begin{array}{l}
3.5261 * 10^{-3} \\
2.4177 * 10^{-3} \\
1.0796 * 10^{-3} \\
2.6350 * 10^{-3}
\end{array}\right] \\
& \mu^{\prime}=\left[\begin{array}{llll}
-1.0028 * 10^{-2} \\
-3.739 * 10^{-3} \\
-1.2105 * 10^{-3} \\
-6.1779 * 10^{-3}
\end{array}\right] \\
& \Sigma=\left[\begin{array}{llll}
2.9068 * 10^{-3} & 6.0378 * 10^{-4} & 6.8643 * 10^{-4} & 5.7658 * 10^{-4} \\
6.0378 * 10^{-4} & 2.1498 * 10^{-3} & 9.8082 * 10^{-4} & 6.0630 * 10^{-4} \\
6.8643 * 10^{-4} & 9.8082 * 10^{-4} & 2.4551 * 10^{-3} & 6.4051 * 10^{-4} \\
5.7658 * 10^{-4} & 6.0630 * 10^{-4} & 6.4051 * 10^{-4} & 2.1930 * 10^{-3}
\end{array}\right]
\end{aligned}
$$

### 2.4.2 Simulation Results

The parameters are then used in conjunction with Monte Carlo simulations to generate $1,000,000$ simulations of the stock return values. Following this, we take the rowsums to evaluate the loss for the portfolio. For this, we assume that we have invested $\$ 10,000$ in each of the four stocks. The results are listed in Table 2.1

| Expected Shortfall | Value at Risk |
| :--- | :--- |
| $\$ 4,793$ | $\$ 4,352$ |

Table 2.1: ES and VaR based on Generalized Hyperbolic Distribution

We will come back to these numbers in Chapter 4, where we will discuss results along with conclusions and inferences.

## Chapter 3

## Risk Modelling via Marginal Distributions

If we have a joint distribution function corresponding to a risk factor random vector, we have a dependence structure that is implicit in it. Copulas help us in the isolation of this dependence structure.

Dependence and correlation may seem synonymous at first glance, however, a deeper look reveals that linear correlations are inadequate to measure dependence. To this end, we can discuss alternative dependence measures including coefficients of tail dependence using copulas. The latter is of particular interest in the study of financial market crashes, since there may be a high degree of dependence in the tails which represent extreme events.

### 3.1 Basic Copula Theory

In this section we review the definition of a copula, some basic properties that a copula must satisfy and some examples of copulas. We refer to [18] for a comprehensive account of copulas. Throughout this chapter, all distributions are continuous, unless specified
otherwise.

### 3.1.1 Definitions

Definition 1. Ad-dimensional copula is a function $C:[0,1]^{d} \rightarrow[0,1]$ satisfying the following conditions:
(i) $C\left(u_{1}, \ldots, u_{d}\right)$ is increasing in each component $u_{i}$;
(ii) $C\left(1, \ldots, 1, u_{i}, 1, \ldots, 1\right)=u_{i}$ for all $i \in\{1, \ldots, d\}$ and $u_{i} \in[0,1]$;
(iii) For any $\left(a_{1}, \ldots, a_{d}\right),\left(b_{1}, \ldots, b_{d}\right) \in[0,1]^{d}$ with $a_{i} \leq b_{i}$ for all $1 \leq i \leq d$, it holds that

$$
\sum_{i_{1}=1}^{2} \cdots \sum_{i_{d}=1}^{2}(-1)^{i_{1}+\ldots+i_{d}} C\left(u_{1 i_{1}}, \ldots, u_{d i_{d}}\right) \geq 0
$$

where $u_{j 1}=a_{j}$ and $u_{j 2}=b_{j}$ for all $j \in\{1, \ldots, d\}$.
As is well-known from distribution theory, copulas can be equivalently defined as follows.

Definition 2. A function $C:[0,1]^{d} \rightarrow[0,1]$ is a copula if and only if there exists a random vector $U=\left(U_{1}, \ldots, U_{d}\right)$ such that $U_{i}$ is a standard uniform distribution on $[0,1]$ for each $1 \leq i \leq d$ and that

$$
C\left(u_{1}, \ldots, u_{d}\right)=\mathbb{P}\left(U_{1} \leq u_{1}, \ldots, U_{d} \leq u_{d}\right), \forall\left(u_{1}, \ldots, u_{d}\right) \in[0,1]^{d}
$$

We call this random vector $U$, although not unique, a generating random vector of $C$. It often plays a useful role in simulations.

### 3.1.2 Basic Properties

Next we proceed to discuss some properties of copulas. Before we do that, we observe the following properties for distribution functions.

Let $F$ be a continuous distribution function and let $F^{-}$denote the left continuous quantile of $F$, i.e.,

$$
F^{-}(y)=\inf \{x: F(x) \geq y\}
$$

Then the following hold:
(a) If $U$ is a random variable following standard uniform distribution, then the distribution function of $F^{-}(U)$ is $F$;
(b) If $X$ is a random variable with distribution function $F$, then $F(X)$ follows the standard uniform distribution.

We refer to [6] for detailed discussions of these properties.
The primary importance of copulas in isolating dependence structure from marginal distributions is illustrated in the celebrated theorem of Sklar.

Theorem 3 (Sklar). Let $F$ be a joint distribution function with marginals $F_{1}, \ldots, F_{d}$. Then there exists a unique copula $C:[0,1]^{d} \rightarrow[0,1]$ such that, for all $x_{1}, \ldots, x_{d} \in \mathbb{R}$,

$$
\begin{equation*}
F\left(x_{1}, \ldots, x_{d}\right)=C\left(F_{1}\left(x_{1}\right), \ldots, F_{d}\left(x_{d}\right)\right) \tag{3.1}
\end{equation*}
$$

Clearly, by (3.1), for any $u_{1}, \ldots, u_{d} \in[0,1]$,

$$
\begin{equation*}
C\left(u_{1}, \ldots, u_{d}\right)=F\left(F_{1}^{-}\left(u_{1}\right), \ldots, F_{d}^{-}\left(u_{d}\right)\right) . \tag{3.2}
\end{equation*}
$$

Definition 4. We call $C$ in Theorem 3 the copula of $F$. If $F$ is the distribution of $a$ random vector $\left(X_{1}, \ldots, X_{d}\right)$, we also call $C$ the copula of $\left(X_{1}, \ldots, X_{d}\right)$, or $\left(X_{1}, \ldots, X_{d}\right)$ an associated random vector of $C$.

We now make two useful remarks regarding generating and associated random vectors, combining which gives the standard procedures to simulate random vectors with a given copula and given marginals.

Remark 5. Let $C$ be a copula with associated random vector $\left(X_{1}, \ldots, X_{d}\right)$. Let $F_{i}$ be the distribution of $X_{i}$. Note that each $F_{i}\left(X_{i}\right)$ is a standard uniform distribution and that

$$
\begin{aligned}
& \mathbb{P}\left(F_{1}\left(X_{1}\right) \leq u_{1}, \ldots, F_{d}\left(X_{d}\right) \leq u_{d}\right) \\
= & \mathbb{P}\left(F_{1}^{-} \circ F_{1}\left(X_{1}\right) \leq F_{1}^{-}\left(u_{1}\right), \ldots, F_{d}^{-} \circ F_{d}\left(X_{d}\right) \leq F_{d}^{-}\left(u_{d}\right)\right) \\
= & \mathbb{P}\left(X_{1} \leq F_{1}^{-}\left(u_{1}\right), \ldots, X_{d} \leq F_{d}^{-}\left(u_{d}\right)\right) \\
= & F\left(F_{1}^{-}\left(u_{1}\right), \ldots, F_{d}^{-}\left(u_{d}\right)\right) \\
= & C\left(u_{1}, \ldots, u_{d}\right) .
\end{aligned}
$$

Therefore, $\left(F_{1}\left(X_{1}\right), \ldots, F_{d}\left(X_{d}\right)\right)$ is a generating random vector of $C$

Remark 6. Let $C$ be a copula with a generating random vector $\left(U_{1}, \ldots, U_{d}\right)$. Let $F_{1}, \ldots, F_{d}$ be given marginals. Note that each $F_{i}^{-}\left(U_{i}\right)$ has distribution function $F_{i}$ and

$$
\begin{aligned}
& \mathbb{P}\left(F_{1}^{-}\left(U_{1}\right) \leq x_{1}, \ldots, F_{d}^{-}\left(U_{d}\right) \leq x_{d}\right) \\
= & \mathbb{P}\left(U_{1} \leq F_{1}\left(x_{1}\right), \ldots, U_{d} \leq F_{d}\left(x_{d}\right)\right) \\
= & C\left(F_{1}\left(x_{1}\right), \ldots, F_{d}\left(x_{d}\right)\right) .
\end{aligned}
$$

Therefore, by 3.1, $\left(F_{1}^{-}\left(U_{1}\right), \ldots, F_{d}^{-}\left(U_{d}\right)\right)$ is an associated random vector of $C$ with prescribed marginals $F_{i}$ 's.

We end this subsection with a useful result that allows us to simplify the associated random vector when determining its copula. This result demonstrates again the fact that copulas capture the essential dependence structure in a joint distribution.

Proposition 7. Let $\left(X_{1}, \ldots, X_{d}\right)$ be a random vector with copula $C$ and let $T_{1}, \ldots, T_{d}$ be strictly increasing functions on $\mathbb{R}$. Then $\left(T_{1}\left(X_{1}\right), \ldots, T_{d}\left(X_{d}\right)\right)$ also has copula $C$.

### 3.1.3 Examples of Copulas

We provide a number of examples of copulas in this section and these are subdivided into three categories: fundamental copulas represent a number of important special dependence structures; implicit copulas are extracted from well-known multivariate distributions using Sklar's Theorem, but do not necessarily possess simple closed- form expressions; explicit copulas have simple closed-form expressions and follow general mathematical constructions known to yield copulas.

## Independence Copula

As a first example of fundamental copulas, the independence copula is defined by

$$
\begin{equation*}
\Pi(u, ., u)=\prod_{i=1}^{d} u_{i} . \tag{3.3}
\end{equation*}
$$

By Sklar's Theorem, it is clear that a random vector has the independence copula if and only if its marginals are independent.

## Comonotonicity Copula

The comonotonicity copula is another basic example of fundamental copulas. It is defined by

$$
\begin{equation*}
M\left(u_{1}, \ldots, u_{d}\right)=\min \left\{u_{1}, \ldots, u_{d}\right\} . \tag{3.4}
\end{equation*}
$$

Observe that this copula has a generating random vector $(U, \ldots, U)$, where $U$ follows a standard uniform distribution on $[0,1]$.

Let $X$ be a random variable and $T_{i}, i=1, \ldots, d$, be strictly increasing functions
on $\mathbb{R}$. Then by Proposition 7, the copula of $\left(T_{1}(X), \ldots, T_{d}(X)\right)$ is the same as that of $(X, \ldots, X)$, which in turn is the same as that of $(U, \ldots, U)$, where $U=F(X)$. Hence, the copula is in fact just the comonotonicity copula.

## Gauss Copula

Gauss copula is one of the most important implicit copulas. By Proposition 7, $X \sim$ $N_{d}(\mu, \Sigma)$ and $Y \sim N_{d}(0, P)$ have the same copula, where $P$ is the corresponding correlation matrix of $\Sigma$. Their copula is called a Gauss copula. Specifically, by Remark 5, we have

$$
\begin{aligned}
& C_{P}^{\mathrm{Ga}}\left(u_{1}, \ldots, u_{d}\right) \\
= & \mathbb{P}\left(\Phi\left(Y_{1}\right) \leq u_{1}, \ldots, \Phi\left(X_{d}\right) \leq u_{d}\right) \\
= & \Phi_{P}\left(\Phi^{-1}\left(u_{1}\right), \ldots, \Phi^{-1}\left(u_{d}\right)\right) \\
= & \int_{-\infty}^{\Phi^{-1}\left(u_{1}\right)} \cdots \int_{-\infty}^{\Phi^{-1}\left(u_{d}\right)} \operatorname{det}(2 \pi P)^{-\frac{1}{2}} \exp \left\{-\frac{1}{2} y^{\prime} P^{-1} y\right\} \mathrm{d} y .
\end{aligned}
$$

Gauss copulas were most popular in risk modelling. But some academics and practitioners attribute the 2007-2008 crisis to its wide use in credit risk modelling and its weakness to capture dependence, and since then its use has been greatly reduced.

## t Copula

In the same way that we can extract a copula from the multivariate normal distribution, we can extract an implicit copula from any other distribution with continuous marginal
dfs. For example, the $d$-dimensional $t$ copula takes the form

$$
\begin{aligned}
& C_{v, P}^{t}\left(u_{1}, \ldots, u_{d}\right) \\
= & t_{v, P}\left(t_{\nu}^{-1}\left(u_{1}\right), \ldots, t_{\nu}^{-1}\left(u_{d}\right)\right) \\
= & \int_{-\infty}^{\Phi^{-1}\left(u_{1}\right)} \cdots \int_{-\infty}^{\Phi^{-1}\left(u_{d}\right)} \frac{\Gamma[(\nu+d) / 2]}{\Gamma(\nu / 2) \nu^{d / 2} \pi^{d / 2}|\Sigma|^{1 / 2}}\left[1+\frac{1}{\nu}(x-\mu)^{\mathrm{T}} \Sigma^{-1}(x-\mu)\right]^{-(\nu+d) / 2} \mathrm{~d} x .
\end{aligned}
$$

$t$ copulas capture both upper and lower tail dependence. Their capture of dependence are stronger than Gauss copula but weaker than the following two copulas each of which only capture one-sided tail dependence for the upper and lower tail dependence.

## Gumbel Copula

The Gumbel copula is a typical explicit copula, which captures upper tail-dependence.

$$
\begin{equation*}
C_{\theta}^{\mathrm{Gu}}\left(u_{1}, \ldots, u_{d}\right)=\exp \left\{-\left(\sum_{i=1}^{d}\left(-\ln u_{i}\right)^{\theta}\right)^{\frac{1}{\theta}}\right\}, \quad 1 \leq \theta<\infty . \tag{3.5}
\end{equation*}
$$

## Clayton Copula

The last example is the Clayton copula, which is also an explicit copula but captures lower tail-dependence.

$$
\begin{equation*}
C_{\theta}^{\mathrm{Cl}}\left(u_{1}, \ldots, u_{d}\right)=\left(u_{1}^{-\theta}+\cdots+u_{d}^{-\theta}-d+1\right)^{-1 / \theta}, 0<\theta<\infty . \tag{3.6}
\end{equation*}
$$

### 3.2 Marginal Distributions of stocks

We will work with 20 years of data for 4 stocks, which are Apple (AAPL), Bank of America (BAC), Ford (F) and Suncor Energy (SU). The reason for choosing these particular stocks was that they are among the largest in their respective sectors and hail from


Figure 3.1: Marginal distributions of individual stocks.
different sectors, thereby allowing for some diversification. We assume that the marginal distributions of the stocks are normal. This assumption is chosen for simplicity, but may be relaxed to include any other distribution with no change in the procedure. Under this assumption, we calibrate the data set and obtain the distributions for the four stocks in Figure 3.1. The corresponding values are listed in Table 3.1.

Next, we assume that we have invested $\$ 10,000$ in each of the four stocks. In accordance with Basel regulations, we use the 99th percentile for VaR and 97.5 th percentile

| Stock | $\boldsymbol{\mu}$ | $\boldsymbol{\sigma}$ |
| :--- | :--- | :--- |
| AAPL | -0.0053 | 0.0585 |
| BAC | -0.0005 | 0.0629 |
| F | 0.0002 | 0.0618 |
| SU | -0.0027 | 0.0500 |

Table 3.1: Mean and standard deviation of log returns of marginal distributions of stocks onwards for ES. The ES and VaR values based on 1,000,000 simulations are tabulated in Table 3.2.

| Stock | Expected Shortfall | Value at Risk |
| :--- | :--- | :--- |
| AAPL | $\$ 1,312$ | $\$ 1,311$ |
| BAC | $\$ 1,415$ | $\$ 1,310$ |
| F | $\$ 1,378$ | $\$ 1,323$ |
| SU | $\$ 1,076$ | $\$ 1,056$ |
| Sum | $\$ 5,178$ | $\$ 4,999$ |

Table 3.2: ES and VaR based on marginal distributions of stocks

### 3.3 Empirical Results

For each of the four copulas defined in Section 2.1.3, and for the empirical marginal distributions determined in Section 2.2, we evaluate the results of plugging in the marginal distributions into the copulas.

### 3.3.1 Gauss Copula

We start with the Gauss Copula. For this, we need a correlation matrix. As suggested in [15], the following is a close approximation to the calibrated correlation matrix:

$$
\rho_{\mathrm{S}}\left(X_{i}, X_{j}\right)=(6 / \pi) \arcsin \frac{1}{2} \rho_{i j} \approx \rho_{i j}
$$

where $\rho_{\mathrm{S}}$ is the matrix formed by taking the pairwise Spearman's rank correlation coefficients between the stock return vectors. The matrix $\rho$ then takes the following form based on the empirical data:

$$
\rho=\left(\begin{array}{llll}
1.0000 & 0.2713 & 0.2843 & 0.2329 \\
0.2713 & 1.0000 & 0.4137 & 0.2439 \\
0.2843 & 0.4137 & 1.0000 & 0.2650 \\
0.2329 & 0.2439 & 0.2650 & 1.0000
\end{array}\right)
$$

We then simulate the copula by the following steps:

## Algorithm 2.1 (simulation of Gauss copula).

(1) Generate $Z \sim N_{d}(0, P)$;
(2) Return $U=\left(\Phi\left(Z_{1}\right), \ldots, \Phi\left(Z_{d}\right)\right)^{\prime}$, where $\Phi$ is the standard normal df. The random vector $U$ has df $C_{P}^{\mathrm{Ga}}$;
(3) Apply Remark 6 with the marginal distributions of the stocks as calculated in section 3.2 to obtain a random vector with $C_{P}^{\mathrm{Ga}}$.

Once we have the copula, and plugged in the marginals, we take the rowsums to evaluate the loss for the portfolio. For this, we assume that we have invested $\$ 10,000$ in each of the four stocks. The results are listed in Table 3.3

| Expected Shortfall | Value at Risk |
| :--- | :--- |
| $\$ 3,388$ | $\$ 3,303$ |

Table 3.3: ES and VaR based on Gauss Copula

As expected, both VaR and ES decrease when the stocks are put in a portfolio. The number for Value at Risk is $\$ 3,303$ and for Expected Shortfall is $\$ 3,388$, down from $\$ 4,999$ and $\$ 5,178$ respectively.

### 3.3.2 t Copula

The next copula that we discuss is the t Copula. For this, we need a correlation matrix and degrees of freedom. We may choose any arbitrary correlation matrix, however, as suggested in [15] the following relation holds:

$$
\rho_{\tau}\left(X_{i}, X_{j}\right)=(2 / \pi) \sin ^{-1} \rho_{i j}
$$

where $\rho_{\tau}$ is the matrix formed by taking the pairwise Kendall's rank relation coefficients between the stock return vectors.

It then follows, that our matrix may be estimated by multiplying both sides by $\pi / 2$ and subsequently taking the sine of both sides leading to the following for the purposes of this simulation:

$$
\rho=\left(\begin{array}{llll}
1.0000 & 0.2866 & 0.3007 & 0.2477 \\
0.2866 & 1.0000 & 0.4415 & 0.2612 \\
0.3007 & 0.4415 & 1.0000 & 0.2858 \\
0.2477 & 0.2612 & 0.2858 & 1.0000
\end{array}\right)
$$

We also choose $\nu$ to be 5 . The reasoning behind this is that we want to differentiate this from the Gauss Copula (since t Copula converges asymptotically to the Gauss Copula as degrees of freedom increase) and in order to observe the strong tail dependence present in the t Copula by design.

We then simulate the copula by the following steps:

## Algorithm 2.2 (simulation of $t$ copula).

(1) Generate $X \sim t_{d}(\nu, 0, P)$.
(2) Return $U=\left(t_{\nu}\left(X_{1}\right), \ldots, t_{\nu}\left(X_{d}\right)\right)^{\prime}$, where $t_{\nu}$ denotes the df of a standard univariate $t$ distribution. The random vector $U$ has df $C_{v, P}^{t}$.
(3) Apply Remark 6 with the marginal distributions of the stocks as calculated in section 3.2 to obtain a random vector with $C_{v, P}^{t}$.

Once we have the copula, and plugged in the marginals, we take the rowsums to evaluate the loss for the portfolio. For this, we assume that we have invested $\$ 10,000$ in each of the four stocks. The results are listed in Table 3.4.

| Expected Shortfall | Value at Risk |
| :--- | :--- |
| $\$ 3,700$ | $\$ 3,660$ |

Table 3.4: ES and VaR based on $t$ Copula

As expected, both VaR and ES decrease when the stocks are put in a portfolio even for the t Copula. The number for Value at Risk is $\$ 3,660$ and for Expected Shortfall is $\$ 3,700$, down from $\$ 4,999$ and $\$ 5,178$ respectively. Also note that while the Value at Risk increasing by using a t Copula as opposed to a Gauss Copula, the Expected Shortfall actually goes down. This is due to the fact that for the same correlation matrix $\rho$, the t Distribution has heavier tails than the Normal Distribution.

### 3.3.3 Gumbel Copula

Another copula that we discuss in greater detail is the Gumbel copula. The interesting feature about this copula is that it captures upper tail-dependence. A Gumbel copula is uniquely characterized by the parameter $\theta \geq 1$.

Calibrating $\theta$ in the two-dimensional case boils down to a one-to-one mapping between the Kendall's rank correlation coefficient and $\theta$, however, it is not trivial to calibrate in the d-dimensional case. For the purpose of our simulations, we choose $\theta=3$, and then run the following algorithm to simulate the Gumbel Copula:

## Algorithm 2.3 (simulation of Gumbel copula).

(1) First, generate a variate $V$ with $\mathrm{df} G$ such that $\hat{G}$, the Laplace-Stieltjes transform of $G$, will be the inverse of the generator $\phi$ of the required copula.
(2) Next, generate independent uniform variates $X_{1}, \ldots, X_{d}$.
(3) Return $U=\left(\hat{G}\left(-\ln \left(X_{1}\right) / V\right), \ldots, \hat{G}\left(-\ln \left(X_{d}\right) / V\right.\right.$ where V is a positive stable variate $V \sim \operatorname{St}(1 / \theta, 1, \gamma, 0)$, where $\gamma=(\cos (\pi /(2 \theta)))^{\theta}$ and $\theta>1$. The resultant df has Laplace transform $\hat{G}(t)=\exp \left(-t^{1 / \theta}\right)$ which is what we wanted.
(4) Apply Remark 6 with the marginal distributions of the stocks as calculated in section 3.2 to obtain a random vector with $C_{\theta}^{G u}$.

Once we have the copula, and plugged in the marginals, we take the rowsums to evaluate the loss for the portfolio. For this, we assume that we have invested $\$ 10,000$ in each of the four stocks. The results are listed in Table 3.5.

| Expected Shortfall | Value at Risk |
| :--- | :--- |
| $\$ 5,613$ | $\$ 5,544$ |

Table 3.5: ES and VaR based on Gumbel Copula

Due to high dependence in the upper tails, we have that the Value at Risk and Expected Shortfall increase from $\$ 4,999$ and $\$ 5,178$ to $\$ 5,544$ and $\$ 5,613$ respectively. This can also partly be explained by the instability of the random number generator for Positive Stable Variates in MATLAB.

We will come back to these numbers in Chapter 4, where we will discuss results along with conclusions and inferences.

## Chapter 4

## Conclusions and Inferences

This chapter is dedicated to summarizing and dissecting the results obtained in Chapters 2 and 3 , and attempting to explain them using both Mathematics and Finance. We will begin with numerical comparisons, and then proceed to inferences. Finally, we will finish by discussing potential for future research.

### 4.1 Numerical Comparisons

First, we summarize the results obtained by the various copulas in Table 4.1.

| Copula Name | Expected Shortfall | Value at Risk |
| :--- | :--- | :--- |
| Gauss Copula | $\$ 3,388$ | $\$ 3,303$ |
| t Copula | $\$ 3,700$ | $\$ 3,660$ |
| Gumbel Copula | $\$ 5,613$ | $\$ 5,544$ |

Table 4.1: ES and VaR based on various Copulas

It is evident from the table that both Gauss Copula and t Copula are able to capture the concept of risk reduction due to diversification [13], whereas the Gumbel Copula amplifies upper tail risk.

Next, we compare this with the numbers obtained by using the Generalized Hyperbolic Distribution. Recall that the values for the Value at Risk and Expected Shortfall were $\$ 4,352$ and $\$ 4,793$ respectively. While these are higher than those obtained by using the Gauss and t Copulas, they are significantly lower than the values of the Gumbel Copula. It is worth stressing again that this is the most generic version of the Generalized Hyperbolic Distribution, with all parameters being free.

Another comparison can be drawn between the values obtained by using the sum of the individual Value at Risk and Expected Shortfall for the various stocks versus the values obtained by putting them in a portfolio first and subsequently applying either Copulas or Generalized Hyperbolic Distributions to the complete portfolio. Recall that the sum of the individual values was $\$ 5,178$ and $\$ 4,999$ for the Expected Shortfall and Value at Risk respectively. While the Generalized Hyperbolic Distribution, and the Gauss and t Copulas result in a reduction of risk, the Gumbel Copula results in an increase in risk.

### 4.2 Inferences

The class of Generalized Hyperbolic Distributions encompasses a wide range of distributions due to the fact that there are six parameters in total. This inherently allows for tweaking and fitting according to real life needs. This customization is useful when a simple explanatory model (for example, the Multivariate Normal Model), may be insufficient to capture the intricacies of the data set.

Copulas are an elegant way of measuring and allowing for dependence in portfolios that may otherwise be difficult to capture. Different copulas have different properties, and we have to find one that suits our needs. Often, the most extreme case (in this case, the Gumbel Copula) can be taken from the point of view of a conservative risk-adverse investor. For other practical purposes (for example, internal and external reporting),
other copulas like the Gauss and the t Copulas may be better fits.

### 4.3 Future Research

The following are questions that can be considered extensions to the work done in this thesis and may be suitable for researching further:

1. Is it worth defining the Generalized Hyperbolic Distribution in its most generic form (using all six parameters), or is there merit to simplifying the problem by fixing one or more parameters, thereby specifying a subset of Generalized Hyperbolic Distributions (several of which are named distributions)?
2. How do these two methods, namely using Copulas or Generalized Hyperbolic Distributions, stack against other known methods of estimation in literature, for example, the Rearrangement Algorithm proposed in [17]?
3. How well do these methods capture risk in real life? This is a question that is quite open-ended and subjective, but having a metric that quantifies and compares the performance of different methods in the existing financial markets can be useful.
4. What other risk measures (for example, the Entropic Risk measure [5]) can be suitable for quantifying financial risk? What properties do these risk measures satisfy when applied to individual stocks? More importantly, what properties still hold (or come into existence) while discussing portfolio risk?

## Appendix A

## Codes for Chapter 2

## A. 1 EM algorithm: Calibration and Simulation

```
% clearvars -except returnmatrix
```

$\mathrm{X}=$ returnmatrix ;
$[\mathrm{n}, \mathrm{d}]=\operatorname{size}($ returnmatrix) ; \% calculate dimensions of return matrix

```
nsimulations = 1000000; % number of simulations to run
```

confidence_level_VaR $=0.99 ; \%$ confidence level for VaR
confidence_level_ES $=0.975 ; \%$ confidence level for ES
dollarvalue $=10000 ; \%$ amount invested in each stock
\% initial values
gamma $=\operatorname{zeros}(1, \mathrm{~d}) ;$
$\mathrm{mu}=\operatorname{mean}(\mathrm{X}) ;$
Sigma $=\operatorname{cov}(X) ;$
psi $=1$;
chi $=1 ;$
lambda $=-0.5$;
iterations $=100$;
klambda $=5 ;$
kchi $=5$;
$\mathrm{kpsi}=5$;
$\mathrm{kW}=5$;
for iter = 1:iterations
\% lambda optimization
chiwx $=$ chi+sum $(((-m u+X) /$ Sigma $) . *(-m u+X), 2) ; \%$ this is $n x 1$
matrix
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
for $\mathrm{j}=1$ :klambda
\%compute the density of W_i|x_i
lambdawx $=$ lambda - d $/ 2$;
\%compute c
$\mathrm{c}=\mathrm{zeros}(\mathrm{n}, 1) ; \%$ this is nx1 matrix: $\mathrm{E}[\ln \mathrm{W}]$
for $\mathrm{i}=1: 1$ : n
fun $1=@(\mathrm{w}) \log (\mathrm{w}) . * \mathrm{w} .{ }^{\wedge}(\operatorname{lambdawx}-1) . * \exp (-0.5 *$
$\operatorname{psiwx} * w) . * \exp \left(-0.5 * \operatorname{chiwx}(\mathrm{i}, 1) * w .^{\wedge}(-1)\right) ;$
$c(i, 1)=0.5 *(\operatorname{psiwx} / \operatorname{chiwx}(\mathrm{i}, 1))^{\wedge}(0.5 * \operatorname{lambdawx}) /$
besselk(lambdawx, sqrt(chiwx (i, 1$) * \operatorname{psiwx})) *$

```
            integral(fun1,0,Inf);
                clear fun1;
            end
            %update lambda, chi, psi
            options = optimoptions('fmincon','Algorithm','sqp','
            Display ','off');
    fun2 = @(t) t*log(chi)-t*log(psi) +2*log(besselk(t,
        sqrt(chi*psi))) - 2*t*mean(c);
    A = [ ];
    B = [ ];
    Aeq = [];
    beq = [];
    lb = [];
    ub = [];
    nonlcon = [];
    t0 = lambda;
    t = fmincon(fun2, t0,A,B,Aeq, beq, lb,ub, nonlcon,
        options);
    clear fun2;
    %print out the result
    step = j;
        lambda = t;
    end
    % chi optimization
    lambdawx = lambda-d/2;
```

```
psiwx = psi+gamma/Sigma*transpose(gamma);
for j=1:kchi %work on Q2 and partially update lambda,chi,psi
        for k times
```

    \%compute the density of W_i|x_i
        chiwx \(=\operatorname{chi}+\operatorname{sum}(((-m u+X) /\) Sigma \() \cdot *(-m u+X), 2) ; \%\) this is
        nx1 matrix
    \%compute a,b, c
        \(\mathrm{b}=\mathrm{sqrt}(\mathrm{psiwx} . /\) chiwx\() . *\) besselk (lambdawx -1, sqrt (chiwx \(*\)
        psiwx))./besselk(lambdawx, sqrt(chiwx*psiwx));\% this
        is nx1 matrix: \(\mathrm{E}\left[\mathrm{W}^{\wedge}(-1)\right]\)
    \%update lambda, chi, psi
        options \(=\) optimoptions('fmincon','Algorithm', 'sqp',
        Display ', 'off');
        fun2 \(=@(\mathrm{t}) \operatorname{lambda} * \log (\mathrm{t})+2 * \log (\) besselk (lambda, sqrt \((\mathrm{t} *\)
        psi)) ) \(+\mathrm{t} *\) mean (b);
    \(\mathrm{A}=-1 ;\)
    \(B=0 ;\)
    Aeq \(=[] ;\)
    beq \(=[] ;\)
    \(\mathrm{lb}=[]\);
    \(\mathrm{ub}=[] ;\)
    nonlcon \(=[] ;\)
    \(\mathrm{t} 0=\mathrm{chi} ;\)
    \(\mathrm{t}=\mathrm{fmincon}(\) fun2, \(\mathrm{t} 0, \mathrm{~A}, \mathrm{~B}, \mathrm{Aeq}, \mathrm{beq}, \mathrm{lb}, \mathrm{ub}\), nonlcon, options);
    \%print out the result
    step \(=\mathrm{j}\);
    chi \(=t ;\)
    end
\% psi optimization

$$
\text { lambdawx }=\text { lambda }-\mathrm{d} / 2 \text {; }
$$

$$
\operatorname{chiwx}=\operatorname{chi}+\operatorname{sum}(((-m u+X) / \operatorname{Sigma}) \cdot *(-\operatorname{mu}+X), 2) ; \% \text { this is nx1 }
$$

matrix

$$
\text { for } j=1: \mathrm{kpsi}
$$

\%compute the density of W_i|x_i
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
\%compute $\mathrm{a}, \mathrm{b}, \mathrm{c}$
$\mathrm{a}=\mathrm{sqrt}($ chiwx $/ \mathrm{psiwx}) . *$ besselk (lambdawx +1 , sqrt (chiwx $*$ psiwx))./besselk(lambdawx, sqrt(chiwx*psiwx));\% this is nx1 matrix: $\mathrm{E}[\mathrm{W}]$
\%update lambda, chi, psi
options $=$ optimoptions ('fmincon', 'Algorithm ', 'sqp ', ' Display ', 'off');
fun2 $=@(\mathrm{t})-\operatorname{lambda} * \log (\mathrm{t})+2 * \log (\mathrm{besselk}(\operatorname{lambda}$, sqrt (chi *t) ) $)+\mathrm{t} * \operatorname{mean}(\mathrm{a})$;
$\mathrm{A}=-1 ;$
$\mathrm{B}=0$;
Aeq $=[] ;$
beq $=[] ;$
$\mathrm{lb}=[] ;$
$\mathrm{ub}=[] ;$
nonlcon $=[] ;$
$\mathrm{t} 0=\mathrm{psi} ;$
$\mathrm{t}=\mathrm{fmincon}(\mathrm{fun} 2, \mathrm{t} 0, \mathrm{~A}, \mathrm{~B}, \mathrm{Aeq}, \mathrm{beq}, \mathrm{lb}, \mathrm{ub}$, nonlcon, options);
\%print out the result
step $=$ j;
psi $=\mathrm{t}$;
end
\% mu, sigma, gamma optimization
for $\mathrm{i}=1: \mathrm{kW}$
lambdawx $=$ lambda - d $/ 2$;
chiwx $=$ chi+sum $(((-m u+X) /$ Sigma $) \cdot *(-m u+X), 2) ; \%$ this is nx1 matrix
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
j;
\%compute the density of W_i|x_i
chiwx $=\operatorname{chi}+\operatorname{sum}(((-m u+X) /$ Sigma $) \cdot *(-\operatorname{mu}+X), 2) ; \%$ this is nx1 matrix
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
\%compute $\mathrm{a}, \mathrm{b}$
$\mathrm{a}=\mathrm{sqrt}(\operatorname{chiwx} / \mathrm{psiw} \mathrm{x}) . *$ besselk(lambdawx+1, sqrt(chiwx$*$ psiwx))./ besselk(lambdawx, sqrt(chiwx*psiwx));\% this is nx1 matrix: E[W]
$b=\operatorname{sqrt}($ psiwx.$/ \operatorname{chiwx}) . *$ besselk (lambdawx -1 , sqrt (chiwx $*$ psiwx))./ besselk(lambdawx, sqrt(chiwx*psiwx));\% this is nx1 matrix: $\mathrm{E}\left[\mathrm{W}^{\wedge}(-1)\right]$
\%update mu, Sigma, gammas
gamma $=\operatorname{mean}(\mathrm{b} \cdot *(\mathrm{X}-\operatorname{mean}(\mathrm{X}))) /(1-\operatorname{mean}(\mathrm{b}) *$ mean $(\mathrm{a}))$;
$\mathrm{mu}=\operatorname{mean}(\mathrm{X})-\operatorname{mean}(\mathrm{a}) *$ gamma;
Sigma $=$ transpose $(X-m u) *(b . *(X-m u)) / n-$ mean $(a) *$ transpose $($ gamma) *gamma;
end

$$
\text { if iter }=\text { iterations }-1
$$

gamma_n_1 = gamma;

$$
\mathrm{mu}_{-\mathrm{n}} \mathrm{1}=\mathrm{mu}
$$

Sigma_n_1 = Sigma;
chi_n_1 $=$ chi $;$
psi_n_1 = psi;
lambda_n_1 = lambda;
end
if iter $=$ iterations
gamma_n = gamma;
mu_n $=\mathrm{mu}$;
Sigma_n $=$ Sigma;
chi_n $=$ chi;
psi_n $=$ psi;
lambda_n $=$ lambda;
end
end
\% error sanity checks
error_gamma $=$ norm $(($ gamma_n./gamma_n_1 -1$))$;
error_mu $=$ norm $(($ mu_n./mu_n_1 -1$))$;
error_Sigma $=$ norm $\left(\left(\operatorname{Sigma\_ n./Sigma\_ n\_ 1-1));~}\right.\right.$
error_chi $=$ norm $(($ chi_n./chi_n_1 -1$))$;
error_psi $=$ norm $(($ psi_n./ psi_n_1 - 1$))$;
error_lambda $=$ norm $\left(\left(\operatorname{lambda\_ n./lambda\_ n\_ 1-1));~}\right.\right.$
\%simulation of final distribution
$\mathrm{W}=$ gigrnd (lambda, psi, chi, nsimulations) ; \% generate W
$Z=\operatorname{mvnrnd}(z e r o s(1, d), ~ S i g m a, ~ n s i m u l a t i o n s) ; ~ \% ~ g e n e r a t e ~ Z ~ w i t h ~$
designed Sigma
$\mathrm{X} \operatorname{sim}=\mathrm{mu}+\mathrm{W} *$ gamma+sqrt(W).*Z; \% combine to give X
Xsimsum $=\operatorname{sum}(X \operatorname{sim}, 2) ; \%$ find rowsums to calculate portfolio
losses
sorted_lossesXsim $=$ sort (Xsimsum, 'descend') ; \% sort portfolio
losses
num_losses $=$ numel (sorted_lossesXsim) ; \% count portfolio losses
VaR_index $=$ floor $((1-$ confidence_level_VaR $) *$ num_losses $)+1 ; \%$
Calculate the index of the sorted losses that will be VaR
ES_index $=$ floor $((1-$ confidence_level_ES $) *$ num_losses $)+1 ; ~ \% ~$

Calculate the index of the sorted losses that will contribute to ES

175 VaRsumXsim $=$ sorted_lossesXsim(VaR_index)*dollarvalue; \% Use the index to extract $V a R$ from sorted losses
${ }_{176}$ ES1Xsim $=($ sorted_lossesXsim (1:ES_index-1)); \% Extract the loss tail

177 ESsumXsim $=$ mean(ES1Xsim) $*$ dollarvalue ; \% average of loss tail to calculate ES

## A. 2 EM algorithm: testing on W

1 \%This is the EM algorithm for estimating lambda, chi and psi while mu, Sigma, gamma are fixed.

2
3 format long;
4
5
6 \%Simulating X for data
7
$8 \mathrm{~d}=4 ; \%$ number of stocks
$9 \mathrm{n}=1000$; \% number of data points
10 \% Important: a larger $n$ allows to see the estimated parameters closer to true values!

11
12 gamma $=[0,1,4,-2]$;
${ }_{13} \mathrm{mu}=[2,1,-1,0]$;
${ }_{14} \mathrm{~A}=\left[\begin{array}{llllllllllllllll}2 & 0 & 0 & 1 ; & 0 & 1 & 0 & 0 ; & 0 & 0 & 4 & 1 ; & 1 & 0 & 1 & 4\end{array}\right] ;$

Sigma $=A /$ nthroot $(\operatorname{det}(A), d) ;$ \%for purpose of comparison, we adopt $\operatorname{det}($ Sigma $)=1$
chi $=2$;
psi $=0.5$;
lambda $=-1 ;$
$\mathrm{W}=$ gigrnd (lambda, psi, chi, n$) ;$ generate W
$Z=\operatorname{mvnrnd}(z e r o s(1, d), ~ S i g m a, ~ n) ; \%$ generate $Z$ with designed Sigma
$\mathrm{X}=\mathrm{mu}+\mathrm{W} *$ gamma+sqrt(W).$* \mathrm{Z}$; \% combine to give X
clear W;
clear Z;
\%initial values
chi $=1 ;$
psi $=1$;
\%settle initial value of lambda
fun $=@($ lambda $)$ besselk(lambda $+1,1)$-besselk(lambda, 1$) *$ nthroot $($ $\operatorname{det}(\operatorname{cov}(X)), d) ;$
lambda $=$ fzero (fun,-0.5$)$;
clear fun;
\%Excute the EM algorithm
$\mathrm{k}=1000 ;$ \% number of partial updates on Q 2
\% Important: a large k will allow us to see whether the
algorithm converges!
for $\mathrm{j}=1: \mathrm{k}$ \%work on Q 2 and partially update lambda, chi, psi for k
times
\%compute the density of W_i|x_i
lambdawx $=$ lambda - d $/ 2$;
chiwx $=\operatorname{chi}+\operatorname{sum}(((-\operatorname{mu}+\mathrm{X}) / \operatorname{Sigma}) \cdot *(-\operatorname{mu}+\mathrm{X}), 2) ; \%$ this is
nx1 matrix
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
\%compute $\mathrm{a}, \mathrm{b}, \mathrm{c}$
$\mathrm{a}=\mathrm{sqrt}($ chiwx $/ \mathrm{psiwx}) . *$ besselk(lambdawx +1 , sqrt (chiwx $*$
psiwx))./besselk(lambdawx, sqrt( $\operatorname{chiwx} * \operatorname{psiwx})) ; \%$ this
is nx1 matrix: $\mathrm{E}[\mathrm{W}]$
$\mathrm{b}=\operatorname{sqrt}(\mathrm{psiwx} . / \operatorname{chiwx}) \cdot *$ besselk(lambdawx-1, sqrt(chiwx $*$
psiwx))./besselk(lambdawx, sqrt(chiwx*psiwx));\% this
is nx1 matrix: $\mathrm{E}\left[\mathrm{W}^{\wedge}(-1)\right]$
$\mathrm{c}=\mathrm{zeros}(\mathrm{n}, 1) ; \%$ this is nx1 matrix: $\mathrm{E}[\ln \mathrm{W}]$
for $i=1: 1: n$
fun1 $=@(\mathrm{w}) \log (\mathrm{w}) . * \mathrm{w} .{ }^{\wedge}(\operatorname{lambdawx}-1) . * \exp (-0.5 * \mathrm{psiw} \mathrm{x} *$
$\mathrm{w}) . * \exp \left(-0.5 * \operatorname{chiwx}(\mathrm{i}, 1) * \mathrm{w} .{ }^{\wedge}(-1)\right)$;
$c(i, 1)=0.5 *(\operatorname{psiwx} / \operatorname{chiwx}(\mathrm{i}, 1))^{\wedge}(0.5 * \operatorname{lambdawx}) /$
besselk (lambdawx, $\operatorname{sqrt}(\operatorname{chiwx}(\mathrm{i}, 1) * \operatorname{psiwx})) *$ integral
(fun1, 0, Inf);
clear fun1;
end
\%update lambda, chi, psi
options $=$ optimoptions('fmincon', 'Algorithm ', 'sqp');
fun2 $=@(\mathrm{t}) \mathrm{t}(1) * \log (\mathrm{t}(2))-\mathrm{t}(1) * \log (\mathrm{t}(3))+2 * \log (\operatorname{besselk}($
$\mathrm{t}(1), \operatorname{sqrt}(\mathrm{t}(2) * \mathrm{t}(3)))+\mathrm{t}(3) * \operatorname{mean}(\mathrm{a})+\mathrm{t}(2) * \operatorname{mean}(\mathrm{~b})-2 * \mathrm{t}$
(1) *mean (c);
$\mathrm{A}=[0,-1,0 ; 0,0,-1] ;$
$\mathrm{B}=[0 ; 0] ;$
Aeq $=[] ;$
beq $=[]$;
$\mathrm{lb}=[] ;$
$\mathrm{ub}=[]$;
nonlcon $=[] ;$
$\mathrm{t} 0=[$ lambda, chi, psi];
$\mathrm{t}=\mathrm{fmincon}(\mathrm{fun} 2, \mathrm{t} 0, \mathrm{~A}, \mathrm{~B}, \mathrm{Aeq}, \mathrm{beq}, \mathrm{lb}, \mathrm{ub}$, nonlcon, options);
clear fun2;
clear a;
clear b;
clear c;
clear lambdawx;
clear chiwx;
clear psiwx;
\%print out the result
fprintf (2, ' $\backslash$ nstep $\left.\backslash n^{\prime}\right)$
step $=$ j

## A. 3 EM algorithm: testing with W as fixed

\%This is the EM algorithm for estimating mu, Sigma, gamma while lambda, chi and psi are fixed.

2
3

4
5 \%Simulating X for data

6
${ }_{7} \mathrm{~d}=4$; \%number of stocks
$8 \mathrm{n}=100000$ \% number of data points.
9 Important: a larger n allows to see the estimated parameters closer to true values!

10
11 gamma $=[0,4,0,4]$;

```
\(\mathrm{mu}=[2,10,-1,0] ;\)
\(\mathrm{A}=\left[\begin{array}{cccccccccccccccc}2 & 0 & 0 & 1 ; & 0 & 1 & 0 & 0 ; & 0 & 0 & 4 & 1 ; & 1 & 0 & 1 & 4\end{array}\right]\);
Sigma \(=A /\) nthroot \((\operatorname{det}(A), d) ;\) for purpose of comparison, we
    adopt \(\operatorname{det}(\) Sigma \()=1\)
    chi=2;
    psi=5;
    lambda \(=-3\);
    \%It seems some combination of values of lambda, chi, psi can
        cause problems for the algorithm.
\(\mathrm{W}=\) gigrnd (lambda, psi, chi, n\()\); \% generate W
```



```
        Sigma
```

$\mathrm{X}=\mathrm{mu}+\mathrm{W} *$ gamma+sqrt(W).*Z; \% combine to give X
\% initial values
gamma $=\operatorname{zeros}(1, \mathrm{~d})$;
$\mathrm{mu}=\operatorname{mean}(\mathrm{X}) ;$
Sigma $=\operatorname{cov}(X) / n t h r o o t(\operatorname{det}(\operatorname{cov}(X)), d) ;$
\%small influence on initial values
\%Excute the EM algorithm
\%compute the density of W_i|x_i
lambdawx $=$ lambda - d $/ 2$;
$\mathrm{k}=10000 ; \%$ number of partial updates on Q1;
\% Important: a large $k$ will allow us to see whether the
algorithm converges!
for $j=1: k$ \% work on Q_1 and partially update mu, Sigma, gamma
for $k$ times
fprintf(2, ' $\left.\operatorname{nstep} \backslash n^{\prime}\right)$
j
\%compute the density of W_i|x_i
chiwx $=\operatorname{chi}+\operatorname{sum}(((-m u+X) /$ Sigma $) \cdot *(-m u+X), 2) ; \%$ this is
nx1 matrix
psiwx $=$ psi+gamma/Sigma*transpose (gamma) ;
\%compute $\mathrm{a}, \mathrm{b}$
$\mathrm{a}=\mathrm{sqrt}($ chiwx $/ \mathrm{psiwx}) . *$ besselk(lambdawx +1 , sqrt (chiwx $*$
psiwx))./besselk(lambdawx, sqrt(chiwx*psiwx));\% this
is nx1 matrix: $\mathrm{E}[\mathrm{W}]$
$b=\operatorname{sqrt}($ psiwx. $/$ chiwx $) . *$ besselk (lambdawx -1, sqrt (chiwx $*$
psiwx))./besselk(lambdawx, sqrt(chiwx*psiwx));\% this
is nx1 matrix: $\mathrm{E}\left[\mathrm{W}^{\wedge}(-1)\right]$
\%update mu, Sigma, gammas
gamma $=\operatorname{mean}(\mathrm{b} . *(\mathrm{X}-\operatorname{mean}(\mathrm{X}))) /(1-\operatorname{mean}(\mathrm{b}) * \operatorname{mean}(\mathrm{a}))$
$\mathrm{mu}=\operatorname{mean}(\mathrm{X})-$ mean $(\mathrm{a}) *$ gamma
Sigma $=\operatorname{transpose}(\mathrm{X}-\mathrm{mu}) *(\mathrm{~b} \cdot *(\mathrm{X}-\mathrm{mu})) / \mathrm{n}-$ mean $(\mathrm{a}) * \operatorname{transpose}($
gamma) *gamma;
Sigma $=$ Sigma $/$ nthroot $(\operatorname{det}(\operatorname{Sigma}), d)$
\%print true values of parameters for comparison
fprintf(2, '\ntrue values $\left.\backslash n^{\prime}\right)$
gamma $=[0,4,0,4]$
$\mathrm{mu}=[2,10,-1,0]$
$\mathrm{A}=\left[\begin{array}{llllllllllllllll}2 & 0 & 0 & 1 ; & 0 & 1 & 0 & 0 ; & 0 & 0 & 4 & 1 ; & 1 & 0 & 1 & 4\end{array}\right] ;$
Sigma $=A / n t h r o o t(\operatorname{det}(A), d)$ \%for purpose of comparison, we
adopt $\operatorname{det}(\operatorname{Sigma})=1 \mathrm{~s}$

## Appendix B

## Codes for Chapter 3

## B. 1 Code for Marginals and Copulas

close all
pricematrix $=$ [AAPL.AdjClose BAC.AdjClose F.AdjClose SU.AdjClose
]; \% load price matrix

```
[n,d] = size(pricematrix); % calculate dimensions of price
``` matrix
```

returnmatrix = zeros(n-1,d); % initializing return matrix

```
names \(=\{\) 'AAPL', 'BAC', 'F', 'SU' \(\} ; \%\) names of stocks
nsimulations \(=10000\); number of simulations to run
confidence_level_VaR \(=0.99 ; \%\) confidence level for VaR
confidence_level_ES \(=0.975\); \% confidence level for ES
dollaramount \(=10000\); dollar amount for each stock
theta \(=3 ; \%\) for Gumbel Copula
for \(\mathrm{i}=1: \mathrm{n}-1 \%\) loop over time
    returnmatrix \((\mathrm{i},:)=-\log (\) pricematrix \((\mathrm{i}+1,:) . /\) pricematrix \((\mathrm{i}\)
end
meanvec \(=\operatorname{mean}(\) returnmatrix \() *\) dollaramount; \(\%\) find mean of dollar losses
stdvec \(=\operatorname{std}(\) returnmatrix \() *\) dollaramount ; \% find standard deviation of dollar losses
figure
counter \(=0 ; \%\) initialized counter
for \(\mathrm{i}=1: \mathrm{d} \%\) loop over stocks
\(\mathrm{x}=[-3 * \operatorname{stdvec}(\mathrm{i})+\) meanvec (i) \(: 0.001: 3 * \operatorname{stdvec}(\mathrm{i})+\operatorname{meanvec}(\mathrm{i})]\); \% marginals
norm \(=\) normpdf( \(x\), meanvec (i), stdvec (i)) ; \% calculate density
counter \(=\) counter \(+1 ; \%\) increase counter
subplot(2, 2 , counter) \% initialize subplot
plot(x, norm) \% plot marginal
xlabel('Returns');
ylabel('Density');
mytitleText \(=[\) 'Normal Distribution plot for ', names(i)];
title (mytitleText, 'Interpreter', 'tex' ) ;
\(\mathrm{y}=\operatorname{meanvec}(\mathrm{i})+\operatorname{randn}(\) nsimulations, 1\() * \operatorname{stdvec}(\mathrm{i}) ; \%\) simulate marginals
sorted_losses \(=\operatorname{sort}(y, ' d e s c e n d ') ; \%\) sort marginal losses
num_losses \(=\) numel (sorted_losses) ; \% count marginal losses
VaR_index \(=\) floor \(((1-\) confidence_level_VaR \() *\) num_losses \()+1 ; \%\) Calculate the index of the sorted losses that will be VaR
ES_index \(=\) floor \(((1-\) confidence_level_ES \() *\) num_losses \()+1 ; ~ \% ~\)
Calculate the index of the sorted losses that will
```

        contribute to ES
    VaR(i) = sorted_losses(VaR_index); % Use the index to
        extract VaR from sorted losses
    ES1 =(sorted_losses(1:ES_index - 1)); % Extract the loss tail
    ES(i) = mean(ES1); % average of loss tail to calculate ES
    end
correlationmatrixZ = corr(returnmatrix ,'type','Spearman'); %
correlation matrix for Gauss Copula
correlationmatrixT = sin (0.5* pi*corr(returnmatrix, 'type','
Kendall')); % correlation matrix for t Copula
gamma = (cos(pi/(2*theta)) )^theta; % gamma value for Gumbel
Copula
Z = mvnrnd(zeros(1,d), correlationmatrixZ,nsimulations); %
simulate multivariate random normal for Gauss copula
T = mvtrnd(correlationmatrixT , 5, nsimulations); % simulate
multivariate random t for t copula
U = rand(nsimulations,4);
Gu = [];
while length(Gu) < nsimulations
Gurand = random('stable', 1/ theta, , ,gamma, 0, 1); %simulate
stable random variates for Gumbel copula
if Gurand > 0
Gu = [Gu ; Gurand ];
end
end

```
```

simulatedmatrixZ = normcdf(Z); % evaluate normal cumulative
distribution function to generate Gauss copula
simulatedmatrixT = tcdf(T,5); % evaluate t cumulative
distribution function to generate t copula
simulatedmatrixGu = exp(-((-log(U)./Gu).^(1/theta))); % get
generating vector for Gumbel copula
Xz=[]; % initialize loss matrix for Gauss copula
Xt = []; % initialize loss matrix for t copula
Xgu = []; % initialize loss matrix for Gumbel copula
for i = 1:d
Xz = [Xz norminv(simulatedmatrixZ(:, i),meanvec(i), stdvec(i))
]; % plug in marginals to get losses for Gauss copula
Xt = [Xt norminv(simulatedmatrixT (:, i) ,meanvec(i),stdvec(i))
]; % plug in marginals to get losses for t copula

```
    \(\mathrm{Xgu}=[\mathrm{Xgu}\) norminv\((\operatorname{simulatedmatrixGu(:,i),meanvec(i),~stdvec~(~}\)
        i)) \(]\); \(\%\) plug in marginals to get losses for Gumbel copula
    end
    Xzsum \(=\operatorname{sum}(\mathrm{Xz}, 2) ; \%\) find rowsums to calculate portfolio losses
        for Gauss copula
    Xtsum \(=\operatorname{sum}(\mathrm{Xt}, 2) ; \%\) find rowsums to calculate portfolio losses
        for \(t\) copula
    Xgusum \(=\operatorname{sum}(\mathrm{Xgu}, 2) ; \%\) find rowsums to calculate portfolio
        losses for Gumbel copula
sorted_lossesz \(=\) sort (Xzsum, 'descend') ; \% sort portfolio losses
    for Gauss copula
\({ }_{71}\) sorted_lossesgu \(=\) sort (Xgusum, 'descend'); \% sort portfolio losses for Gumbel copula

2 2 num_losses \(=\) numel (sorted_lossesz) ; \% count portfolio losses
\({ }_{73}\) VaR_index \(=\) floor \(((1-\) confidence_level_VaR \() *\) num_losses \()+1\); \% Calculate the index of the sorted losses that will be VaR

ES_index \(=\) floor \(((1-\) confidence_level_ES \() *\) num_losses \()+1 ; ~ \% ~\)
Calculate the index of the sorted losses that will contribute to ES

VaRsumz \(=\) sorted_lossesz (VaR_index) ; \% Use the index to extract VaR from sorted losses for Gauss copula

ES1z \(=\left(\right.\) sorted_lossesz \(\left.\left(1: E S \_i n d e x-1\right)\right) ;\) Extract the loss tail for Gauss copula

ESsumz \(=\) mean(ES1z); \% average of loss tail to calculate ES for Gauss copula

VaRsumt \(=\) sorted_lossest (VaR_index) ; \% Use the index to extract VaR from sorted losses for t copula

ES1t \(=(\) sorted_lossest (1:ES_index -1\())\); Extract the loss tail for \(t\) copula

ESsumt \(=\) mean \((E S 1 t) ;\) average of loss tail to calculate ES for t copula

VaRsumgu \(=\) sorted_lossesgu(VaR_index); \% Use the index to extract VaR from sorted losses for Gumbel copula

ES1gu \(=(\) sorted_lossesgu(1:ES_index -1\())\); \% Extract the loss tail for Gumbel copula
B.1. COPULA CODES
for Gumbel copula

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