Model uncertainty and VaR aggregation

Paul Embrechts^a, Giovanni Puccetti^{b,*}, Ludger Rüschendorf^c

^aSenior SFI Professor, RiskLab, Department of Mathematics, ETH Zurich, 8092 Zurich, Switzerland
^bDepartment of Mathematics for Decision Theory, University of Firenze, 50127 Firenze, Italy
^cDepartment of Mathematical Stochastics, University of Freiburg, 79104 Freiburg, Germany

Abstract

Despite well-known shortcomings as a risk measure, Value-at-Risk (VaR) is still the industry and regulatory standard for the calculation of risk capital in banking and insurance. This paper is concerned with the increase of the VaR for a portfolio position as a function of different dependence scenarios on the factors of the portfolio. Besides summarizing the most relevant analytical bounds, including a discussion of their sharpness, we introduce a new numerical algorithm which allows for the computation of reliable (sharp) bounds for the VaR of high-dimensional portfolios with dimensions *d* in the several hundreds. We show that additional positive dependence information (like positive correlation) will typically not improve the upper bound substantially. In contrast higher order marginal information on the model, when available, may lead to strongly improved bounds. Several examples of practical relevance show how explicit VaR bounds can be obtained. These bounds can be interpreted as a measure of model uncertainty induced by possible dependence scenarios.

Keywords: Copula, Fréchet class, Model Uncertainty, Operational Risk, Positive Dependence, Rearrangement Algorithm, Risk Aggregation, Value-at-Risk, VaR-bounds.

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1. Introduction

Since the early nineties, Value-at-Risk (VaR) has established itself as a (if not the) key metric for the calculation of regulatory capital within the financial industry. Furthermore, VaR is increasingly used as a risk management constraint within portfolio optimization. Whereas books like Jorion (2006) prize VaR as the industry standard, numerous papers have pointed out many of the (most obvious) shortcomings of VaR as a risk measure; see for instance McNeil et al. (2005) and the references therein, but also the recent Basel Committee on Banking Supervision (2012), already refereed to as Basel 3.5. A very informative overview on the use of VaR technology within the banking industry is Pérignon and Smith (2010). As so often, a middle-of-the-road point of view is advisable: there is no doubt that the construction and understanding of the P&L of a bank's trading book is of the utmost importance. The latter includes the availability of data warehouses, independent pricing tools, a complete risk factor mapping, proper corporate governance structures around the risk committee, etc... In that sense, VaR, as a number, is just the peak of the risk management iceberg. Nonetheless, once the number leaves the IT system of the CRO, all too often it starts a life of its own and one often forgets the numerous warnings about its proper interpretation. Moreover, once several VaRs are involved, the temptation is

To set the scene, consider the calculation of the Value-at-Risk (VaR) at a confidence level α for an aggregate loss random variable L having the form

$$L = \sum_{i=1}^{d} L_i,$$

where L_1, \ldots, L_d , in the case of Operational Risk, correspond to the loss random variables for given business lines or risk types, over a fixed time period T. The VaR of the aggregate position L, calculated at a probability level $\alpha \in (0,1)$, is the α -quantile of its distribution, defined as

$$VaR_{\alpha}(L) = F_L^{-1}(\alpha) = \inf\{x \in \mathbb{R} : F_L(x) \ge \alpha\},\tag{1}$$

where $F_L(x) = P(L \le x)$ is the distribution function of L. Here, as a statistical quantity and for α typically close to 1, $\operatorname{VaR}_{\alpha}(L)$ is a measure of extreme loss, i.e. $P(L > \operatorname{VaR}_{\alpha}(L)) = 1 - \alpha$ is small

The current regulatory framework for banking supervision, referred to as Basel II (becoming Basel III), allows large inter-

Email addresses: embrechts@math.ethz.ch (Paul Embrechts), giovanni.puccetti@unifi.it (Giovanni Puccetti), ruschen@stochastik.uni-freiburg.de (Ludger Rüschendorf)

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there to calculate functions of them (like adding) forgetting the considerable model uncertainty underlying such constructions; see Basel Committee on Banking Supervision (2010) for a regulatory overview on risk aggregation. A typical such example is to be found in the realm of Operational Risk as defined under Basel II and III. Throughout the paper we will use the latter as a motivating example and consider the organization of an Operational Risk database in business lines and risk types; for a background to this and for further references, see for instance McNeil et al. (2005, Chapter 10).

^{*}Corresponding author

national banks to come up with internal models for the calculation of risk capital. For Operational Risk, under the so-called Loss Distribution Approach (LDA) within Basel II, financial institutions are given full freedom concerning the stochastic modeling assumptions used. The resulting risk capital must correspond to a 99.9%-quantile of the aggregated loss data over the period of a year; we leave out the specific details concerning internal, external and expert opinion data as they are less relevant for the results presented in this paper. Using the notation introduced above, the risk capital for the aggregate position L is typically calculated as $VaR_{0.999}(L)$. Concerning interdependence of risks, no specific rules are given beyond the statement that explicit and implicit correlation assumptions between loss random variables used have to be plausible and need to be well founded; in the case of Operational Risk see Cope and Antonini (2008) and Cope et al. (2009). For the sequel of this paper, we leave out statistical (parameter) uncertainty.

In order to calculate $VaR_{\alpha}(L)$, one needs a joint model for the random vector $(L_1, \ldots, L_d)'$. This would require a d-variate dataset for the past occurred losses, which often is not available. Typically, only the marginal distribution functions F_i of L_i are known or statistically estimated, while the dependence structure between the L_i 's is either completely or partially unknown.

In standard practice, the total capital charge C to be allocated is derived from the addition of the VaRs at probability level $\alpha = 0.999$ for the marginal random losses L_i , namely

$$\operatorname{VaR}_{\alpha}^{+}(L) = \sum_{i=1}^{d} \operatorname{VaR}_{\alpha}(L_{i}) = \sum_{i=1}^{d} F_{i}^{-1}(\alpha).$$

Indeed, industry typically reports

$$C = \delta \operatorname{VaR}_{\alpha}^{+}(L), \quad 0 < \delta \le 1; \tag{2}$$

the value of δ is often in the range (0.7,0.9) and reflects socalled diversification effects. A capital charge based on (2) would imply a subadditive regime for VaR, i.e.

$$VaR_{\alpha}(L) = VaR_{\alpha} \left(\sum_{i=1}^{d} L_i \right) \le \sum_{i=1}^{d} VaR_{\alpha}(L_i) = VaR_{\alpha}^{+}(L).$$
 (3)

The case $\delta=1$ (no diversification) in (2) can be mathematically justified by the assumption of perfect positive dependence (which implies maximal correlation) among marginal risks. Indeed, under this so-called *comonotonic dependence* scenario, $VaR_{\alpha}(L) = VaR_{\alpha}^{+}(L)$; see McNeil et al. (2005, Proposition 6.15). Practitioners criticize this assumption as not being realistic, and remark that random losses are not perfectly correlated in view of their heterogeneous nature. Though the $\delta=1$ maximal-correlation scenario is often considered as highly conservative, the inequality in (3) is typically violated for very heavy-tailed, very skewed losses, possibly exhibiting special dependencies, situations which are no doubt present in Operational Risk data; see for instance Moscadelli (2004), Panjer (2006) and Shevchenko (2011).

Based on the above example from the capital charge calculation of Operational Risk it is clear that there exists considerable model uncertainty underlying the *diversification* factor

 δ , which for practically relevant models may take values well above the additive case $\delta = 1$. It is exactly this kind of model uncertainty that the present paper addresses.

Recently, a number of numerical and analytical techniques have been developed in order to calculate conservative values for $VaR_{\alpha}(L)$ under different dependence assumptions regarding the loss random variables L_i . In this paper we describe these methodologies and give insight in the worst-case dependence structure (copula) describing the worst-VaR scenario. A main contribution of this paper is the introduction of an algorithm which allows to calculate sharp dependence bounds for the VaR of high-dimensional portfolios allowing us to consider inhomogeneous portfolios with dimension d in the several hundreds. We also discuss the possible influence of additional positive dependence information as well as information on higher dimensional sub-vectors of marginals. The main message coming from our paper is that currently a whole toolkit of analytical and numerical techniques is available to better understand the aggregation and diversification properties of non-coherent risk measures such as Value-at-Risk.

We very much hope that our paper is both accessible to the academic researcher as well as to the more quantitative practitioner. With this goal in mind, we have strived at keeping the technical details to a minimum, stressing more the algorithmic numerical aspects of the results discussed. Of course, we will direct the reader interested in more mathematical details to the relevant research papers. We strongly believe that the results and techniques summarized are sufficiently novel and will benefit the wider financial industry.

With financial/actuarial applications in mind, and without loss of generality, in almost all the examples contained in the paper we use power law models for the marginal distributions of the risks such as the Pareto distribution. In particular, we often use a Pareto distribution with tail parameter $\theta=2$ in order to represent marginal risks with finite mean but infinite variance. This choice is pedagogical and does not affect the computational properties of the methodologies discussed.

In Section 2, we study the case where the marginal distribution functions F_i of L_i are fixed while the dependence structure (copula) between the L_i 's is completely unknown. In the homogeneous case where the risk factors L_i are identically distributed, a simple analytical formula allows to compute the worstpossible VaR for portfolios of arbitrary dimensions when the marginal distributions F_i are continuous. For *inhomogeneous* portfolios having arbitrary marginals, a new numerical algorithm allows to compute best- and worst-possible VaR values up to dimensions $d \approx 600$. Under the restriction of the dependence structure to positive dependence, possible improvements of the bounds are discussed in Section 3. Finally, in Section 4, we consider a more general case where extra information is known about sub-vectors of the marginal risks. In the Subsections 1.1-1.4 below we first gather some definitions, notation and basic methodological tools, together with some key references.

1.1. Fréchet classes

Denote $L = (L_1, \dots, L_d)'$. The Value-at-Risk for the aggregate position $L = L_1 + \dots + L_d$ is certainly not uniquely determined by the marginal distributions F_1, \dots, F_d of the risks L_i . In fact, there exist infinitely many joint distributions on \mathbb{R}^d which are consistent with the choice of the marginals F_1, \dots, F_d . We denote by $\mathfrak{F}(F_1, \dots, F_d)$ the *Fréchet class* of all the possible joint distributions F_L on \mathbb{R}^d having the given marginals F_1, \dots, F_d . For $\alpha \in (0, 1)$, upper and lower bounds for the Value-at-Risk of L are then defined as

$$\overline{\mathrm{VaR}}_{\alpha}(L) = \sup \left\{ \mathrm{VaR}_{\alpha}(L_1 + \dots + L_d) : F_L \in \mathfrak{F}(F_1, \dots, F_d) \right\},$$
(4a)

$$\underline{\operatorname{VaR}}_{\alpha}(L) = \inf \left\{ \operatorname{VaR}_{\alpha}(L_1 + \dots + L_d) : F_L \in \mathfrak{F}(F_1, \dots, F_d) \right\}. \tag{4b}$$

The above definitions directly imply the VaR range for L given by

$$\underline{\operatorname{VaR}}_{\alpha}(L) \le \operatorname{VaR}_{\alpha}(L_1 + \dots + L_d) \le \overline{\operatorname{VaR}}_{\alpha}(L).$$
 (5)

We refer to the bounds $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\alpha}(L)$ as the worst-possible and, respectively, the best-possible VaR for the position L, at the probability level α . When attained, the upper and lower bounds in (4) are sharp (best-possible): they *cannot be improved* if further dependence information on $(L_1, \ldots, L_d)'$ is not available. We call any joint model for $(L_1^*, \ldots, L_d^*)'$ with prescribed marginals F_1, \ldots, F_d such that

$$\overline{\operatorname{VaR}}_{\alpha}(L) = \operatorname{VaR}_{\alpha}(L_1^* + \dots + L_d^*)$$

a worst-case dependence or worst-case coupling. Analogously, any joint model for $(L_1^*, \ldots, L_d^*)'$ with the prescribed marginals such that

$$VaR_{\alpha}(L) = VaR_{\alpha}(L_1^* + \cdots + L_d^*)$$

is a *best-case dependence* or *best-case coupling*. Of course, the choice of wording *best* versus *worst* is arbitrarily and depends on the specific application at hand. Problems related to (4) with moment information have always been relevant in actuarial mathematics. One of the early contributors was De Vylder (1996).

1.2. Copulas

To make this paper self-contained, we give a brief introduction to some copula concepts that we will need in the following. The reader not familiar with the theory of copulas is referred to Nelsen (2006), McNeil et al. (2005, Chapter 5) and Durante and Sempi (2010).

A copula C is a d-dimensional distribution function (df) on $[0,1]^d$ with uniform marginals. Given a copula C and d univariate marginals F_1, \ldots, F_d , one can always define a df F on \mathbb{R}^d having these marginals by

$$F(x_1, \dots, x_d) = C(F_1(x_1), \dots, F_d(x_d)), \ x_1, \dots, x_d \in \mathbb{R}.$$
 (6)

Sklar's Theorem states conversely that we can always find a copula C coupling the marginals F_i of a fixed joint distribution F through the above expression (6). For continuous marginal

dfs, this copula is unique. Hence Sklar's Theorem states that the copula C of a multivariate distribution F contains all the dependence information of F.

A first example of a copula is the independence copula

$$\Pi(u_1,\ldots,u_d)=\Pi_{i=1}^d u_i.$$

The name of this copula derives from the fact that the risk vector $(L_1, \ldots, L_d)'$ has copula Π if and only if its marginal risks L_i are independent. Under independence among the marginal risks, (6) reads as

$$F(x_1, \dots, x_d) = \Pi(F_1(x_1), \dots, F_d(x_d)) = F_1(x_1) \cdot \dots \cdot F_d(x_d).$$

Any copula C satisfies the so-called Fréchet bounds

$$\max \left\{ \sum_{i=1}^{d} u_i - d + 1, 0 \right\} \le C(u_1, \dots, u_d) \le \min\{u_1, \dots, u_d\},$$

for all $u_1, \ldots, u_d \in [0, 1]$. The upper Fréchet bound

$$M(u_1,\ldots,u_d)=\min\{u_1,\ldots,u_d\}$$

is the so-called *comonotonic* copula, which represents perfect positive dependence among the risks. In fact, a risk vector $(L_1, \ldots, L_d)'$ has copula M if and only if its marginal risks are all almost surely (a.s.) increasing functions of a common random factor. For a detailed discussion of the concept of comonotonicity within quantitative risk management we refer to Dhaene et al. (2002) and Dhaene et al. (2006); see also McNeil et al. (2005, Section 6.2.2). The lower Fréchet bound

$$W(u_1, \ldots, u_d) = [u_1 + \cdots + u_d - d + 1]^+$$

is sharp. However, it is a well-defined copula only in dimension d=2. In this case, it is called the *countermonotonic* copula and represents perfect negative dependence between two risks. A risk vector $(L_1, L_2)'$ has copula W if and only if its marginal risks are a.s. decreasing functions of each other.

The upper and lower Fréchet bounds play the role of optimal copulas in many optimization problems of interest in quantitative risk management. For instance it is well known that the maximal variance for the sum of risks with given marginals is attained when the risks are comonotonic, that is when they have copula C = M. Analogously, the minimal variance for the sum of two risks with given marginals is attained when they are countermonotonic, C = W. These results derive from the classical Hoeffding-Fréchet bounds and can be seen as particular cases of a more general ordering theorem; see Corollary 3 in Rüschendorf (1983).

1.3. Worst and best VaR

Sklar's Theorem allows us to reformulate (4) as optimization problems over C_d , the set of all d-dimensional copulas:

$$\overline{\mathrm{VaR}}_{\alpha}(L) = \sup \left\{ \mathrm{VaR}_{\alpha}(L_{1}^{C} + \dots + L_{d}^{C}) : C \in C_{d} \right\}, \tag{7a}$$

$$\underline{\operatorname{VaR}}_{\alpha}(L) = \inf \left\{ \operatorname{VaR}_{\alpha}(L_{1}^{C} + \dots + L_{d}^{C}) : C \in C_{d} \right\}. \tag{7b}$$

Here the vector $(L_1^C, \ldots, L_d^C)'$ has the same marginal distributions as (L_1, \ldots, L_d) and copula C. In general, it is difficult to evaluate the bounds in (4) or in (7) in explicit form especially when one has to deal with $d \geq 3$ risks. This is related to the fact that in general Value-at-Risk is *non-subadditive*. As a consequence, the comonotonic copula M is in general *not* a solution to the <u>problem $\overline{\mathrm{VaR}}_{\alpha}(L)$ in (7a)</u>. Equivalently, the worst-VaR value $\overline{\mathrm{VaR}}_{\alpha}(L)$ in (4) is *not* attained when all the risks are perfectly positively dependent. Analogously, the countermonotonic copula W is in general *not* a solution to the problem $\overline{\mathrm{VaR}}_{\alpha}(L)$ in (7b) for d=2.

As already stated above, in the comonotonic case C = M we have that

$$\operatorname{VaR}_{\alpha}^{+}(L) = \operatorname{VaR}_{\alpha}(L_{1}^{M} + \dots + L_{d}^{M}) = \sum_{i=1}^{d} \operatorname{VaR}_{\alpha}(L_{i}) = \sum_{i=1}^{d} F_{i}^{-1}(\alpha).$$
(8)

It is not difficult to provide examples of interest in quantitative risk management where, for a copula C, necessarily $C \neq M$, we have that

$$\operatorname{VaR}_{\alpha}(L_{1}^{C} + \cdots + L_{d}^{C}) > \sum_{i=1}^{d} \operatorname{VaR}_{\alpha}(L_{i}).$$

For instance, in the presence of infinite-mean marginal distributions, it is typical to have

$$\operatorname{VaR}_{\alpha}(L_{1}^{\Pi} + \dots + L_{d}^{\Pi}) > \sum_{i=1}^{d} \operatorname{VaR}_{\alpha}(L_{i})$$

for sufficiently high levels of α ; see Embrechts and Puccetti (2010b, Section 5.3), Mainik and Rüschendorf (2010), Mainik and Embrechts (2012) and the numerous references therein.

1.4. Complete mixability

When dealing with extremal values for Value-at-Risk, the ideas of perfect positive and negative dependence as represented by the Fréchet bounds M and W can be deceiving. Handling non-subadditive risk measures requires the knowledge of alternative dependence concepts; *complete mixability* turns out to be such a concept. It turns out to be highly useful towards the calculation of VaR bounds.

Definition 1. A distribution function F on \mathbb{R} is d-completely mixable (d-CM) if there exist d random variables X_1, \ldots, X_d identically distributed as F such that

$$P(X_1 + \dots + X_d = c) = 1,$$
 (9)

for some constant $c \in \mathbb{R}$. Any vector $(X_1, \ldots, X_d)'$ satisfying (9) with $X_i \sim F, 1 \leq i \leq d$, is called a d-complete mix. If F has finite first moment μ , then $c = \mu d$.

Complete mixability is a concept of negative dependence. In dimension d=2 complete mixability implies countermonotonicity. Indeed, a risk vector $(L_1, L_2)'$ is a 2-complete mix if and only if $L_1 = k - L_2$ a.s, and this implies that its copula is the

lower Fréchet bound W (the converse however does not hold). In higher dimensions $d \geq 3$ a completely mixable dependence structure minimizes the variance of the sum of risks with given marginal distributions. In fact, a risk vector $(L_1, \ldots, L_d)'$ with identically distributed marginals is a d-complete mix if and only if the variance of the sum of its components is equal to zero. Not all univariate distributions F are d-CM. As an example, it is sufficient to take F as the two-point distribution giving probability mass p > 0 to x = 0 and (1 - p) to x = 1. Since the only way to make $L_1 + L_2$ a constant is to choose $L_2 = 1 - L_1$, F is not 2-CM for $p \neq 1/2$.

The structure of dependence (copula) representing complete mixability is not so intuitive and, at the moment, does not have an easy mathematical formulation like in the case of the Fréchet bounds. We illustrate this with a discrete example. We choose F to give mass 1/5 to any of the first five integers. A 3-complete mix of F can be represented by the following matrix, in which any row is to be seen as a vector in \mathbb{R}^3 having probability mass 1/5:

$$\left[\begin{array}{cccc}
1 & 5 & 3 \\
2 & 3 & 4 \\
3 & 1 & 5 \\
4 & 4 & 1 \\
5 & 2 & 2
\end{array}\right].$$

Since the sum of each row in the above matrix is equal to k = 9 (note that the mean of F is equal to 3), F turns out to be 3-completely mixable. It is useful to compare the above matrix with the one representing comonotonicity among three F-distributed risks:

$$\left[\begin{array}{cccc}
1 & 1 & 1 \\
2 & 2 & 2 \\
3 & 3 & 3 \\
4 & 4 & 4 \\
5 & 5 & 5
\end{array}\right].$$

In this latter case, the variance of the row-wise sums is maximized. Some other examples of completely mixable distributions, as well as an insight into the theory of complete mixability, are given in Rüschendorf and Uckelmann (2002), Wang and Wang (2011) and Puccetti et al. (2012). The concept of complete mixability will play an important role in the optimization problems (7) in the homogeneous case where the L_i 's are identically distributed.

2. Computing the VaR range with given marginal information

In this section, we consider the case when the risk vector $(L_1, \ldots, L_d)'$ has given marginal distribution functions F_1, \ldots, F_d while its dependence structure is completely unknown. Recently, some new numerical and analytical tools have been developed to calculate the VaR range in (5) under these assumptions. First, we study the *homogeneous* case where the marginal risks are all identically distributed. Then, we will consider the more general *inhomogeneous* framework in which the marginal distributions are allowed to differ.

2.1. Identically distributed marginals

Throughout this section we assume that the marginal risks L_i are all identically distributed as F, that is $F_1 = \cdots = F_d = F$. In the case d = 2, the calculation of the sharp VaR bounds in (4) reduces to a simple formula if F satisfies some regularity conditions.

Proposition 2. In the case d=2 with $F_1=F_2=F$, let F be a continuous distribution concentrated on $[0,+\infty)$ with an ultimately decreasing density on $(\overline{x}_F,+\infty)$, for some $\overline{x}_F \geq 0$. Then

$$\underline{\operatorname{VaR}}_{\alpha}(L) = F^{-1}(\alpha) \quad and \quad \overline{\operatorname{VaR}}_{\alpha}(L) = 2F^{-1}\left(\frac{1+\alpha}{2}\right), (10)$$

for all $\alpha \in [F(\overline{x}_F), 1)$.

Remark 3. 1. If $\bar{x}_F = 0$, e.g. in the case F is Pareto distributed, that is

$$F(x) = 1 - (1+x)^{-\theta}, x > 0,$$
(11)

for some tail parameter $\theta > 0$, the sharp bounds in (10) hold for any level of probability $\alpha \in (0, 1)$.

2. For d=2, the sharp bounds $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\alpha}(L)$ are known for *any* type of marginal distributions F_1, F_2 . The slightly more complicated formulas to compute the bounds in the general case are given in Rüschendorf (1982, Proposition 1).

For a given α , a worst-case dependence vector (L_1^*, L_2^*) such that $VaR_{\alpha}(L_1^* + L_2^*) = \overline{VaR}_{\alpha}(L)$ is given by

$$\begin{cases} L_2^* = L_1^* & \text{a.s., when } L_1 < F^{-1}(\alpha), \\ L_2^* = F^{-1} \left(1 + \alpha - F(L_1^*) \right) & \text{a.s., when } L_1 \ge F^{-1}(\alpha). \end{cases}$$

In Figure 1, left, we show the copula of the risk vector (L_1^*, L_2^*) . In the right part of the same figure, we show the support of the risk vector (L_1^*, L_2^*) when L_1^* and L_2^* are both Pareto(2)-distributed. The *support* of a random vector X is the smallest closed set A such that $P(X \notin A) = 0$. It is interesting to note the interdependence of (L_1^*, L_2^*) . In the upper $(1 - \alpha)$ part of their supports, the marginal risks L_1^* and L_2^* are countermonotonic. This means that the variance of the sum of the upper $(1 - \alpha)$ parts of their supports is minimized. In the lower α -part of their supports, the marginal risks L_1^* and L_2^* are a.s. identical and hence comonotonic. This is however not relevant since the interdependence in this lower part of the joint distribution can be chosen arbitrarily; see Puccetti and Rüschendorf (2012a, Theorem 2.1).

The case d=2 is mainly pedagogical. The typical dimensions used in practice may vary from d=7 or 8 to 56, say, for the aggregation of Operational Risk factors; see Moscadelli (2004), but may go up to d in the several hundreds for typical aggregation models for other applications like hierarchical risk aggregation models; see for instance Arbenz et al. (2012). In the case case d>2, the sharp bound $\overline{\mathrm{VaR}}_{\alpha}(L)$ has been obtained only recently in the homogeneous case under different

sets of assumptions. For a distribution function F, define the dual bound D(s) as

$$D(s) = \inf_{t < s/d} \frac{d \int_t^{s - (d-1)t} \overline{F}(x) dx}{(s - dt)}, \tag{12}$$

where $\overline{F}(x) = 1 - F(x)$. The dual bound D(s) in (12) is an upper bound on the tail function of L, that is

$$P(L_1 + \cdots + L_d > s) \leq D(s);$$

see for instance Puccetti and Rüschendorf (2012a). This directly implies that

$$\overline{\operatorname{VaR}}_{\alpha}(L) \le D^{-1}(1 - \alpha) = \inf\{s \in \mathbb{R} : D(s) \ge 1 - \alpha\}. \tag{13}$$

The VaR bound $D^{-1}(1-\alpha)$ is numerically easy to evaluate independently of the size d of the portfolio (L_1, \ldots, L_d) . Under some extra assumptions, we have that the inequality in (13) becomes an equality.

Proposition 4 (Dual bound). In the homogeneous case $F_i = F, 1 \le i \le d$, with $d \ge 3$, let F be a continuous distribution with an unbounded support and an ultimately decreasing density. Suppose that for any sufficiently large threshold s the infimum in (12) is attained at some a < s/d, that is assume that

$$D(s) = \frac{d \int_a^b \overline{F}(x) dx}{(b-a)} = \overline{F}(a) + (d-1)\overline{F}(b), \tag{14}$$

where b = s - (d - 1)a, with $F^{-1}(1 - D(s)) \le a < s/d$. Then, for any sufficiently large threshold α we have that

$$\overline{\text{VaR}}_{\alpha}(L) = D^{-1}(1 - \alpha). \tag{15}$$

Remark 5. The above proposition is a particular case of Puccetti and Rüschendorf (2012c, Theorem 2.5). We refer to the latter paper and references therein for mathematical details in addition to the following points:

- 1. Under the assumptions of Proposition 4, the infimum in (12) is attained at a < s/d if and only if the first order condition (14) holds. In order to calculate the worst-VaR value $\overline{\text{VaR}}_{\alpha}(L)$ it is sufficient to compute the function D(s) by solving numerically the univariate equation (14) and hence to compute numerically its inverse D^{-1} at the level $(1-\alpha)$. The treatment of an *arbitrary* number of identically distributed risks is then made possible; see Figure 2 and Table 2.
- 2. For the Pareto distribution (11) with tail parameter $\theta > 0$ we have that

$$\overline{\text{VaR}}_{\alpha}(L) = D^{-1}(1 - \alpha),$$

for any $\alpha \in (0, 1)$. Portfolios of Pareto distributed risks are studied in Table 2.

3. The sharpness of the bound $D^{-1}(1-\alpha)$ in (15) can be stated under different sets of assumptions for the distribution function F. To cite a most useful case, sharpness holds for distributions F having a concave density on the interval (a,b). This allows for instance to compute the sharp bound $\overline{\mathrm{VaR}}_{\alpha}(L) = D^{-1}(1-\alpha)$ in case of Gamma and LogNormal distributions; see Figure 2.

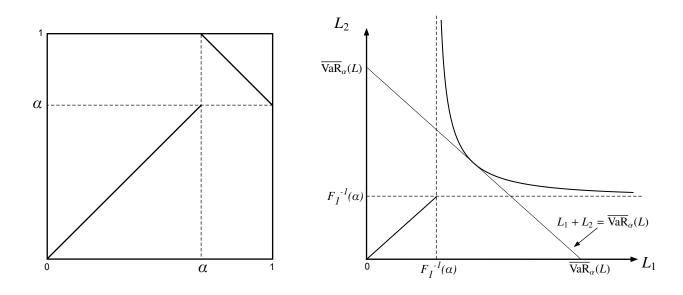


Figure 1: Bivariate copula (left) and support (right) of the vector (L_1^*, L_2^*) attaining the worst-possible VaR for $L_1 + L_2$ when L_1 and L_2 are both Pareto(2)-distributed.

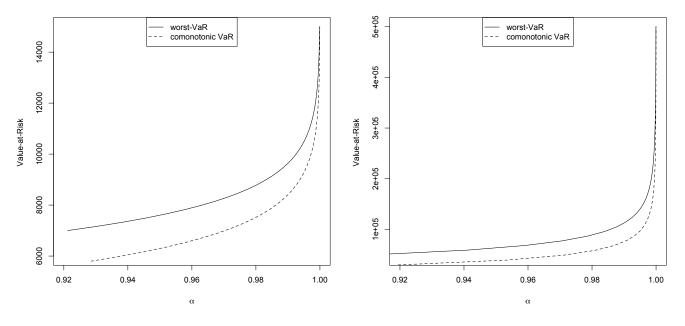


Figure 2: Value for $\overline{\text{VaR}}_{\alpha}(L)$ (see (15)) and $\text{VaR}_{\alpha}^{+}(L)$ (see (8)), for the sum of d=1000 Gamma(3,1)- (left) and LogNormal(2,1)- (right) distributed risks.

- 4. The equation (15) holds typically for distributions F and confidence levels α standardly used in quantitative risk management, also in the case of heavy tailed, infinite-mean models.
- 5. So far, there does not exist a method which allows to compute $VaR_{\alpha}(L)$ analytically for $d \ge 3$.

When the distribution F satisfies the assumptions of Proposition 4, a worst-case dependence vector (L_1^*,\ldots,L_d^*) such that $\overline{\mathrm{VaR}}_\alpha(L) = \mathrm{VaR}_\alpha(L_1^*+\cdots+L_d^*)$ has been described in Wang et al. (2011) and Puccetti and Rüschendorf (2012c). Here the concept of complete mixability is crucial. The risk vector (L_1^*,\ldots,L_d^*) satisfies the following two properties:

(a) When one of the L_i^* 's lies in the interval (a,b), then all the L_i^* 's lie in (a,b) and are a d-complete mix, i.e. for all $1 \le i \le d$,

$$P(L_1^* + \cdots + L_d^* = s | L_i \in (a, b)) = 1;$$

(b) For all $1 \le i \le d$, we have that

$$P\left(L_j = F_{a^*}^{-1}\left((d-1)\overline{F}_{a^*}(L_i)\right)\Big|L_i \ge b\right) = 1, \text{ for all } j \ne i,$$

where $a^* = F^{-1}(1 - D(s))$ and $F_{a^*}(x) = (F(x) - F(a^*))/\overline{F}(a^*)$. $\overline{F}(a^*)$ is the distribution of the random variable $Y_{a^*} \stackrel{\text{d}}{=} (L_1|L_1 \ge a^*)$. The interdependence described by the two properties above can be summarized as:

The two properties (a) and (b) determine the behavior of the worst-case dependence only in the upper $(1-\alpha)$ parts of the marginal supports where $L_i \geq a^*, 1 \leq i \leq d$. Analogous to the case d=2, the interdependence coupling in the α lower parts of the marginal supports can be set arbitrarily.

In Figure 3 we show a two-dimensional projection of the d-variate copula merging the upper $(1-\alpha)$ parts of the optimal risks L_i^* . In practice, only two situations can occur: either one of the risks is large (above the threshold b) and all the others are small (below the threshold a), or all the risks are of medium size (they lie in the interval (a,b)) with their sum being equal to the threshold $\overline{\mathrm{VaR}}_{\alpha}(L)$. This is a negative dependence scenario analogous to the one underlying Figure 1. In fact the worst-VaR scenario contains a part where the risks are d-completely mixable, with the variance of their sum being equal to zero. In the remaining part, it exhibits mutual exclusivity: only one risk can be large at one time.

For a risk portfolio $(L_1, ..., L_d)'$ it is of interest to study the *superadditivity ratio*

$$\delta_{\alpha}(d) = \frac{\overline{\text{VaR}}_{\alpha}(L)}{\text{VaR}_{\alpha}^{+}(L)}$$

between the worst-possible VaR and the comonotonic VaR, at some given level of probability $\alpha \in (0, 1)$. The value $\delta_{\alpha}(d)$ measures how much VaR can be superadditive as a function of the

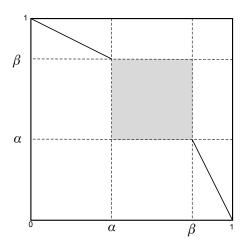


Figure 3: One of the identical two-dimensional projections of the d-variate copula merging the upper $(1-\alpha)$ parts of the optimal risks L_i^* . In the figure, we have $\alpha=1-D(s)$ and $\beta=\alpha/(d-1)$. The grey area represents a completely mixable part.

dimensionality d of the risk portfolio under study. For instance, for elliptically distributed risks it is well known that $\delta_{\alpha}(d) = 1$ for any $d \geq 1$; see McNeil et al. (2005, Theorem 6.8). In Figure 4 and Figure 5, left, we plot the function $\delta_{\alpha}(d)$ for a number of different homogeneous portfolios. In these cases, $\delta_{\alpha}(d)$ seems to settle down to a limit in d fairly fast. We therefore define

$$\delta_{\alpha} = \lim_{d \to +\infty} \delta_{\alpha}(d),$$

whenever this limit exists. For large dimensions d one can then approximate the worst-possible VaR value as

$$\overline{\operatorname{VaR}}_{\alpha}(L) \approx \delta_{\alpha} \operatorname{VaR}_{\alpha}^{+}(L) = d\delta_{\alpha} \operatorname{VaR}_{\alpha}(L_{1}).$$

We study the superadditivity constant δ_{α} for some homogeneous risk portfolios of interest in finance and insurance. For portfolios of LogNormal(2,1)-distributed risks, we have $\delta_{0.99} \cong$ 1.49 and $\delta_{0.999} \cong 1.37$; see Figure 4, left. For portfolios of Gamma(3,1)-distributed risks, we have $\delta_{0.99} \cong 1.15$ and $\delta_{0.999} \cong$ 1.11; see Figure 4, right. For portfolios of Pareto(2)-distributed risks, we have $\delta_{0.99} \cong 2.11$ and $\delta_{0.999} \cong 2.03$; see Figure 5, left. In Figure 5, right, one can see that the limit constant δ_{α} depends on the tail parameter θ of the Pareto marginals: the smaller the tail parameter θ , the more superadditive the VaR of the sums of the risks can be. It is also interesting that, in the examples studied, the superadditivity ratio is larger for smaller levels of α . A figure analogous to Figure 5 cannot be obtained analytically for the ratio $VaR_{\alpha}^{+}(L)/VaR_{\alpha}(L)$; see point 5 in Remark 5. Risk portfolios showing an analogous behavior can be found in other studies like Mainik and Rüschendorf (2010), Mainik and Embrechts (2012) and Mainik and Rüschendorf (2012).

2.2. Inhomogeneous marginals

If one drops the assumption of identically distributed risks, the bounds given in (10) and (15) cannot be used. For d = 2,

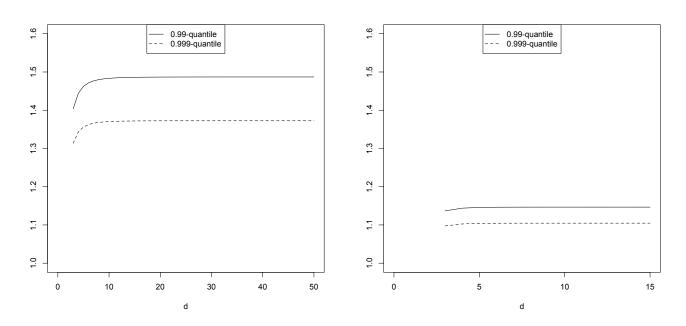


Figure 4: Plot of the function $\delta_{\alpha}(d)$ versus the dimensionality d of the portfolio for a risk vector of LogNormal(2,1)-distributed (left) and Gamma(3,1)-distributed (right) risks, for two different quantile levels.

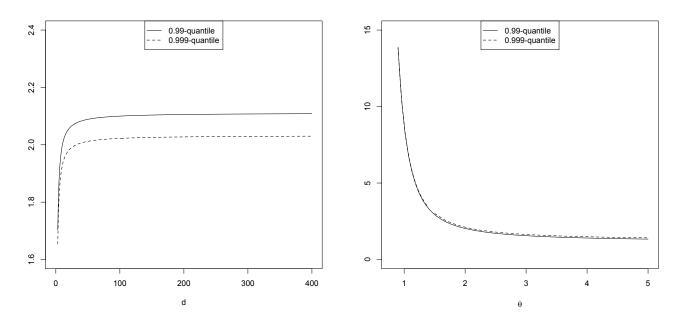


Figure 5: Left: plot of the function $\delta_{\alpha}(d)$ versus the dimensionality d of the portfolio for a risk vector of Pareto(θ)-distributed risks, for two different quantile levels and $\theta = 2$. Right: plot of the limit constant δ_{α} versus the tail parameter θ of the Pareto distribution.

the sharp bounds $VaR_{\alpha}(L)$ and $\overline{VaR}_{\alpha}(L)$ can be calculated easily also in the inhomogeneous case using Rüschendorf (1982, Proposition 1); see also Puccetti and Rüschendorf (2012a, Theorem 2.7). In higher dimension $d \ge 3$ the computation of the dual functional D(s) with different marginal distributions may become numerically cumbersome. The numerical complexity of the dual bound D(s) typically increases with the number of blocks of marginals with identical distributions. For instance, if all the d marginal distributions are different, the computation of dual bounds is manageable up to small dimension d = 10. An example with d = 8 is illustrated in Embrechts and Puccetti (2006a). However, it is possible to compute the dual bound D(s) for relatively large dimensions d if the inhomogeneous risks L_i can be divided in n sub-groups having homogeneous marginals within. In this case, the numerical complexity of the dual bound D(s) only depends on n, and is independent of the cardinality of each of the sub-groups of homogeneous marginals. It is also important to remark that the sharpness of dual bounds in dimension $d \ge 3$ has not been proved for inhomogeneous marginals.

For the computation of bounds on distribution functions Puccetti and Rüschendorf (2012b) introduced a *rearrangement algorithm* (RA) working well for dimension $d \leq 30$. In order to compute sharp bounds for the VaR, we adapt and modify this RA in order to compute the sharp bounds $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\alpha}(L)$ in the inhomogeneous case. Our modification allows to apply the algorithm to high-dimensional inhomogeneous portfolios up to dimension d=600 which previously were well out of the range of numerical and analytical methods. The RA can compute the worst and best VaR values in (4) with good accuracy for *any* set of marginals F_i and relatively large dimensions d.

In the following, we say that two vectors $\mathbf{a}, \mathbf{b} \in \mathbb{R}^N$ are oppositely ordered if $(a_j - a_k)(b_j - b_k) \le 0$ holds for all $1 \le j, k \le N$.

Rearrangement Algorithm (RA) to compute $VaR_{\alpha}(L)$.

- 1. Fix an integer N and the desired level of confidence α .
- 2. Define the matrices $X^{\alpha} = (x_{i,j}^{\alpha})$ and $Y^{\alpha} = (y_{i,j}^{\alpha})$ as

$$x_{i,j}^{\alpha} = F_j^{-1} \left(\alpha + \frac{(1-\alpha)(i-1)}{N} \right), \quad y_{i,j}^{\alpha} = F_j^{-1} \left(\alpha + \frac{(1-\alpha)i}{N} \right),$$
 (16)

for $1 \le i \le N$, $1 \le j \le d$.

- 3. Permute randomly the elements in each column of X^{α} and Y^{α} .
- 4. Iteratively rearrange the j-th column of the matrix X^{α} so that it becomes oppositely ordered to the sum of the other columns, for $1 \le j \le d$.
- 5. Repeat Step 4. until no further changes occur, that is until a matrix $X^* = (x_{i,j}^*)$ is found with each column oppositely ordered to the sum of the others.

- 6. Apply Steps 4.–5. to the matrix Y^{α} until a matrix $Y^* = (y_{i,j}^*)$ is found with each column oppositely ordered to the sum of the others.
- 7. Define

$$\underline{s}_N = \min_{1 \le i \le N} \sum_{1 \le j \le d} x_{i,j}^*$$
 and $\overline{s}_N = \min_{1 \le i \le N} \sum_{1 \le j \le d} y_{i,j}^*$.

Then we have $\underline{s}_N \leq \overline{s}_N$ and

$$\lim_{N\to\infty} \overline{s}_N = \lim_{N\to\infty} \underline{s}_N = \overline{\mathrm{VaR}}_{\alpha}(L).$$

In practice we find the algorithm to give excellent approximations for moderately large N.

Rearrangement Algorithm (RA) to compute $\underline{\text{VaR}}_{\alpha}(L)$.

- 1. Fix an integer N and the desired level of confidence α .
- 2. Define the matrices $X^{\alpha} = (x_{i,j}^{\alpha})$ and $Y^{\alpha} = (y_{i,j}^{\alpha})$ as

$$x_{i,j}^{\alpha} = F_j^{-1} \left(\frac{\alpha(i-1)}{N} \right), \quad y_{i,j}^{\alpha} = F_j^{-1} \left(\frac{\alpha i}{N} \right),$$

for $1 \le i \le N$, $1 \le j \le d$.

- 3. Permute randomly the elements in each column of X^{α} and Y^{α} .
- 4. Iteratively rearrange the j-th column of the matrix X^{α} so that it becomes oppositely ordered to the sum of the other columns, for $1 \le j \le d$.
- 5. Repeat Step 4. until no further changes occur, that is until a matrix $X^* = (x_{i,j}^*)$ is found with each column oppositely ordered to the sum of the others.
- 6. Apply Steps 4.–5. to the matrix Y^{α} until a matrix $Y^* = (y_{i,j}^*)$ is found with each column oppositely ordered to the sum of the others.
- 7. Define

$$\underline{t}_N = \max_{1 \leq i \leq N} \sum_{1 \leq j \leq d} x_{i,j}^* \quad \text{ and } \overline{t}_N = \max_{1 \leq i \leq N} \sum_{1 \leq j \leq d} y_{i,j}^*.$$

Then we have $\underline{t}_N \leq \overline{t}_N$ and

$$\lim_{N\to\infty} \overline{t}_N = \lim_{N\to\infty} \underline{t}_N = \underline{\mathrm{VaR}}_{\alpha}(L).$$

In practice we find the algorithm to give excellent approximations for moderately large N.

Remark 6. For mathematical details about the RA, we refer the reader to Puccetti and Rüschendorf (2012b). Here we limit our attention to the following, more practical points:

1. We call the interval $(\underline{s}_N, \overline{s}_N)$ the rearrangement range for $\overline{\text{VaR}}_{\alpha}(L)$. The length $(\overline{s}_N - \underline{s}_N)$ of this interval depends on the dimensionality d of the risk portfolio under study and on N. For a fixed d, the sequence $(\overline{s}_N - \underline{s}_N)$ asymptotically goes to zero as o(1/N). For sufficiently large N, we also have that $\underline{s}_N \leq \overline{\text{VaR}}_{\alpha}(L)$. Analogous considerations can be made for the rearrangement range $(\underline{t}_N, \overline{t}_N)$ for $\underline{\text{VaR}}_{\alpha}(L)$. For sufficiently large N we have that $\overline{t}_N \geq \underline{\text{VaR}}_{\alpha}(L)$.

- 2. The randomization Step 3. is introduced in order to avoid pathological examples in which the sequences $(\overline{s}_N \underline{s}_N)$ and $(\overline{t}_N \underline{t}_N)$ do not converge to zero (we thank Robert Weismantel for pointing this out). In Table 2, we check the accuracy of the RA for some Pareto(2) risk portfolios for which we know, by Proposition 4, the exact value of $\overline{\text{VaR}}_{\alpha}(L)$. This table also highlights the possibly large difference between the comonotonic $\text{VaR}_{\alpha}^+(L)$ and the worst-possible $\overline{\text{VaR}}_{\alpha}(L)$. In Table 2 we use different dimensions d as well as values of N which represent a good compromise between computational time used and accuracy obtained. In order to perform all the computations in the remainder of the paper we use an Apple MacBook (2 GHz Intel Core 2 Duo, 2 GB RAM). Computation times can no doubt be dramatically reduced on a more powerful machine.
- 3. As a numerical algorithm, the RA can be used with *any* type of marginal distributions. The figures in Table 2 are obtained for a homogeneous portfolio so as to be able to check the accuracy of the RA via the dual bound in Proposition 4. In general, both the accuracy and the computation time of the RA are *not* affected by the type of the marginal distributions used. We apply the RA to inhomogeneous marginals in Section 4.

The probabilistic idea behind the RA is easy. For a fixed $\alpha \in [0,1]$, the j-th columns of the matrices X^{α} and Y^{α} represent two stochastically ordered N-point discretizations of the $(1-\alpha)$ upper parts of the supports of the marginal risk L_j . The RA rearranges the columns of X^{α} into the matrix X^* in order to find the maximal value \underline{s}_N such that the componentwise sum of any row of X^* is larger than \underline{s}_N . Analogously, the RA rearranges the columns of Y^{α} into the matrix Y^* in order to find the maximal value \overline{s}_N such that the componentwise sum of any row of Y^* is larger than \overline{s}_N . For N large enough we have that $\underline{s}_N \leq \overline{\mathrm{VaR}}_{\alpha}(L) \simeq \overline{s}_N$ as a direct consequence of Puccetti and Rüschendorf (2012b, Theorem 3.1). An analogous mechanism yields $\mathrm{VaR}_{\alpha}(L)$.

We illustrate the RA in an example with homogeneous marginals. We consider d=3 marginals of Pareto type with identical tail parameters $\theta=2.5$. Then, we set N=50 and compute $\overline{\text{VaR}}_{\alpha}(L)$ for $\alpha=0.99$ via the RA. The initial matrix X^{α} defined in (16) for $\alpha=0.99$ is shown in Table 1 (A). The j-th column of X^{α} represents a 50-point discretization of the upper 1% of the support of the j-th marginal distribution. In the same (A) part of the table, we also show the N-dimensional vector of the rowwise sums of X^{α} , as well as the d-dimensional vector having as components the aggregate sums of the columns of X^{α} .

During the iteration of the algorithm (Steps 3.–5.), the elements within each column of X^{α} are re-shuffled until a matrix X^* is found with each column oppositely ordered to the sum of the others; see Table 1 (B). This rearrangement procedure of the columns of X^{α} aims at maximizing the minimal component of the vector of the row-wise sums of X^* . Indeed, note how the minimal component of the row-wise sums (15.92872) is increased (to 24.46538) when passing from X^{α} to X^* , while the column-wise sums remain unchanged (the marginals are still

the same). Compared to X^{α} , the matrix X^* represents a different coupling (copula) of the same marginals in which the variance of the marginal numbers (rows) is reduced. The minimal component of the vector of the sums of the rows of X^* is $\underline{s}_{50} = 24.46538$ and represents a lower bound on $\overline{\text{VaR}}_{\alpha}(L)$. Performing an analogous rearrangement of the column of the matrix Y^{α} one finds $\overline{s}_{50} = 25.12000$, which is instead approximately an upper bound on $\overline{\text{VaR}}_{\alpha}(L)$. Note that the estimates \underline{s}_{50} and \bar{s}_{50} are actually random as in Step 3. the RA performs a randomization of the column of X^{α} . This random uncertainty becomes negligible for values of N large enough. From the application of the RA described above for N = 50 one obtains $VaR_{\alpha}(L) \in [24.47, 25.12]$. It is sufficient to run the algorithm with N = 1.0e05 to obtain the first two decimals of $\overline{\text{VaR}}_{\alpha}(L) =$ 24.93 in about 0.2 seconds. Of course, in this pedagogical case one could instantly obtain the exact value $\overline{\text{VaR}}_{\alpha}(L) = 24.93$ from Proposition 4. The power of the RA is that it can be applied also to inhomogeneous portfolios of risks and is able to compute numerically also $VaR_{\alpha}(L)$.

It is interesting to see that already for N=50, the final matrix X^* in Table 1 (B) approximates the worst-case dependence for the sum of continuous homogeneous marginals shown in Figure 3. Indeed, one can check that basically two structures occur the rows of X^* : either all the components of the row are close to each other, and sum up to a value which is just above the threshold $\underline{s}_{50}=24.47$ (e.g. row 9), or one of them is large and all the others are small (e.g. row 46). Of course, this structural dichotomy becomes much clearer when N increases.

As a more realistic example stemming from an application of Extreme Value Theory, we study a risk portfolio where the marginal losses are distributed by a Generalized Pareto Distribution (GPD), that is we assume

$$F_i(x) = 1 - \left(1 + \xi_i \frac{x}{\beta_i}\right)^{-1/\xi_i}, x \ge 0,$$

for all i = 1, ..., d. Note that for a GPD distribution, whenever $0 < \xi_i \le 1$, $E(L_i) = \infty$, and for $1 < \xi_i < 2$, $E(L_i) < \infty$ but $var(L_i) = \infty$. We choose for the dimensionality d = 8 of the portfolio and the parameters of the GPD distributions, the values reported in the quantitative impact study data of Moscadelli (2004). Under these marginal assumptions, the risk portfolio $(L_1, \ldots, L_d)'$ shows a very heavy-tailed behavior, with six out of eight losses L_i exhibiting an infinite mean marginal model. In the other two cases where the mean is finite, the loss distributions do not have finite variance. In Figure 6, we plot the VaR range (5) as well as the estimate for $VaR_{\alpha}^{+}(L)$, versus the confidence level α . For practice, the wide dependence range for the values of α typically used, that is $\alpha = 0.99, 0.999$, should raise some concerns. The large difference between the worst- and best-possible VaR under the same marginal assumptions can be better appreciated on a log-scale in the same figure.

For the dimension d=8 in the Moscadelli example, the RA algorithm produces accurate estimate of $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\alpha}(L)$ in about 30 seconds. However, the RA algorithm can handle large portfolios as described in Table 2. The results in these examples imply a considerable model uncertainty issue underlying VaR calculations for confidence levels close to 1. Of course,

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                                                           8.102821
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                                                          13.734232
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                                                                                   8.952679 24.51256
     29.170882
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                            29.170882 87.51265
                                                           6.845648
                                                                       8.714231
    444.710518 444.710518 444.710518
                                                         444.710518 444.710518 444.710518
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Table 1: (A): The matrix X^{α} defined in (16) for $\alpha = 0.99$ and N = 50 (representing comonotonicity among the discrete marginals); (B): The matrix X^* derived by the iterative rearrangement of the columns of X^{α} .

statistical uncertainty at the level of the marginal distributions further compounds the problem.

3. Positive dependence information

The worst-VaR copulas given in Section 2, Figures 1 and 3, are probably considered as unrealistic due to their minimal variance parts in which the risks are countermonotonic (for d=2) or completely mixable (in the case $d \geq 3$). Of course, a positive dependence structure combined with the knowledge of the marginal distributions of $(L_1, \ldots, L_d)'$ will tighten the interval of admissible VaRs in (5). However, assuming that the risks are positively dependent does not eliminate countermonotonicty and completely mixable parts from the worst-VaR scenarios and does not necessarily lower the estimate of $\overline{\text{VaR}}_{\alpha}(L)$. This latter point is the object of this section. We start by introducing a natural concept of positive dependence.

Definition 7. The risk vector $(L_1, ..., L_d)'$ is said to be positively lower orthant dependent (PLOD) if for all $(x_1, ..., x_d)' \in \mathbb{R}^d$

$$P(L_1 \le x_1, \dots, L_d \le x_d) \ge \prod_{i=1}^d P(X_i \le x_i) = \prod_{i=1}^d F_i(x_i).$$
 (17)

The risk vector $(L_1, ..., L_d)'$ is said to be positively upper orthant dependent (PUOD) if for all $(x_1, ..., x_d)' \in \mathbb{R}^d$

$$P(L_1 > x_1, \dots, L_d > x_d) \ge \prod_{i=1}^d P(X_i > x_i) = \prod_{i=1}^d \overline{F}_i(x_i).$$
 (18)

Finally, the risk vector $(L_1, ..., L_d)'$ is said to be positively orthant dependent (POD) if it is both PLOD and PUOD.

For d=2, conditions (17) and (18) are equivalent. However, this is *not* the case for $d\geq 3$. In higher dimensions the PLOD and PUOD concepts are distinct; see for instance Nelsen (2006, Section 5.7). If $(L_1,\ldots,L_d)'$ has copula C, condition (17) can be equivalently expressed as $C\geq \Pi$, the independence copula. Analogously, condition (18) can be written as $\overline{C}\geq \overline{\Pi}$, where \overline{C} denotes the joint tail function of a copula C, also referred to as the survival copula; see Nelsen (2006, Section 2.6).

Under the addition of a positive dependence restriction, VaR bounds for the sum of risks have been derived in Theorem 3.1 in Embrechts et al. (2005); see also Rüschendorf (2005) and Puccetti and Rüschendorf (2012a). We state this result here for the case of identical marginals using the same notation as in the unconstrained case with no dependence information.

Proposition 8. In the homogeneous case $F_i = F$, $1 \le i \le d$, let F be a distribution with decreasing density on its entire domain. If the risk vector $(L_1, \ldots, L_d)'$ is PLOD then, for any fixed real threshold s, we have

$$\overline{\operatorname{VaR}}_{\alpha}(L) \le dF^{-1}\left((1-\alpha)^{\frac{1}{d}}\right). \tag{19}$$

Remark 9. In Embrechts et al. (2005), the bound (19) is given in a slightly more complicated form for any set of marginal distributions. In the same reference, an analogous bound for $\overline{\text{VaR}}_{\alpha}(L)$ is given if the risk vector is assumed to be PUOD.

In the case d=2 the inequality given in (19) is sharp. In Figure 7, left, we show the copula of a PLOD risk vector $(L_1^*, L_2^*)'$ for which $\text{VaR}_\alpha(L_1^* + L_2^*) = \overline{\text{VaR}}_\alpha(L)$. Even if the structure of dependence of this vector is PLOD, its geometry is not so different if compared to the optimal copula in the unconstrained case (Figure 1, left). Again, the copula of $(L_1^*, L_2^*)'$ contains a countermonotonic part, in which the risks are a.s. decreasing functions of each other. Thus, the assumption of positive dependence *does not* eliminate the possibility of such optimal (*unrealistic*) copulas. In particular, assuming positive correlation does not improve the bound. The reason for this is not to be found in the concept of VaR but rather raises some questions about the appropriateness of PLOD (PUOD) as a concept of positive (negative) dependence.

Given the shape of the copula attaining the bound (19) under additional positive dependence restrictions, one cannot expect an essential improvement of the VaR bound given in the unconstrained case when only the marginals of the L_i 's are known. Indeed, in Figure 7, right, we plot the worst-VaR value $\overline{\text{VaR}}_{\alpha}(L)$ (see (19)) and the comonotonic $\text{VaR}_{\alpha}^+(L)$ (see (8)), for the sum of two Pareto(2) distributions. The improvement of the bound given by the additional information is negligible.

The situation gets more involved in higher dimensions ($d \ge$ 3), as the bound (19) fails to be sharp. The dual bound given in (15) for the unconstrained case actually turns out to be better than (19) with positive dependence information; this can be seen in Figure 8. This is not so suprising, as the dual bound given in (15) derives from a different methodology based on the powerful tool offered by the theory of mass transportation; see Embrechts and Puccetti (2006b) on this. As a matter of fact, the bound (19) is not useful for higher dimensions ($d \ge 3$) where the search for a sharp bound with marginal and positive dependence information is still open. However, we do not expect much improvement over the dual bounds even for optimal ones in the positive dependence case. Take for instance the problem of maximizing the covariance of $(L_1, L_2)'$ when d = 2 and the marginals F_1 and F_2 are given. By Hoeffding's covariance representation formula, see McNeil et al. (2005, Lemma 5.2.4), one has

$$Cov(L_1, L_2) = \int (F(x_1, x_2) - F_1(x_1)F_2(x_2)) dx_1 dx_2,$$

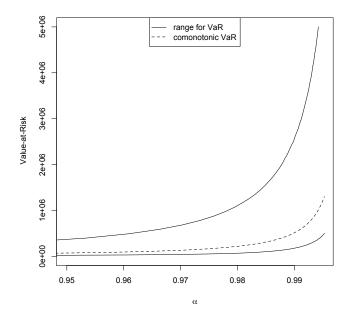
where F is the joint distribution of $(L_1, L_2)'$. It is clear that here the PLOD constraint $F(x_1, x_2) \ge F_1(x_1)F_2(x_2)$ does not help to improve an upper bound on $Cov(L_1, L_2)$.

4. Higher dimensional dependence information

For a vector $(L_1, ..., L_d)'$ for which one only knows the marginal distributions $F_1, ..., F_d$, we have that

$$\operatorname{VaR}_{\alpha}(L) \le \operatorname{VaR}_{\alpha}(L_1 + \dots + L_d) \le \overline{\operatorname{VaR}}_{\alpha}(L).$$
 (20)

If one adds PLOD/PUOD information on top of the knowledge of the marginals, the worst VaR in (20) is only minimally affected. It is clear that in practice more dependence information on the vector $(L_1, \ldots, L_d)'$ may be available. Such a case would



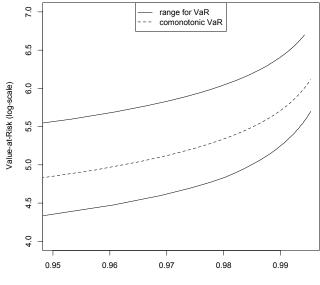


Figure 6: VaR range (5), and comonotonic VaR (8) (in log-scale on the right) for the sum of d = 8 GPD risks with parameters following Moscadelli (2004), based on RA for N = 1.0e05.

be where specific assumptions on sub-vectors of $(L_1, \ldots, L_d)'$ are made. One reason for this could be that the individual risk factors may be grouped in economically relevant sectors. This would lead to a narrowing of the range on $VaR_{\alpha}(L_1 + \cdots + L_d)$ in (20).

Thus, we consider the case that not only the one-dimensional marginal distributions of the risk vector are known, but also that for a class \mathcal{E} of sets $J \subset \{1,\ldots,d\}$, the joint marginal distributions $F_J, J \in \mathcal{E}$ are fixed. In this case, we get the generalized Fréchet class

$$\mathfrak{F}_{\mathcal{E}}=\mathfrak{F}(F_J,J\in\mathcal{E})$$

of all probability measures on \mathbb{R}^d having sub-vector models F_J on \mathbb{R}^J , for all $J \in \mathcal{E}$. W.l.o.g. we assume that $\bigcup_{J \in \mathcal{E}} J = \{1, \ldots, d\}$. Thus, we have

$$\mathfrak{F}_{\mathcal{E}} \subset \mathfrak{F}(F_1,\ldots,F_d),$$

that is $\mathfrak{F}_{\mathcal{E}}$ is a sub-class of the class of all possible dependence structures. The knowledge of higher dimensional joint distributions is in general not sufficient to determine the joint model of $(L_1,\ldots,L_d)'$. Nevertheless, having higher dimensional information restricts the class of possible dependence structures and thus leads to improved upper and lower bounds for the VaR of the joint portfolio.

As an example we consider, for d even, a class \mathcal{E} of particular interest in actuarial applications: we set $\mathcal{E} = \{\{2j - 1, 2j\} : j = 1, \dots, d/2\}$, defining the Fréchet class

$$\mathfrak{F}_{\mathcal{E}} = \mathfrak{F}(F_{12}, F_{34}, \dots, F_{d-1d}).$$

Hence in this case risk estimates on the global position $L_1 + \cdots + L_d$ have to be obtained based on distributional information

for all two-dimensional sub-vectors $(L_{2j-1}, L_{2j})'$. Other examples of marginals classes \mathcal{E} have been treated in Puccetti and Rüschendorf (2012a) and Embrechts and Puccetti (2010a).

Our aim is to find bounds for the tail risks

$$\overline{\mathrm{VaR}}_{\alpha}^{\mathcal{E}}(L) = \sup \left\{ \mathrm{VaR}_{\alpha}(L_1 + \dots + L_d) : F_L \in \mathfrak{F}_{\mathcal{E}} \right\}, \qquad (21a)$$

$$\underline{\operatorname{VaR}}_{\alpha}^{\mathcal{E}}(L) = \inf \left\{ \operatorname{VaR}_{\alpha}(L_1 + \dots + L_d) : F_L \in \mathfrak{F}_{\mathcal{E}} \right\}, \quad (21b)$$

which improve the corresponding bounds $\overline{\mathrm{VaR}}_{\alpha}(L)$ and $\underline{\mathrm{VaR}}_{\alpha}(L)$ defined in (4). If $F_L \in \mathfrak{F}_{\mathcal{E}}$, we have

$$\underline{\mathrm{VaR}}_{\alpha}(L) \leq \underline{\mathrm{VaR}}_{\alpha}^{\mathcal{E}}(L) \leq \mathrm{VaR}_{\alpha}(L) \leq \overline{\mathrm{VaR}}_{\alpha}^{\mathcal{E}}(L) \leq \overline{\mathrm{VaR}}_{\alpha}(L). \tag{22}$$

A reduction method introduced in Puccetti and Rüschendorf (2012a) allows to find reduced bounds $\overline{\mathrm{VaR}}_{\alpha}^{\mathcal{E}}(L)$ and $\underline{\mathrm{VaR}}_{\alpha}^{\mathcal{E}}(L)$ using Proposition 4 and the RA introduced in Section 2. The reduction method consists in associating to the risk vector $(L_1,\ldots,L_d)'$ with $\mathfrak{F}_L \in \mathfrak{F}_{\mathcal{E}}$ the random vector $(Y_1,\ldots,Y_n)'$ defined by

$$Y_i = L_{2i-1} + L_{2i}, \ j = 1, \dots, n,$$
 (23)

where n = d/2. If we also denote by H_j the distribution of Y_j , the risk vector $(Y_1, \ldots, Y_n)'$ has fixed marginals H_1, \ldots, H_n . Therefore, it is possible to apply the techniques introduced in Section 2 to compute the *reduced* VaR bounds:

$$\overline{\mathrm{VaR}}_{\alpha}^{r}(L) = \sup \left\{ \mathrm{VaR}_{\alpha}(Y_1 + \dots + Y_n) : F_{Y} \in s\mathfrak{F}(H_1, \dots, H_n) \right\},$$
(24a)

$$\underline{\operatorname{VaR}}_{\alpha}^{r}(L) = \inf \left\{ \operatorname{VaR}_{\alpha}(Y_{1} + \dots + Y_{n}) : F_{Y} \in \mathfrak{F}(H_{1}, \dots, H_{n}) \right\}. \tag{24b}$$

d = 8	N = 1.0e05	avg time: 30 secs		
α	$\underline{\operatorname{VaR}}_{\alpha}(L)$ (RA range)	$VaR^+_{\alpha}(L)$ (exact)	$\overline{\operatorname{VaR}}_{\alpha}(L)$ (exact)	$\overline{\text{VaR}}_{\alpha}(L) \text{ (RA range)}$
0.99	9.00 - 9.00	72.00	141.67	141.66–141.67
0.995	13.13 - 13.14	105.14	203.66	203.65-203.66
0.999	30.47 - 30.62	244.98	465.29	465.28-465.30
d = 56	N = 1.0e05	avg time: 9 mins		
α	$\underline{\operatorname{VaR}}_{\alpha}(L)$ (RA range)	$VaR^+_{\alpha}(L)$ (exact)	$\overline{\operatorname{VaR}}_{\alpha}(L)$ (exact)	$\overline{\text{VaR}}_{\alpha}(L) \text{ (RA range)}$
0.99	45.82 - 45.82	504	1053.96	1053.80-1054.11
0.995	48.60 - 48.61	735.96	1513.71	1513.49-1513.93
0.999	52.56 - 52.58	1714.88	3453.99	3453.49-3454.48
d = 648	N = 5.0e04	avg time: 8 hrs		
α	$\underline{\operatorname{VaR}}_{\alpha}(L)$ (RA range)	$VaR^+_{\alpha}(L)$ (exact)	$\overline{\text{VaR}}_{\alpha}(L) \text{ (exact)}$	$\overline{\text{VaR}}_{\alpha}(L) \text{ (RA range)}$
0.99	530.12 - 530.24	5832.00	12302.00	12269.74-12354.00
0.995	562.33 - 562.50	8516.10	17666.06	17620.45-17739.60
0.999	608.08 - 608.47	19843.56	40303.48	40201.48-40467.92

Table 2: Estimates for $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\gamma}(L)$ for random vectors of Pareto(2)-distributed risks.

Using the key fact that $L_1 + \cdots + L_d = Y_1 + \cdots + Y_n$, Proposition 3.3 in Puccetti and Rüschendorf (2012a) states that

$$\overline{\operatorname{VaR}}_{\alpha}^{r}(L) = \overline{\operatorname{VaR}}_{\alpha}^{\mathcal{E}}(L)$$
 and $\operatorname{VaR}_{\alpha}^{r}(L) = \operatorname{VaR}_{\alpha}^{\mathcal{E}}(L)$

for the particular class $\mathcal E$ introduced above. Therefore, we can rewrite (22) as

$$\underline{\mathrm{VaR}}_{\alpha}(L) \leq \underline{\mathrm{VaR}}_{\alpha}^{r}(L) \leq \mathrm{VaR}_{\alpha}(L) \leq \overline{\mathrm{VaR}}_{\alpha}^{r}(L) \leq \overline{\mathrm{VaR}}_{\alpha}(L).$$

A reduction method similar to the one described above has also been given in Puccetti and Rüschendorf (2012a) in the case of a general marginal system \mathcal{E} . The corresponding reduced bounds $\overline{\text{VaR}}_{\alpha}^{r}(L)$ and $\underline{\text{VaR}}_{\alpha}^{r}(L)$ however may fail to be sharp.

We illustrate how to calculate the bounds in (24) in two examples. First we assume the bivariate distributions $F_{2j-1,2j}$, $1 \le j \le n$ to be identical and generated by coupling two Pareto marginals having tail parameter $\theta > 0$ by a Pareto copula with parameter $\gamma \ne 0$. The bivariate Pareto copula with parameter $\gamma > 0$ is given by

$$C_{\gamma}^{Pa}(u,v) = ((1-u)^{-1/\gamma} + (1-v)^{-1/\gamma} - 1)^{-\gamma} + u + v - 1.$$

Under these assumption, the bivariate distribution function F_{12} is given by

$$F_{12}(x_1, x_2) = 1 + ((1+x_1)^{\theta/\gamma} + (1+x_2)^{\theta/\gamma} - 1)^{-\gamma} - (1+x_1)^{-\theta} - (1+x_2)^{-\theta},$$
(25)

while the n = d/2 random variables Y_j defined in (23) are identically distributed as

$$H(x) = P(Y_j \le x) = P(Y_1 \le x) = P(L_1 + L_2 \le x), j = 2, ..., n.$$

Here, we have that

$$H(x) = \int_0^x F_{2|x_1}(x - x_1) dF_1(x_1), \tag{26}$$

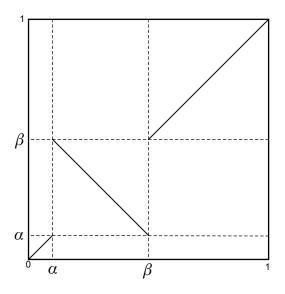
where we denote by $F_{2|x_1}$ the conditional distribution of $(L_2|L_1 = x_1)$. For this example, the conditional distribution $F_{2|x_1}$ is available in closed form and

$$F_{2|x_1}(x) = 1 - (1+x_1)^{\theta/\gamma + \theta} \left((1+x)^{(\theta/\gamma)} + (1+x_1)^{(\theta/\gamma)} - 1 \right)^{-\gamma - 1}.$$

Since the risk vector $(Y_1, \ldots, Y_n)'$ is homogeneous, we can apply the dual bound methodology introduced in Proposition 4 to compute $\overline{\text{VaR}}_{\alpha}^{r}(L)$ via (24a). In Proposition 4, we simply use n = d/2 (the number of the Y_r 's) instead of d and set F = H.

In Figure 9, we plot the unconstrained sharp VaR bound $\overline{\text{VaR}}_{\alpha}(L)$ and the reduced bound $\overline{\text{VaR}}_{\alpha}^{r}(L)$ for a random vector of d = 600 Pareto(2)-distributed risks under the marginal system described above. In the left figure the parameter of the Pareto copula is set to $\gamma = 1.5$. This implies a strong positive dependence between consecutive marginals. In the right figure we assume instead that the marginals are pairwise independent. The higher dimensional information reduces the conservative estimate of VaR in both cases, the larger reduction occurring in the case of the bivariate independence constraints. Recall that the calculation of the bound $\overline{\text{VaR}}_{\alpha}^{r}(L)$ in a homogeneous setting is independent from the dimensionality n of the risk vector $(Y_1, \ldots, Y_n)'$, confirming that the dual bound methodology is very effective for homogeneous settings. In Table 3 we compare the estimates for $VaR_{\alpha}(L_1 + \cdots + L_d)$ in the case of a homogeneous portfolio of Pareto(2) marginals and under different dependence scenarios.

In order to compute the improved bounds in (24) for *inhomogeneous* portfolios, one has to rely on the RA. We assume to have a portfolio of $d = 2n^2$ Pareto distributed risks, divided into n sub-groups of 2n risks. Risks within the same sub-group are assumed to be homogeneous, but risks in different sub-groups may have a different Pareto tail parameter. Within the i-th group, $1 \le i \le n$, we assume that each risk is Pareto(θ_i)-



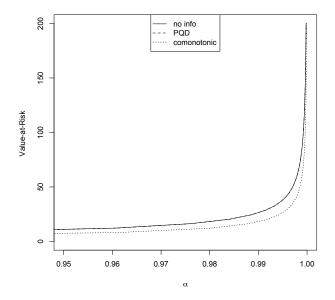


Figure 7: Bivariate copula of the vector (L_1^*, L_2^*) attaining the worst-VaR bound $M^{-1}(1-\alpha)$ under additional positive dependence restrictions (left). Values for $\overline{\text{VaR}}_{\alpha}(L)$ in the unconstrained case (no info), under additional positive dependence information (PQD) and comonotonic $\text{VaR}_{\alpha}^+(L)$ (see (8)), for the sum of two Pareto(2) distribution (right).

α	$\operatorname{VaR}^+_{\alpha}(L)$	$\overline{\operatorname{VaR}}_{\alpha}^{r}(L),$ (A)	$\overline{\operatorname{VaR}}_{\alpha}^{r}(L)$, (B)	$\overline{\mathrm{VaR}}_{\alpha}(L)$
0.99	5400.00	8496.13	10309.14	11390.00
0.995	7885.28	12015.04	14788.71	16356.42
0.999	18373.67	26832.2	33710.3	37315.70

Table 3: Estimates for $VaR_{\alpha}(L)$ for a random vector of d=600 Pareto(2)-distributed risks under different dependence scenarios: $VaR_{\alpha}^{+}(L)$ ($(L_{1},\ldots,L_{600})'$ has copula C=M); $\overline{VaR}_{\alpha}^{r}(L)$, (A): the bivariate marginals $F_{2j-1,2j}$ are independent; $\overline{VaR}_{\alpha}^{r}(L)$, (B): the bivariate marginals $F_{2j-1,2j}$ have Pareto copula with $\gamma=1.5$; $\overline{VaR}_{\alpha}(L)$: no dependence assumptions are made.

distributed and that the bivariate distributions $F_{2j-1,2j}$, $1 \le j \le$ *n* are of the form (25). A vector $\theta = (\theta_1, \dots, \theta_n)'$ then gives a full description of the marginals of the risk portfolio. The copula parameter is set to $\gamma = 1.5$ in each of the sub-groups. In Table 4, we give RA ranges for $\overline{\text{VaR}}_{\alpha}^{r}(L)$ and $\underline{\text{VaR}}_{\alpha}^{r}(L)$, as well as for $\overline{\text{VaR}}_{\alpha}(L)$ and $\underline{\text{VaR}}_{\alpha}(L)$ for different values of n, and at the quantile level $\alpha = 0.999$. In Table 4, computation times are indicated for the computation of the reduced bounds $\overline{\text{VaR}}'_{\alpha}(L)$ and $\underline{\mathrm{VaR}}_{\alpha}^{r}(L)$. These times are in general larger if compared to the homogeneous case with number of marginal distributions $d = n^2$. Indeed, in order to apply the RA to the marginals H_i , one has to compute the quantiles of the distribution H in (26) which is in general a more time consuming operation especially considering that one has to handle different tail parameters. If one has an efficient procedure to obtain these latter quantiles, then the RA computation times of the reduced VaR intervals are approximately the same as in the homogeneous case with $d = n^2$ marginal distributions.

To summarize, the same techniques introduced in Section 2, where one only knows the marginal distributions of the risk vector $(L_1, \ldots, L_d)'$ can be applied to the case where higher dimensional information is available. In order to use the reduction method one only needs to have the conditional distribution function $F_{i|x_1}$ available in closed form, for any $x_1 \in \mathbb{R}$. This conditional distribution is typically available for bivariate models derived from continuous marginals and a continuous copula, but it might be difficult to compute for higher dimensional subgroups of marginals.

Worst-case dependence structures for the problems (21) are in general not available. However, some approximation results given in Embrechts and Puccetti (2010a, Section 5) indicate that they still contain a minimal variance component.

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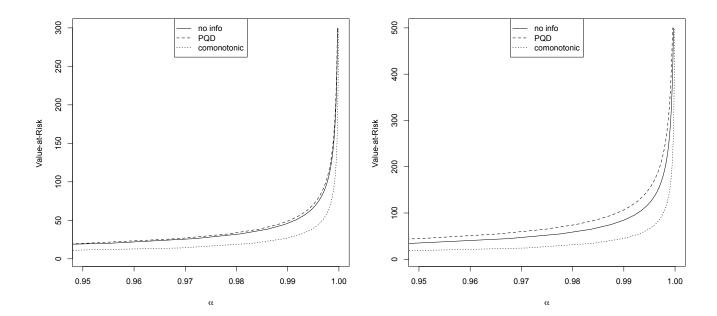


Figure 8: Values for $\overline{\text{VaR}}_{\alpha}(L)$ in the unconstrained case (*no info*), under additional positive dependence information (*PQD*) and comonotonic $\overline{\text{VaR}}_{\alpha}^{+}(L)$ (see (8)) for the sum of d = 3 (left) and d = 5 (right) Pareto(2) marginals.

			$\underline{\text{VaR}}_{0.999}(L)$	$\underline{\operatorname{VaR}}_{0.999}^{r}(L)$	VaR _{0.999} ⁺ (L)	$\overline{\text{VaR}}_{0.999}^{r}(L)$	$\overline{\text{VaR}}_{0.999}(L)$
n	d	comp. time	(RA range)	(RA range)	(exact)	(RA range)	(RA range)
2	8	4 mins	30.47 - 30.62	55.11 - 55.40	158.49	226.07 - 226.11	277.27 - 277.28
5	50	1.5 hrs	30.47 - 30.62	55.15 - 55.44	652.92	1024.36 - 1024.47	1152.64 - 1152.90
18	648	4 hrs	339.86 - 340.00	341.09 - 341.22	7373.01	11408.63 - 11439.61	12626.80 - 12695.52

Table 4: Estimates for $\text{VaR}_{\alpha}(L_1 + \cdots + L_d)$ for random vectors of Pareto-distributed risks with different tail parameters. The vector of tail parameters are $\theta = (2, 3)'$ (first portfolio), $\theta = (2, 2.5, 3, 3.5, 4)'$ (second portfolio) and $\theta = (2, 2.125, \dots, 4, 4.125)'$ (third portfolio). Under the additional dependence scenario, the bivariate marginals $F_{2j-1,2j}$ of the risk vector have Pareto copula with $\gamma = 1.5$. We set N = 1.0e05 in the first two portfolios and N = 5.0e04 in the last.

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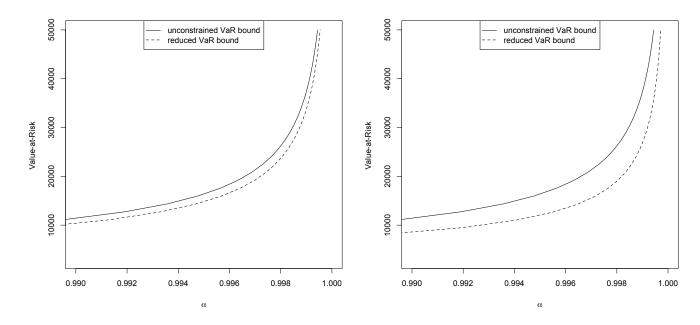


Figure 9: VaR bounds $\overline{\text{VaR}}_{\alpha}(L)$ (see (5)) and reduced bounds $\overline{\text{VaR}}_{\alpha}^{r}(L)$ (see (24a)) for a random vector of d = 600 Pareto(2)-distributed risks with fixed bivariate marginals $F_{2j-1,2j}$ generated by a Pareto copula with $\gamma = 1.5$ (left) and by the independence copula (right).

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