

Informative Risk Measures in the Banking Industry: A Proposal based on the Magnitude-Propensity Approach

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Abstract

Despite decades of research in risk management, most of the literature has focused on scalar risk measures (like e.g. Value-at-Risk and Expected Shortfall). While such scalar measures provide compact and tractable summaries, they provide a poor informative value as they miss the intrinsic multivariate nature of risk. To contribute to a paradigmatic enhancement, and building on recent theoretical work by [Faugeras and Pagès, 2024], we propose a novel multivariate representation of risk that better reflects the structure of potential portfolio losses, while maintaining desirable properties of interpretability and analytical coherence. The proposed framework extends the classical frequency–severity approach and provides a more comprehensive characterization of extreme events. Several empirical applications based on real-world data demonstrate the feasibility, robustness and practical relevance of the methodology, suggesting its potential for both regulatory and managerial applications.

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1 Introduction and motivation

The pursuit of effective risk measures in the financial industry is an intriguing and long-standing endeavor that has garnered significant attention from both scientific and regulatory perspectives. An adequate comprehensive risk management process typically consists of the following key steps: (i) Identify risks, to detect all sources of risk; (ii) Assess risks, i.e., measure them; (iii) Manage risks, practically to avoid, transfer, hedge, or set limits and warnings; (iv) Monitor and review risks, with ongoing activity. While some steps, such as identification and management, relate to business knowledge and the enterprise’s organization and processes, the measurement step (ii) is highly technical, involving several related challenging components, such as: the choice of a "good" risk measure, the setup of its parameters, a probabilistic model for random adverse events, the statistical calibration of model parameters, and finally, the statistical estimation of risk measures given empirical data. From a conceptual perspective, the choice of risk measures, i.e., how to represent or summarize the randomness of business variables (Profits and Losses in financial markets, Credit Default events, Operational Losses), is a highly intriguing task, as the adopted risk measures should satisfy certain quantitative and qualitative properties. Briefly, given a univariate random variable X describing the potential profits and losses of a portfolio (PnL for brevity), one aims to summarize in a single scalar value, the risk measure, the uncertainty that the bank (or insurance companies) must manage. It is noteworthy

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that the risk evaluation goal is often combined with return evaluation, as most capital allocation choices are driven by the objective of optimizing the risk-versus-return trade-off in some sense. Extensive work has been devoted in corporate finance research to topics such as *RORAC* (Return on Risk-adjusted Capital) and *RAROC* (risk-adjusted return on capital) indicators, i.e., how to allocate capital, the scarce resource, by seeking the best allocation on the efficient frontier concerning the enterprise's risk appetite. A seminal reference in this field, recently updated, is [Brealey et al., 2019]. Adopting these models requires reliable input in terms of (expected) return and risk estimation. While return on an investment primarily poses statistical forecasting challenges, capturing risk concisely for a portfolio is a more arduous task involving subtle mathematical and conceptual points. We recall that risk, namely the profits and losses behavior, is typically described by a random variable. Furthermore, in many real-world models, the distribution of this random variable is unknown, as it is the outcome of a sophisticated model involving several risk factors, non-linear relationships, non-Gaussian elementary distributions, and so on. This extends the problem from risk quantification to model uncertainty, where the distribution itself is not certain, usually known as *model risk*, see [Danielsson et al., 2016] for a deep empirical survey. An extensive literature on risk measures exists, stemming from both academic research and financial practitioners. The subtle distinction between risk and uncertainty, along with a review of risk measures, is provided in [Biglova et al., 2008]. The banking and insurance sectors exhibit a high level of regulation; hence, regulators (i.e., governments, central banks, banking authorities) have devoted significant effort to defining appropriate risk measures to disclose banks' risks to the market and all stakeholders. In particular, for over a decade now, international regulations and authorities have been emphasizing the concept of a *Risk Appetite Framework* (RAF), which banks are required to establish, and the promotion of Risk Culture. In this context, the concepts of *Risk Capacity* (or *Tolerance*) – the maximum risk that can be borne, especially *Risk Appetite* – the desired risk level, and Risk Profile – the actual risk over time, are crucial, see [European Central Bank, 2021] (specifically the "Definitions" on page 15), and the Basel Committee's paper No. 328 for further guidance. Risk culture pertains to the dissemination and awareness of risk among the bank's top figures (Board, Top Management), cascading throughout the organization, encompassing risk assumptions, risk types, models, and measures. Section 9 of the [European Central Bank, 2021] guidelines provides insights on this aspect. From a practical standpoint, the RAF is further articulated into a *Risk Appetite Statement* (RAS), which contains high-level (strategic) indicators, known as Tier 1, for various risk types, followed by managerial indicators (Tier 2), and finally, operational or warning indicators at Tier 3 level. Of course, different regulations prescribe different risk measures. The general ambitious goal of a risk measure is to summarize the risky side of the uncertainty concept, i.e., how much extreme losses could affect the expected return of any investment strategy. As the extreme losses must be faced by the own capital of the bank (or the insurance company), it is quite common to say that the risks absorb capital. At a very general level, we define X_θ as the random variable describing the phenomenon to be investigated, where (θ) represents the parameter (scalars or arrays) characterizing the variable. Let $F(x; \theta)$ be its cumulative distribution, i.e., $F(x; \theta) = P(X_\theta \leq x)$. More generally, let \mathcal{F} be the set of all random variables on the real space (typically, $\mathcal{F} = L^\infty(\mathbb{R})$). Following [Embrechts et al., 2015b], the ways to measure risk can be grouped into four categories: the notional-amount approach, sensitivity measures, risk measures based on scenarios (*stress tests*), and risk measures based on the *PnLs* distribution. Focusing on the last category, most portfolio risk measures are statistical quantities describing the conditional or unconditional loss distribution over some predetermined horizon. The most popular include volatility σ , *Value-at-Risk* (VaR), and *Expected Shortfall* (ES). More formally, a risk measure is an application $\rho(X) : \mathcal{F} \rightarrow \mathbb{R}_+$ mapping the random variable X into a positive scalar value, representing its uncertainty. A first point is to recall that risk relates to uncertainty, but in most situations, banks focus on downside risk, i.e., the adverse side of risk, where extreme losses in the business might occur with a given frequency. For this reason, volatility σ (estimated standard deviation of portfolio returns), although widely popular due to its simplicity and ubiquitous use in the asset management field to rank products by their risk, is not adopted in financial regulation, as it captures randomness from a symmetric perspective. To define

what a good risk measure is, the notion of *coherent* risk measures has become prevalent, as outlined in the seminal paper by [Artzner et al., 1997], see the next section for a detailed overview. If we change perspective, we could wonder if the traditional approach, i.e., seeking "the best" risk measure, is the proper approach, or if an alternative strategy should be adopted. Can a one-dimensional approach, i.e., a scalar measure summarizing adverse *PnLs* results, truly provide sufficient information about risk? Or should we move to a multidimensional set of values to represent it? To apply this perspective, which key theoretical or conceptual ideas need to be adapted? Are there alternative approaches to risk management beyond the search for an appropriate scalar risk measure? Describing risk by a couple (or a few) indicators is quite common in fields other than financial risk management. In the insurance sector, particularly in the incident claims area, it is popular to summarize each event category by two parameters: the frequency (λ) and the claim amount (S). If the two are independent or approximately independent, one can easily obtain the expected value of the claims to be managed as $C = \lambda \cdot S$. The loss (L) in a given time horizon (T) is represented in the simplest case by a Compound Poisson process, namely: $L(T) = \sum_{i:t(i) \leq T} S(i)$, where the sum of random losses (lognormal, gamma, or other appropriate random variable) is extended to a random number of events driven by the Poisson random variable. A similar approach prevails in the operational risk sector in banks, i.e., the risk of losses arising from incidents, errors, failures, or fraud. In the Loss Distribution Approach (LDA), a random number of negative events are recorded over time, each with its severity. Even in the simplest lognormal approach for severity, we do not know the exact closed-form distribution of a sum of lognormal random variables, *a fortiori* if the number of events itself is random. To this extent, several analytical approximations have been suggested, some from the mathematical finance field, see e.g. [Turnbull and Wakeman, 1991], others specifically developed in the operational risk context, like e.g. [Peters and Shevchenko, 2015]. If we move to enterprise risk management, even in large corporate contexts, the methodology relies less on quantitative tools, as expert assessment is the most popular approach. Once the risks in the company have been identified and listed, for each risk, the process owner or the expert panel is asked to complete a questionnaire to assign some Key Performance Indicators (KPIs), such as: (i) The probability that no losses are observed; (ii) The expected loss in normal situations; (iii) The magnitude of losses in extreme cases, where "extreme" is qualified by some percentile level or as the worst case over a very long horizon, etc. A detailed guide to this approach is outlined in [Institute of Operational Risk, 2019]. To summarize the above alternative approach and to avoid any confusion, we point out that while the definition of a good (scalar) risk measure focuses on the final output of uncertainty, i.e., the extreme losses for the institution, the risk representation by a few parameters (probability of the event, its severity, etc.) concerns the components, i.e., the inputs to the final risk measure. The general idea underlying our work is to combine these two aspects, namely the classical scalar approach generally adopted in practice for defining a risk measure in the risk management process, along with the multivariate perspective based on its typical parameters on the input side.

From a theoretical standpoint, we refer to the recent work on the magnitude-propensity approach in risk management by [Faugeras and Pagès, 2024]. The authors' groundbreaking work employs two closely related approaches in their theoretical framework: optimal transport and quantization techniques. Their methodology can be distilled into three key points. Firstly, they argue that traditional single-valued risk measures, such as *VaR* or *ES*, are inadequate for fully capturing the underlying uncertainty. Secondly, they acknowledge the binary nature of losses - they either occur or they don't. Lastly, they propose approximating the original risk model (represented by the random variable X) with a binary variable, which offers a more nuanced perspective on risk, balancing simplicity with a richer representation of uncertainty. This simplified model is characterized by two parameters: p , the probability of incurring no losses, and m , the magnitude of losses when they do occur (with probability $1 - p$). Such distribution approximation problem is mathematically formulated as mass transportation in Wasserstein metric space. The optimal transportation to a discrete measure is analogous to the optimal quantization problem, a well-established concept in Engineering and Signal Processing literature. Consequently, the proposed approach for quantifying risk on both magnitude and propensity scales can be characterized as a specialized, constrained optimal quantization problem. The

formulation of [Faugeras and Pagès, 2024] provides a rigorous mathematical framework for risk assessment, bridging the gap between theoretical optimal transport and practical risk quantification. It leverages the Wasserstein metric’s properties to capture the multidimensional nature of risk, while the quantization aspect ensures computational tractability and interpretability of the results. We immediately note that the approach of [Faugeras and Pagès, 2024] differs from that of vector risk measures, first introduced by [Jouini et al., 2004] and later revisited and developed by [Ben Tahat and Lépinette, 2014], [Feinstein and Rudloff, 2015], and [Ararat and Feinstein, 2024], in order to account for proportional transaction costs in multi-asset markets.

Numerous attempts have been made to introduce additional properties aimed at improving these measures, identifying desirable characteristics, and clarifying what constitutes a well-defined notion of risk at the vector level. However, extending the same axioms used for scalar coherence often renders vector risk measures difficult to interpret and, in particular, to implement, since an additional selection procedure is typically required to determine a specific capital requirement or allocation rule within the set-valued framework, see [Ararat and Feinstein, 2024] for a recent overview of these methods. The multidimensional approach we adopt here is of a different nature and not directly comparable to theirs.

Regarding our contribution to this field, from a theoretical point of view, we endeavor to extend the framework of [Faugeras and Pagès, 2024] to the 3-point (i.e., zero, moderate, and extreme losses). In this multidimensional context, this contribution is significant. While the Risk Appetite Framework encompasses not only market and credit risk (the focus of our examples) but also liquidity, operational, reputational, and other risks, one of the pressing issues for banks, as requested by the European Central Bank during inspections, is "How did you determine the set of limits, i.e., the level of capacity/tolerance or warning levels?". There is no common practice or guidelines on this matter. Most banks adopt the following approach: they consider the historical Value-at-Risk data, take the highest value, perhaps add a 10 – 20% buffer, and set that as the tolerance level (keeping in mind that, at the bank level, as opposed to sub-portfolios, there may also be regulatory limits). In this regard, our optimal upper threshold m_2 corresponding to extreme losses could serve as an excellent "automatic machine" (recall that ECB highly appreciates objectivity in the determination of the thresholds) to determine and support the setting of tolerance and/or capacity levels. For risks that involve only losses (operational, financial), the lower threshold m_1 , corresponding to moderate losses, could also be an equally automatic tool for defining the risk appetite as observed in the past or as a "floor" for such appetite. In the context of Risk Culture and top management’s awareness of risks, a description using (m_1, m_2) with the associated probabilities (p_1, p_2) could provide an intuitive and complementary alternative to VaR and, more importantly, to the more complex Expected Shortfall, due to its familiarity with concepts such as severity and frequency. Finally, for "financial conglomerates" under European directives, i.e., groups that include banks, insurance companies, and asset management firms, the regulations mandate the integration (effectively, uniformity and convergence) of measurement methods and risk culture across the various legal entities within the group. Our approach, which combines ideas from finance and insurance logic, could be a viable solution in this context.

Furthermore, our contribution is also focused on applications, as we discuss certain constraints that must be incorporated into the optimization problem to render the solution more realistic in practical contexts. We apply the technique to the historical simulation approach for *VaR* calculation and to the Montecarlo simulation for the Default Risk Charge (DRC) measure, which are largely the most adopted by major banks, precisely 19 out of 31 in the 2021 ECB review, see [European Central Bank, 2021]. To accomplish our objective, we utilize the *PnLs* derived from the real portfolio of a significant European bank. This makes our work complementary to [Faugeras and Pagès, 2024], which primarily pertains to the insurance sector, where we implement the third application case. The paper is organized as follows: Section 2 reviews the relevant literature on risk measures; Section 3, after a brief background on quantization and the magnitude-propensity approach, elucidates the extensions we implement to make this proposal more suitable in the field of risk management; Section 4 illustrates the application and the dataset employed for calibrating the models; Section 5 contains our results, along with a discussion of the key points; Section 6 summarizes the conclusions and

outlines further research avenues. We gather in the Appendix the technical proofs and some material on the numerical optimization procedures.

2 Scalar Risk Measures: Review and Limitations

Risk quantification and measurement constitute essential elements in the domains of finance, insurance, and decision-making processes. Throughout the years, a diverse array of risk measures has been proposed and analyzed, each exhibiting distinct advantages and limitations. Coherent risk measures represent a fundamental and well-established framework for quantifying risk in financial and insurance contexts. These measures adhere to a set of axioms that ensure a consistent and rational approach to risk assessment. The coherence properties, including subadditivity, monotonicity, translation invariance, and positive homogeneity, provide a robust foundation for comparing and aggregating risks across diverse portfolios and financial instruments, see e.g. [Artzner et al., 1997]. The field of risk measurement has witnessed the development and application of numerous theories, ranging from the widely utilized Value-at-Risk and Expected Shortfall to *Stress Tests*. Researchers and practitioners have extensively explored these methodologies to evaluate potential losses and manage uncertainties in various financial contexts. *VaR*, a ubiquitously employed risk metric within the financial sector, provides a probabilistic estimate of potential losses at a predetermined confidence level over a specified temporal horizon. This measure enables financial institutions to quantify and articulate downside risk with precision, thereby facilitating informed decision-making and risk management strategies. The diffusion of *VaR* lies in its ability to distill complex risk profiles into a single, comprehensible figure, rendering it a useful tool for risk communication among stakeholders, regulatory compliance, and internal risk control mechanisms. For a given holding period and a level $\alpha \in (0, 1)$ the *VaR* of the *PnL* distribution X is defined as

$$VaR_\alpha(X) := -\inf\{x \in \mathbb{R} : \mathbb{P}(X \leq x) > \alpha\},$$

or, equivalently, $VaR_\alpha = \inf\{x \in \mathbb{R} : \mathbb{P}(Loss \geq x) \leq \alpha\}$. One of the primary merits of *VaR* lies in its simplicity and interpretability, offering a lucid and succinct estimation of potential financial losses. This characteristic enables risk managers to swiftly evaluate and juxtapose risks across diverse portfolios or trading positions. Nevertheless, *VaR* is not without its limitations, particularly in its sensitivity to extreme events, where its single-point estimate may inadequately capture tail risk. The work of [Embrechts et al., 2015b] elucidates how *VaR* embodies the propensity aspect of risk by determining the leftmost quantile, yet fails to encapsulate the magnitude of potential losses. Furthermore, *VaR* has been subject to substantial criticism due to its *non-subadditivity*, a property that contravenes the axioms of coherent risk measures. This deficiency renders the aggregation of *VaR* values across portfolios problematic. To illustrate this point, one can readily construct an example utilizing two sub-portfolios, P_1 and P_2 , where the following inequality holds: $VaR_\alpha(P = P_1 \cup P_2) > VaR_\alpha(P_1) + VaR_\alpha(P_2)$. This mathematical representation underscores the inherent limitations of *VaR* in accurately reflecting the cumulative risk of combined portfolios, thereby highlighting the need for more robust risk measurement methodologies in complex financial environments.

Conversely, the essential supremum, denoted as $\rho_\infty(X) := \text{ess sup } X$, represents the antithetical extreme to *VaR*, quantifying the maximal magnitude of a potential loss without estimating the associated probabilities. This risk measure, also known as the worst-case scenario or *tail Value-at-Risk*, encapsulates the most extreme potential loss, irrespective of its probability of occurrence. As such, it provides a conservative estimate of the worst possible outcome but does not offer a comprehensive risk assessment that incorporates both magnitude and propensity aspects. One of the primary advantages of $\rho_\infty(X)$ lies in its simplicity and interpretability. By focusing solely on the maximum potential loss, it enables risk managers to identify worst-case scenarios and allocate capital reserves accordingly. In situations where extreme events can have severe consequences, $\rho_\infty(X)$ provides a valuable upper bound for risk exposure, ensuring that institutions are adequately prepared for the most adverse outcomes. However, this extreme focus on maximum loss is

not without drawbacks. By disregarding probabilities, $\rho_\infty(X)$ neglects the likelihood of less extreme but still significant losses. Consequently, risk managers relying exclusively on this measure might overlook the impact of moderately severe yet more probable events, potentially leading to suboptimal risk management strategies. Furthermore, $\rho_\infty(X)$ exhibits high sensitivity to outliers and extreme observations, rendering it vulnerable to estimation errors and model mis-specifications. In practice, financial data often display heavy-tailed distributions, implying that extreme events occur more frequently than a Normal distribution would suggest. As a result, $\rho_\infty(X)$ might overstate worst-case scenarios, leading to excessively conservative risk assessments and potential over-allocation of capital reserves. The theoretical underpinnings and practical implications of the essential supremum in risk measurement have been extensively explored in the seminal works of [Rockafellar and Uryasev, 2002] and [Pichler and Shapiro, 2021].

Expected Shortfall (*ES*), also known as Conditional Value-at-Risk, emerges as the most promising alternative to address *VaR*'s limitations. *ES* quantifies the expected value of losses beyond the *VaR* threshold, providing information on extreme event severity and tail behavior. It is formally defined as:

$$ES_\alpha = \mathbb{E}[Loss | Loss \geq VaR_\alpha].$$

Expected Shortfall exemplifies a coherent risk measure, adhering to the fundamental axioms of coherent measures (translation invariance, subadditivity, positive homogeneity, and monotonicity, see [Artzner et al., 1997]). By quantifying potential losses beyond the *VaR* threshold, *ES* provides a more comprehensive and robust risk assessment, particularly for heavy-tailed distributions, see [Acerbi and Tasche, 2002]. This attribute renders *ES* preferable in financial risk management, where tail events can precipitate significant systemic consequences. Regulatory bodies and financial institutions increasingly adopt *ES*, acknowledging its capacity to address *VaR*'s limitations and more accurately represent extreme risk scenarios. Then *ES* is going to replace *VaR* ($\alpha = 99\%$, $h = 10$ days) in the impending Fundamental Review of the Trading Book (FRTB) regulation for market risk minimum capital requirement calculations, see [Basel Committee of Banking Supervision, 2019]. In the forthcoming regulatory framework, Expected Shortfall is employed to quantify market risk at a $\alpha = 97.5\%$ confidence level and different time horizons h ($h = 10, 20, 40, 60, 120$ days depending on the liquidity of the risk factors categories). However, financial institutions are still required to validate their models using the back-testing procedure based on *VaR* calculated at both $\alpha = 99\%$ and $\alpha = 95\%$ confidence levels. This unconventional regulatory setup stems from a decade-long debate surrounding the *elicitability* property of risk measures. Elicitability, derived from point forecasting, requires statistical functionals to maintain consistency with their evaluation through historical averaging, akin to M -functionals (see [Gneiting, 2011]). In this context, *VaR* is elicitable but not coherent, while Expected Shortfall is coherent and law-invariant but not elicitable. Notably, *ES* exhibits joint elicibility with *VaR* ([Fissler and Ziegel, 2016]). Furthermore, *ES* possesses the crucial property of backtestability, enabling rigorous performance evaluation in historical simulations and real-world scenarios, as elucidated by [Acerbi and Székely, 2017], who underscore *ES*'s capacity to provide robust and reliable risk assessments, rendering it an invaluable tool for risk managers in evaluating potential losses across diverse financial scenarios¹. More recently, the competition among the different proposals for good risk measures has been enriched by some works about the concept of *observability* of a risk measure. Quite surprisingly, the popular volatility indicator and the *ES* show to be more observable than *VaR*. To this extent, an insightful analysis is provided in [Acerbi and Székely, 2023], by leveraging the concept of *sharp backtest*.

Finally, it is worth to note that in the FRTB regulation the banks must also calculate another risk measure, the *Default Risk Charge* (DRC), namely a 1-Year *VaR* with $\alpha = 99.9\%$. This capital requirement aims to capture the default risk in the trading book. Until the 2008-2011 financial crisis, the so called *market*

¹Our magnitude-propensity-based approach can be subjected to backtesting by extending the regulatory framework prescribed for *VaR*, see [Kupiec, 1995], as well as the more sophisticated joint *VaR*–*ES* testing procedures proposed by [Acerbi and Székely, 2017]. A comprehensive analysis of backtestability and the extension of the notion of elicibility to our multivariate framework requires a dedicated investigation and will therefore be developed in a forthcoming paper.

risk building block in the Basel regulation included only the *spread risk*, i.e. the price uncertainty coming from the spread level, e.g in the bond evaluation. According to the Basel 2.5 Reform, it was requested to estimate the potential losses coming from the default (or migration) events also for the trading book of the bank, see [European Banking Authority, 2012]. While the *PnLs* of the traditional market risk come from the movements of the risk factors (equities, interest rates, forex rates, spreads, commodities) that are commonly assumed to have continuous distribution, in the credit risk case the portfolio losses are originated by the default binary events D_n associated to the obligors ($n = 1, \dots, N$):

$$Loss_{DRC} = \sum_n EAD_n \cdot \mathbf{1}_{D_n} \cdot LGD_n,$$

where EAD_n and LGD_n represent, respectively, the amount of exposure and the fraction that the bank is not able to recover once the default happens, modeled as a deterministic coefficient in the range $[0, 1]$, or by a random variable, usually a *Beta* (a, b). It is important to notice that the combinatorial effects associated with default events lead to an intrinsic instability of any risk measure, due to the fact that extreme values of the distribution are so pronounced that the VaR no longer provides an adequate representation of tail risk. We will illustrate this effect in the Appendix B.

Stress tests represent a pivotal tool for risk analysis, enabling institutions to assess resilience against adverse market conditions and extreme events. These tests simulate hypothetical crisis scenarios, such as economic recessions, market shocks, or event-related crises, by subjecting portfolios or financial systems to various stress factors. By modeling extreme conditions, stress tests identify vulnerabilities and quantify potential losses under adverse circumstances. The results provide insights into an institution's risk exposure, capital adequacy, and overall financial stability. This risk management tool has become increasingly crucial for regulatory compliance and is widely adopted by financial institutions to ensure robust risk management practices. Notably, stress tests have been instrumental in enhancing the financial sector's resilience following the 2008 global financial crisis. By simulating scenario outcomes, stress tests empower institutions to implement appropriate risk mitigation strategies and maintain financial stability in turbulent times. From a conceptual standpoint, stress tests diverge significantly from measures such as Value-at-Risk and Expected Shortfall. On one hand, stress tests are more objective in their results and are based on a simple "*what-if*" logic: given a scenario, banks can calculate the related *PnL* with high accuracy. On the other hand, regulations mandate the use of *extreme yet plausible* scenarios, a goal that is arduous to achieve even for regulatory authorities. Furthermore, stress test results lack information about the associated probabilities.

In this respect, the common practice consists of some heuristic methods. To determine thresholds for risk tolerance or risk appetite in the their risk appetite framework, typically some banks consider the past time average of the Value-at-Risk and increase it by some subjective quantity (e.g. 20%), other banks calculate directly the *Stressed VaR*, i.e. *VaR* under stressed parameters. Such approach is quite reasonable but exhibits some drawbacks, being too judgmental without a rational basis. Furthermore, these methods do not add any relevant additional information about risk beyond what is provided by the *VaR* itself. Additional information is required, which only a multidimensional approach can provide, as will be explained in the following section. *VaR* does not provide insights into the underlying risk factors or the potential severity and frequency of extreme events. By simply scaling the *VaR* up or down by a fixed percentage, the resulting stressed scenarios fail to capture the complex dynamics and interdependencies of risk factors, as well as the potential for tail events that may deviate significantly from the assumed distribution. A multidimensional approach, on the other hand, can incorporate additional risk dimensions, such as severity and frequency, as well as the potential for non-linear interactions between risk factors. In the following section, we will explore a multidimensional approach that incorporates severity and frequency dimensions to provide a more comprehensive and informative framework for risk management and scenario analysis.

3 From Scalar to Vectorial Risk Measures and from Continuous to Discrete Distributions

[Faugeras and Pagès, 2024] observe that VaR and ES are traditionally calculated for popular confidence levels α (e.g., 95%, 99%, 97.5%), introducing a degree of subjectivity to these risk measures and highlighting potential weaknesses in comparing distinct PnL s distributions. The magnitude-propensity approach aims to address these limitations. Given the random variable for the PnL , say X , the approach consists in defining a discrete binary random variable Y that works as an informative summary of the whole X , whose distribution is given by

$$P^Y = (1 - p) \cdot \delta_0 + p \cdot \delta_m,$$

where p is the probability that the loss is zero, m represents the magnitude of the loss and δ_x denotes the Dirac distribution concentrated at the point x . Various methodologies have been developed to discretize continuous risk distributions, providing decision-makers with robust tools for risk assessment and strategic planning. Two main approaches emerge in the literature - the *optimal transport* approach and the *optimal quantization* approach. The *optimal transport* method approximates continuous distributions by minimizing mass transportation costs between distributions, as explained by [Villani, 2009]. On the other hand, *optimal quantization*, described by [Graf and Luschgy, 2000], [Pagès et al., 2004] and [Luschgy and Pagès, 2023], represents risk data using finite point grids².

While both the optimal transport and quantization approaches aim to disentangle the magnitude and propensity aspects of risk, they employ distinct mathematical frameworks to achieve this goal. The optimal transport approach represents the search for the binary random variable Y as a mass transportation problem, whereas the quantization approach treats it as an optimal discrete representation problem. However, optimal transportation to discrete measures also corresponds to a special case of optimal quantization, in fact the proposed approach to quantify risk on both the magnitude and propensity scales amounts to a special, constrained, optimal quantization problem. Specifically, we briefly review the formulation of the bivariate magnitude-propensity risk measure and its optimization using *Wasserstein distance* for the optimal transport approach. Furthermore, we explore the intricacies of the optimal quantization problem and its relevance in approximating risk distributions with discrete measures. We then attempt to establish a solid theoretical foundation for our research, which aims at extending the original 2-points distribution framework of [Faugeras and Pagès, 2024] by developing a novel 3-points distribution approach. Our extended methodology incorporates an additional quantile, enabling a more informative and comprehensive risk quantification process.

3.1 Optimal Transport Approach

The optimal transport (OT) problem, formally introduced by [Kantorovich, 1942], addresses the transformation of a probability distribution P^X of the random variable X to a distribution P^Y of the variable Y , while minimizing the associated transition cost function $c(x, y) : X \times Y \rightarrow [0, +\infty]$. The Monge-Kantorovich optimal transportation problem has a clear physical interpretation: considering the random variables as material locations, then $c(x, y)$ represents the cost of transporting a unit of material from x to y . Optimal cost functions typically represent the transport cost as the product of mass and inter-location distance. The problem is constrained by complete material relocation. This framework has diverse applications, spanning economics, image processing, and machine learning, providing a robust methodology for distribution transformation analysis and optimization. The formulation maintains scientific rigor while offering concise

²We also mention other methods like K -means clustering, studied by [Liu and Pagès, 2020], which groups similar risk data points, then revealing underlying risk patterns, and the quantile-based methods, explored by [Chernozhukov and Hong, 2004], that partition risk distributions based on predefined quantiles, focusing on specific risk thresholds.

elegance, emphasizing the concept's broad academic and practical relevance. This version further condenses the information while preserving the scientific tone and elegant expression. It highlights the key points about cost function formulation, problem constraints, and the wide-ranging applicability of the optimal transport framework. Mathematically, the Monge-Kantorovich optimal transport problem is formulated as:

$$\inf_{\pi \in \Pi(P^X, P^Y)} \int_{\mathbb{R} \times \mathbb{R}} c(x, y) \pi(dx, dy)$$

where $\Pi(P^X, P^Y)$ is the set of all transport plans, i.e., joint probability measures on $\mathbb{R} \times \mathbb{R}$ with marginals P^X and P^Y , respectively. Under regularity conditions, the optimal transportation plan is defined as a Monge map T , namely $\pi(x, y) = \pi(x, T(x))$, see e.g. [Rachev and Rüschendorf, 1998], [Villani, 2009]. In the degenerate case where $P^Y = \delta_m$, with $m \in \mathbb{R}$, $\Pi(P^X, P^Y)$ reduces to the singleton product measure $\{P^X(dx) \times \delta_m(dy)\}$ and the OT problem becomes

$$\inf_{m \in \mathbb{R}} \int c(x, m) P^X(dx) = \inf_{m \in \mathbb{R}} \mathbb{E}[c(X, m)].$$

For $c(x, y) = (x - y)^2$ one gets the squared L^2 -Wasserstein metric W_2 and the OT problem reads

$$W_2^2(P^X, \delta_m) = \inf_{m \in \mathbb{R}} \mathbb{E}[(X - m)^2] = \text{Var}(X),$$

which is minimized for the mean $m = \mathbb{E}[X]$. For the L^1 distance $c(x, y) = |x - y|$ one gets the median, while for the asymmetric cost $c(x, y) = (x - y)\alpha 1_{x-y \geq 0} + (y - x)(1 - \alpha)1_{y-x > 0}$ with $0 < \alpha < 1$ one gets the (left)- α quantile $m = q_\alpha(X)$, i.e. the Value-at-Risk (see [Koenker, 2005]). Finally, for $c(x, y) = y + \frac{(x-y)1_{x \geq y}}{1-\alpha}$ one gets the Conditional Value-at-Risk, namely the ES, see [Rockafellar and Uryasev, 2002].

In the magnitude-propensity approach investigated by [Faugeras and Pagès, 2024], the optimal transport approach scrutinizes traditional risk measure limitations by reframing them as mass transportation from the original risk distribution P^X to a binary distribution P^Y , encapsulating both risk magnitude and propensity. The corresponding risk measure (m, p) is derived by minimizing the Wasserstein W_2 -distance between P^X and P^Y within a constrained distribution set $\mathcal{P}_0 := \{P^Y = p\delta_m + (1 - p)\delta_0, p \in (0, 1), m \in \mathbb{R}_+\}$. This methodology provides a novel perspective on risk quantification, integrating the multidimensional nature of risk into a cohesive framework, where a loss of magnitude m occurs with probability p , and no loss (i.e. the loss amount equals zero) occurs with probability $(1 - p)$. It then offers a more nuanced approach to risk assessment, potentially enhancing decision-making processes in financial and economic contexts. If we denote with Q_X (resp. Q_Y) the quantile function associated with the distribution P^X (resp. P^Y), then the L^2 -Wasserstein W_2 -distance between P^X and P^Y reads (see e.g. [Rachev and Rüschendorf, 1998]):

$$W_2(P^X, P^Y) = \left(\int_0^1 (Q_X(x) - Q_Y(x))^2 dx \right)^{1/2}$$

where $Q_X(z) := \inf\{x : F_X(x) \geq z\}, 0 < z < 1$ and F_X represents the cumulative distribution function of the distribution P^X , where X is assumed to have finite variance. For the two-point distribution P^Y , the quantile function reads $Q_Y(z) = m1_{1-p < z \leq 1}$, and the optimal quantities (m, p) can be found explicitly by direct optimization, see [Faugeras and Pagès, 2024]. We omit the details of their results as our work employs the quantization method, which proves to be significantly more efficient in our context. In Appendix C we provide some additional details on the algorithmic implementation of the Optimal Transport procedure that we adopted in our analysis.

3.2 The Quantization Approach

Quantization, rooted in engineering and signal processing, offers an alternative paradigm for achieving analogous objectives. This approach involves the optimal discretization of continuous risk distributions, akin to analog-to-digital conversion and data compression methodologies. The primary aim is to identify a finite set of points (codebook) that minimizes the mean squared error or distortion between the original risk distribution P^X and its quantized counterpart P^Y . This process effectively transforms complex, continuous risk landscapes into more manageable, discrete representations.

Following [Graf and Luschgy, 2000] and [Luschgy and Pagès, 2023], an N -vector quantizer on $(\mathbb{R}^d, \|\cdot\|)$ is a mapping $T : \mathbb{R}^d \rightarrow \{x_1, \dots, x_N\}$ where $\{x_1, \dots, x_N\}$ is a codebook of size N , i.e. there is a partition (called Voronoi tessellation) $\{A_i\}_{1 \leq i \leq N}$ with $A_i = \{x \in \mathbb{R}^d : T(x) = x_i\}$, so that

$$T(x) = \sum_{i=1}^N x_i \mathbf{1}_{A_i}(x).$$

For a given $N \in \mathbb{N}$, an N -tuple of elementary quantizers (x_1, \dots, x_N) is optimal if it minimizes over $(\mathbb{R}^d)^N$ the quantization error:

$$\|X - \hat{X}\|_r = \min_{(y_1, \dots, y_N) \in (\mathbb{R}^d)^N} \mathbb{E} \left[\min_{1 \leq i \leq N} \|X - y_i\|^r \right]^{1/r} \quad (1)$$

induced by replacing X by \hat{X} . Then, instead of transmitting the complete signal $X(\omega)$ itself, one first selects the closest x_i in the quantizer set and transmits its (binary or Gray coded) label i . After reception, a proxy $\hat{X}(\omega)$ of $X(\omega)$ is reconstructed using the code book correspondence $i \rightarrow x_i$. Typically r is fixed to be equal to 2, leading to a quadratic quantization error. In this case, an N -optimal quantizer for a distribution P^X is a N -quantizer that minimizes the mean squared error (also called the *distortion* function):

$$\inf_T \mathbb{E}[(X - T(X))^2].$$

In d dimensions, the minimal quantization error converges to zero at a rate of $N^{-\frac{1}{d}}$ as $N \rightarrow \infty$, according to the so-called Zador theorem. Several stochastic optimization procedures based on simulation have been developed to compute these optimal quantizers. For a comprehensive exposition of mathematical aspects of quantization, we refer to [Graf and Luschgy, 2000] and [Luschgy and Pagès, 2023]. Remarkably, it can be shown (e.g. [Graf and Luschgy, 2000]) that an optimal quantizer is a Monge map minimizing the Wasserstein metric $W_2(P^X, P^Y)$ between P^X and P^Y , where P^Y is a discrete measure with N points. In the case where P^Y has two values, with respect to the simplest quantization approach, we impose additional constraints on the codebook points, with one point mass at zero and another at a positive magnitude m , to indicate the presence of loss. The quantization problem is then defined as the constrained two-points quantizer with centers $\{x_1, x_2\} := \{0, m\}$, i.e. as a mapping $T : \mathbb{R}^+ \rightarrow \{0, m\}$ with

$$T(x) = \begin{cases} m & x \geq a \\ 0 & x < a, \end{cases}$$

where a is a threshold to determine. Then, the optimal quantization problem for $X \sim P^X$ with constrained knot at zero writes

$$\inf_{a, m \in \mathbb{R}^+} \mathbb{E}[(X - T(X))^2],$$

and, as the boundary of the Voronoi cells separating the two quantizers x_1, x_2 is given by the center of the interval $[0, m]$, it turns out that the Voronoi partition is given by the two regions $A_1 = \{x \in \mathbb{R}_+ : 0 \leq x \leq m/2\}$ and $A_2 = \{x \in \mathbb{R}_+ : x_2 \geq m/2\}$. Thus, in the case where $r = 2, N = 2, x_1 = 0, x_2 = m$, (1) reads

$$\|X - \hat{X}\|_2 = \min_{m \in \mathbb{R}_+} \mathbb{E} [\min\{X^2, (X - m)^2\}]^{1/2} \quad (2)$$

For the two-point distribution for P^Y , [Faugeras and Pagès, 2024] proved the following result.

Theorem 3.1. (i) If $\mathbb{E}(X^2) < \infty$ and the support of P^X contains at least two points, then there exists a magnitude-propensity pair (p_X, m_X) minimizing the distortion.

(ii) An optimal pair (p_X, m_X) is characterized by solving for a the equation

$$2a = \mathbb{E}[X|X > a], a > 0$$

and then setting $m_X = 2a, p_X = P^X(X > a)$.

Note that m_X can be interpreted either as (twice) a VaR_α (resp. as an ES_α) for a special value of the confidence level α :

$$m_X = CVaR_{p_X}(X) = 2VaR_{1-p_X}(X).$$

What is more, the confidence level α is no longer a subjective choice dependent on the user. This is an improvement from a technical perspective, possibly a drawback in the application playground, as the prudential level is usually regulation-driven, i.e. assigned by the authorities: typically, $\alpha = 99\%$ in the market risk field, $\alpha = 99.9\%$ in the credit risk, finally $\alpha = 99.5\%$ in the *Solvency* insurance regulation.

3.3 Extension to the case of 3-Points Optimal Constrained Quantization

In this subsection we are going to prove the main theoretical contribution of the paper, namely the extension of the results of [Faugeras and Pagès, 2024] to the case of three points. The 3-Points distribution introduces a sophisticated and comprehensive approach to risk quantification, transcending the limitations of conventional methods. This distribution represents risk through three distinct points. The first point is the no loss case, $m_0 = 0$. The second point aims to designate moderate risk (m_1, p_1) , where m_1 represents the magnitude of potential moderate losses (corresponding to the most likely scenarios), and p_1 represents the probability of occurrence for these losses. The third point must capture extreme losses (m_2, p_2) , with m_2 denoting their magnitude and p_2 indicating the probability of occurrence for these severe events. As an example, the so-called 3-point risk analysis consists of summarizing the risk with 3 relevant values, namely: the best case, the worst case, and the most likely case. The three values are usually stimated (assessed) by submitting a standardized questionnaire to a panel of experts. For an extended review of these qualitative risk assessment techniques, see e.g., [Aven, 2016]. The quantitative approach of our work differs from this kind of methodology for its more rigorous and objective framework. The 3-Points distribution with a specific constraint introduces an additional criterion that influences the optimal selection of the discrete distribution parameters. This modification aims to avoid that the pure mathematical optimal solution does not highlight high magnitude losses. Furthermore, the stability of a risk measure over time is a general requirement of any risk measure in the banking regulation.

Following [Graf and Luschgy, 2000], a 3-vector quantizer on $\mathbb{R}, ||\cdot||$) is a mapping $T : \mathbb{R} \rightarrow \{x_0, x_1, x_2\}$, where we relabeled the codebook $(\{x_1, x_2, x_3\} \rightarrow \{x_0, x_1, x_2\})$ for notational convenience. More specifically, one can refine a measure of risk into a classification between a “moderate” and a “large” loss, by using a three points discrete measure, $P^Y = (1 - p_1 - p_2)\delta_0 + p_1\delta_{m_1} + p_2\delta_{m_2}$, with $0 < m_1 < m_2$ and where $x_0 = m_0 = 0$. Utilizing a three-point discrete measure enables the encoding and quantification of both moderate and large losses on the magnitude and propensity scales. This is achieved through the pairs $(m_1; p_1)$ and $(m_2; p_2)$, respectively. One thus introduces the constrained three-point quantizers with centers $\{0, x_1, x_2\} := \{0, m_1, m_2\}$, i.e. as a mapping $T : [0, \infty) \rightarrow \{0, m_1, m_2\}$ with $T(x) = 0\mathbf{1}_{x \leq a_1} + m_1\mathbf{1}_{a_1 < x \leq a_2} + m_2\mathbf{1}_{x > a_2}$, where $0 \leq a_1 < a_2$ are thresholds to be determined. Then, the optimal quantization problem with constrained knot at zero for a random variable X with $\mathbb{E}[X^2] < \infty$ (and whose support contains at least three points) writes

$$\inf_{(a_1, m_1, a_2, m_2) \in [0, +\infty)^4} \mathbb{E}[(X - T(X))^2]. \quad (3)$$

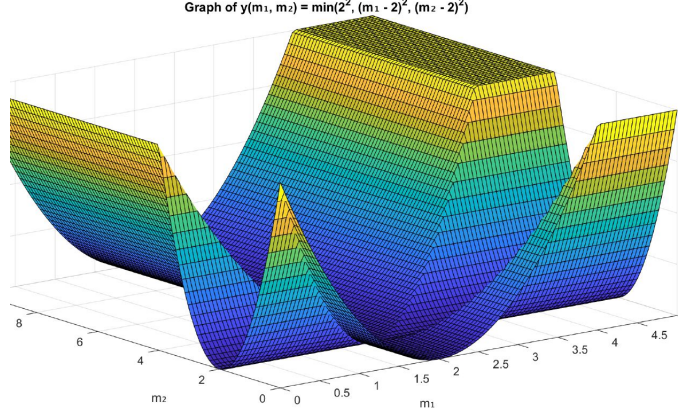


Figure 1: Graphic of the function $H(m_1, m_2) = 2^2 \wedge (m_1 - 2)^2 \wedge (m_2 - 2)^2$.

Using standard arguments, see e.g. [Graf and Luschgy, 2000] and [Luschgy and Pagès, 2023], one already knows that $a_1 = m_1/2$ and $a_2 = (m_1 + m_2)/2$, i.e. that the Voronoi regions write

$$A_0 = \{x : 0 \leq x \leq m_1/2\}, \quad (4)$$

$$A_1 = \{x : m_1/2 \leq x \leq (m_1 + m_2)/2\}, \quad (5)$$

$$A_2 = \{x : x \geq (m_1 + m_2)/2\}. \quad (6)$$

As a consequence, the distortion/objective function writes as a sole function of the magnitudes m_1, m_2 as

$$D(m_1, m_2) = \mathbb{E}[X^2 \wedge (X - m_1)^2 \wedge (X - m_2)^2] = \mathbb{E}[X^2 \wedge (X - m_2)^2] \mathbf{1}_{m_2 \geq m_1}, \quad (7)$$

since $m_2 > m_1$. Introduce now for $x = (x_1, x_2) \in \mathbb{R}_+^2$

$$H(x_1, x_2) = h_0(x_1, x_2) \wedge h_1(x_1, x_2) \wedge h_2(x_1, x_2) := X^2 \wedge (X - x_1)^2 \wedge (X - x_2)^2.$$

Note that h_0, h_1, h_2 are differentiable functions, so that the function H is (at least) piecewise differentiable, and it has directional derivatives everywhere, as shown in Figure 1 where we display the function H for $X = 2$.

As shown in [Faugeras and Pagès, 2024], a convenient tool in the setting of optimization of piecewise smooth functions is the concept of Bouligand derivative (B-derivative), see e.g. [Scholtes, 2012]. By dropping the requirement of linearity of the differential, it represents a first-order approximation and allows to have a single-valued notion of differential.

Definition 3.1. A function $f : \mathbb{R}^2 \rightarrow \mathbb{R}$ is B-differentiable at $\tilde{x} \in \mathbb{R}^2$ if there exists a positive homogeneous function $\nabla^B f(\tilde{x}) : \mathbb{R}^2 \rightarrow \mathbb{R}$ s.t.

$$f(\tilde{x} + v) = f(\tilde{x}) + \nabla^B f(\tilde{x})(v) + o(\|v\|), \quad \forall v \in \mathbb{R}^2. \quad (8)$$

We have the following lemma on the B-differentiability of the function H .

Lemma 3.2. Let $H(x) = \min(h_0(x); h_1(x); h_2(x))$, where $h_0, h_1, h_2 : \mathbb{R}^2 \rightarrow \mathbb{R}$ are differentiable functions. Then H is B-differentiable, and its B-differential is given as follows.

1. For i, j, k distinct indexes in $\{0, 1, 2\}$, for x such that $h_i(x) < \min(h_j(x); h_k(x))$

$$\nabla^B H(x)(v) = \nabla h_i(x)v = \frac{\partial h_i}{\partial x_1}(x)v_1 + \frac{\partial h_i}{\partial x_2}(x)v_2.$$

2. For x such that $h_i(x) = h_j(x) < h_k(x)$

$$\nabla^B H(x)(v) = \min(\nabla h_i(x)v; \nabla h_j(x)v).$$

3. For x such that $h_0(x) = h_1(x) = h_2(x)$

$$\nabla^B H(x)(v) = \min(\nabla h_0(x)v; \nabla h_1(x)v; \nabla h_2(x)v).$$

Proof. See Appendix A.

Before proving the B-differentiability of the distortion function D , we recall that a point x is critical (that is, a potential minimizing point) for D if $\nabla^B D(x)(v) \geq 0 \forall v \in \mathbb{R}^2$.

Theorem 3.3. *The distortion function D defined in (7) for a random variable $X \in L^2$ is B-differentiable on $x \in \mathbb{R}_+^2$, with B-derivative given by*

$$\begin{aligned} \nabla^B D(x_1, x_2)(v) &= \mathbb{E}[2(x_2 - X)v_2 \mathbf{1}_{x_1+x_2 < 2X}] \\ &\quad + 2\mathbb{E}[\min((x_1 - X)v_1 \mathbf{1}_{x_1 < 2X} \mathbf{1}_{x_1+x_2 > 2X})] \\ &\quad + 2\mathbb{E}[\min(0; 2Xv_1) \mathbf{1}_{x_1=2X}]. \end{aligned} \quad (9)$$

Moreover, the points $(0, 0)$, $(0, 2X)$, $(2X, 0)$, $(2X, 2X)$, $(0, x_2)$ (with $x_2 > 2X$) and $(x_1, 0)$ (with $x_1 > 2X$) are not critical for the distortion function, since $\nabla^B D(x)(v) < 0$ for some $v \in \mathbb{R}^2$.

Proof. See Appendix A.

We have now all the ingredient to characterize the critical points, namely we can find the quantizers of the 3-points distribution by solving a system of equations. From Equation (9), by taking separately $v_1 = 0$, $v_2 = 0$ we get

$$P\left(X = \frac{x_1 + x_2}{2}\right) = P\left(X = \frac{x_1}{2}\right) = 0,$$

and we find the following system

$$\begin{aligned} \mathbb{E}[(x_1 - X) \mathbf{1}_{\frac{x_1}{2} < X < \frac{x_1+x_2}{2}}] &= 0, \\ \mathbb{E}[(x_2 - X) \mathbf{1}_{X > \frac{x_1+x_2}{2}}] &= 0, \end{aligned}$$

which leads to the optimal quantizers:

$$m_1 = \mathbb{E}\left[X \mid \frac{m_1}{2} < X < \frac{m_1 + m_2}{2}\right], \quad (10)$$

$$m_2 = \mathbb{E}\left[X \mid X > \frac{m_1 + m_2}{2}\right]. \quad (11)$$

Eventually, once the system is solved using a fixed-point technique, the probabilities p_1, p_2 associated with the quantizers are given by integrating the probability of X on the Voronoi regions given by (5)–(6). Finally, the probability mass of the initial (constrained) quantizer $m_0 = 0$ is given by $(1 - p_1 - p_2)$.

Remark 3.1. The points m_1 and m_2 optimally represent the underlying loss distribution, independently of any *a priori* thresholds such as those described by *VaR* or *ES* associated with some confidence level. Consequently, there is no guaranty that m_2 , intended to represent an extreme loss, can be directly compared to *VaR* or *ES*, nor that the associated probability p_2 has any direct relationship to the confidence level α . Therefore, it is not surprising that, as we show in our numerical experiments, the value of m_2 can differ from *VaR*. In other words, comparing m_2 with *VaR* is not a meaningful exercise, as we only have one m_2 while

we can compute many VaR according to the relevant confidence levels adopted by the financial regulation used in practice (0.95, 0.99, 0.995, 0.999, etc.). Nevertheless, it is advisable to introduce a constraint to prevent the purely mathematical optimum from neglecting high-magnitude losses and to ensure the temporal stability of the risk measure, which is a requirement in banking regulation. In the context of market risk, where m_2 typically turns out to be slightly lower than VaR , it seems reasonable to impose an additional restriction on m_2 , ensuring that it reaches a value at least equal to VaR . This preserves its interpretation as an indicator of extreme risk while maintaining compliance with regulatory standards, which are expressed in terms of VaR . From an analytical standpoint, the introduction of such a constraint increases the complexity of the optimization procedure. However, from a numerical perspective, the procedure remains highly efficient (see the appendix for further details on the optimization procedures), and the resulting performance is fully satisfactory.

4 Case Study based on Real Datasets

4.1 Practical Risk Management

In this section we try to exploit the features of the magnitude-propensity approach in relevant different fields. To achieve significant insights, some concepts about practical risk management are needed. First, the *risk category* must be well defined. In a high level classification, we distinguish market risk vs. credit risk. Market risk is given by the PnLs (profits and losses) uncertainty due to the price dynamics of the assets in the portfolio. Financial portfolios of large banks typically are not concentrated; even if we may observe extreme events due to volatility peaks, the PnLs typical distribution is quite smooth. Credit risk is related to the losses implied by any default in the portfolio, where once the default occurs, the recovery rate drives the amount that one can get back. In this case, we have also concentrated portfolio, and the binary nature of the outcome (no default, default) implies a loss distribution that sometimes shows some peaks in the extreme losses region. Due to high quality, high frequency data and finally a long history of the risk management in the financial area, the regulation typically asks for more conservative confidence level in the credit risk (e.g. 0.999), as we have low frequency data, low data coverage (e.g. many issuers in the portfolio might be unrated, with missing default probability), and the default correlations are not strict sense observable. The above issues are summarized in the model risk definition, i.e. the risk arising from misspecifications in the model itself or in its parameters. The boundary between market and credit risk is quite flexible. A relevant example is given by the *migration risk*, i.e. the risk of losses due to downgrade (rating worsening) of one or more issuers in the portfolio. Furthermore, the well known *spread risk*, i.e. the risk of losses due to the increase of the spread level (e.g. in the government bonds in portfolios) is clearly assigned to the market risk discipline in the banking context, see [Basel Committee of Banking Supervision, 2019], while it can be allocated to market or credit risk filed according to the Solvency regulation for the insurance companies, see [European Commission, 2015]. See [Embrechts et al., 2015a] for a comprehensive survey of the main modeling methodologies related to market and credit risk. Taking into account the above practical and regulatory concepts, we decided to exploit the capabilities of the magnitude-propensity 3-points proposal in three different application contexts, trying to cover a relevant area of the broad market and credit risk fields, as summarized in Table 1.

4.2 Applications Overview and Preliminary Concepts

Financial Institutions commonly adopt two main approaches to measure the risk of a financial position, which they select based on their specific purposes and regulatory requirements: *sensitivity* measures and risk measures based on the profit and loss (PnLs) distribution. Both these approaches are prescribed by the regulations and adopted by the risk management departments.

ID Case	Description	Risk Category	Regulation	PnLs Model	Confidence level	Horizon
1	Market Risk Trading book	Market Risk	Basel	Historical non parametric	99%	1D
2	Credit Risk Trading Book	Credit Risk	Basel	Monte Carlo parametric	99.9%	1Y
3	Market Risk Insurance	Market Risk	Solvency	Monte Carlo parametric	99.5%	1Y

Table 1: Overview of VaR calculation according to regulatory risk cases: model type, regulation, model calculation, confidence level and horizon.

To have a better understanding, let us denote the value of a generic portfolio at the evaluation time t_0 as V_{t_0} . Following standard risk-management practice, V_{t_0} is modeled as a function of time t (the index t_0 represents the current time, so that e.g. $t_0 - 1$ indicates the previous time and so on) and a d -dimensional random vector $\mathbf{Z}_t = (Z_t^1, Z_t^2, \dots, Z_t^d)$ of risk factors, i.e., underlying variables or drivers like e.g. interest rates, stock prices, implied volatility, or exchange rates. Hence,

$$V_{t_0} = f(t_0, \mathbf{Z}), \quad (12)$$

with $f : \mathbb{R} \times \mathbb{R}^d \rightarrow \mathbb{R}$ and where \mathbf{Z} informally represents the heterogeneous set of risk factors (scalar, array, surface, cubes etc.) recorded at past observation times, depending on specific payoffs of the instruments of the portfolio.

For the sake of simplicity, we omitted in the expression the various parameters on which depend the stochastic processes that describe the risk factors \mathbf{Z}_t dynamics. *Sensitivity* based risk measures rely on each risk factor impact, which can be defined as the change in value of an instrument (position) given a small, predetermined (hypothetical) movement in a risk factor that affects the instrument's value: from a mathematical perspective, sensitivities are computed as partial derivatives of the function f with respect to \mathbf{Z} . By defining $\Delta\mathbf{Z} = \mathbf{Z}_t - \mathbf{Z}_{t-1}$ the vector of observed shocks applied to the risk factors, this approach enables the estimation, assuming linearity between V and \mathbf{Z} , of the portfolio response to the new market conditions by multiplying $\Delta\mathbf{Z}$ with the corresponding factor sensitivities, namely $\Delta V \cong \nabla f \cdot \Delta\mathbf{Z}$. Technically speaking, it is just a differential of the value function f evaluated for some small increment $\Delta\mathbf{Z}$. Of course, the second order impact could be added to improve the accuracy of the approximation. The most popular second order approximations are referred according to some naming conventions, such as the *gamma* effect, i.e. the second order derivative of an option the the underlying price, and the *convexity*, namely the second order impact of the interest rate level change on a bond or interest derivative (e.g swaps, caps and floor options). Such risk measures, though valuable in providing information about the robustness of a portfolio value to specific events, have limitations when making capital-adequacy decisions: they do not deal with the dependency properties of the risk drivers, hence they can not create a picture of the overall riskiness of the portfolio of a financial institution, and they do not have any information about the likelihood of the approximated *PnLs*, see [McNeil et al., 2005]. Furthermore, sensitivity approach is a typical *what-if* approach, where there is some subjectivity in defining the level of the extreme market shock to be applied. For this reason, the so called *probabilistic* measures, such as *VaR* and *ES*, are prominent in the financial regulation, to ensure that banks and the insurance companies have enough own capital to face the potential losses under "extreme yet plausible" scenarios, a statement adopted very often by the financial authorities to define the scope of the stress test exercise. The following subsections give more details on the context of the three applications of the methodology.

4.2.1 Market Risk with Historical Simulation

In the introduction, we recalled that the historical simulation is the most popular approach in the large banks to evaluate the financial risk, by *VaR*, *ES* or any other measure. Let us give a brief formal explanation, mainly of

the historical simulation through the *full evaluation* methodology. We can define the daily portfolio *PnLs* as the change in value of the portfolio, driven by the series of risk factor changes \mathbf{Y}_t where $\mathbf{Y}_t := \mathbf{Z}_t - \mathbf{Z}_{t-1} = \Delta \mathbf{Z}$ or alternatively $\mathbf{Y}_t := \ln(\mathbf{Z}_t / \mathbf{Z}_{t-1})$. The *additive* (or absolute) vs *multiplicative* (or percent) definition of a shock is calculated depending on the asset class of the risk factor: usually, banks adopt the additive convention for interest rates and spreads, while the multiplicative convention for returns on equities, foreign exchange rates and funds. For a useful survey, see e.g. [Hudson and Gregoriou, 2010]. For each $t = t_0 - 1, \dots, t_0 - S$ in our array of historical scenarios, the portfolio value is calculated by summing up the contributions over all proper pricing functions f_m (related to the instrument $m = 1, \dots, M$) that select just the required risk factors, typically a small subset of the whole market data \mathbf{Z} .

If we adopt the *full evaluation* approach, the portfolio's *PnL* at a given calculation time (t_0) and for any observed scenario $[t, t_0]$ with $t = t_0 - 1, \dots, t_0 - S$, is given by:

$$PnL_{[t, t_0]} = \sum_m f_m(t_0, \mathbf{Z} + \mathbf{Y}_t) - f_m(t_0, \mathbf{Z}). \quad (13)$$

The value (or *PnLs*) of the portfolio is then the sum of a huge variety of pricing functions with heterogenous risk factor inputs.

Remark 4.1. The above stylized expression uses the additive convention for the shock. In the multiplicative case the part $f_m(t_0, \mathbf{Z} + \mathbf{Y}_t)$ is replaced by $f_m(t_0, \mathbf{Z} \cdot \exp(\mathbf{Y}_t))$.

Remark 4.2. In the above expression we refer to absolute (eur, dollar, etc.) *PnLs*, i.e. the standard convention for bank portfolios. In the asset management field it is common to switch to relative (percent) *PnLs*, by a scaling factor, i.e. the current value V_{t_0} of the portfolio.

Remark 4.3. If *PnL* is calculated using the sensitivities approach, the function $f_m(\cdot)$ is replaced by the gradient vector, and we have $PnL \cong \nabla f_m \cdot \Delta \mathbf{Z}$, possibly enriched with the second order term based on the Hessian matrix.

Remark 4.4. One could replace the term $f_m(t_0, \mathbf{Z} + \mathbf{Y}_t)$ by the more general expression $f_m(t_0 + 1, \mathbf{Z} + \mathbf{Y}_t)$ to properly take into account the “ageing effect”, meaning that if the market shocks are supposed to prevail in a given time horizon h ($h = 1$ for simplicity in this case), then the hypothetical *PnLs* must take into account such time shift in order to shorten the time to maturity of the instruments, such as bonds or derivatives. Despite this seems an obvious concept, very often the ageing is not actually considered, as it would require that all the pricing libraries of the bank are able to price in the future, i.e. not only taking into account the time decay (such as the decreasing maturity for a bond) but also all the other possible events, such as dividend payments and so on. In most cases, a risk factor shock is applied, but not the time shift, so applying the *instantaneous shock* assumption. For the insurance sector, see the survey in [EIOPA, 2024], where it is outlined that only 6 out of 20 insurance companies manage the ageing effect, while the remaining ones apply the instantaneous shock model.

In other words, *PnLs* distribution consists in the range of possible values X (potential profits and losses) that a portfolio may experience over a specific time horizon. The historical simulation approach is data driven, as no parametric distribution family is assumed for the returns. Furthermore, we do not need any correlation parameters or dependence modeling as it is assumed that the dependencies among the risk factors are divided into small pieces inside each scenario (t). Recalling that in measuring portfolio's risk the fundamental aspect is estimating the cumulative density function of the *PnLs*, namely $F(x) = \mathbb{P}(X \leq x)$, or functionals of it, see [McNeil et al., 2005] and that Value at Risk is essentially a quantile (typically 95%, 99%, 99.9%) of the above mentioned *PnLs* distribution, the final step consists in estimating a suited percentile to be applied to the vector of *PnLs*. For the number S of scenarios, banks usually adopt $S = 250$ or $S = 500$, i.e. they collect about 1 or 2 years of daily changes in all the risk factors to which the portfolio is sensitive. The regulatory time horizon h differs from the daily popular horizon, e.g. $h = 10$ days in the current Basel regulation. Collecting a sample of historical non-overlapping risk factors returns ($\mathbf{Y}_{t,h} = \mathbf{Z}_{t+h} - \mathbf{Z}_t$) could be quite difficult. As an

example, to achieve $S = 250$ non overlapping scenarios with $h = 10$ days, we need $h \cdot S = 2500$ observations, i.e. around 10 years of full time series for all the risk factors. Hence, the authorities generally accept the *square root rule*, namely $VaR(\alpha, h) \equiv VaR(\alpha, 1) \cdot \sqrt{h}$.

In daily practice, VaR can be estimated in many ways: from the basic *empirical* percentile to some more robust estimators, such as L -estimators and Harrel-Davis estimator, that smooth the estimation by averaging the $PnLs$ in a neighborhood of the empirical quantile, see[Harrel, 1982]. In this field a comprehensive reference is given by [David and Nagaraja, 2004]. These more advanced estimators of course may be used also in the other simulation contexts, such the Monte Carlo simulation. The historical simulation approach is very popular in the banking industry due to some key reasons. First, it is very intuitive, as the empirical past (recent) distribution of the risk factors returns is accepted (assumed) to be the best estimation of the unknown exact distribution. No functional subjective assumption is made about the distribution shape. Being purely data-driven, it does not rely on any parameters (volatility, correlations, etc). In real world, portfolios may contain many thousands of risk factors: estimating the related parameters and updating periodically the estimates is a very challenging task. For the sake of simplicity, in the following applications we will refer to the basic empirical quantile (e.g. the 5th worst result in an array of $S = 500$ PnL).

4.2.2 Credit Risk in the Trading Book: the Default Risk Charge by Montecarlo approach

The Default Risk Charge (DRC) is a regulatory measure designed to capture default risk within the trading portfolio, as required by Basel standards, particularly within the framework of the Fundamental Review of the Trading Book (FRTB) outlined in the [Basel Committee of Banking Supervision, 2019] document. This model is designed to quantify the risk of loss resulting from the failure of a counterparty or issuer of financial instruments, including equity, bond, and derivative exposures. The Default DRC is specified in Chapter 7, and in the updated Basel Framework, in Paragraphs MAR 33.18- 33.38. These documents establish the criteria for calculating default risk, specifying that

- It must be calculated over a one-year horizon.
- It must reflect a 99.9% confidence level.
- It must include all exposures sensitive to default risk within the trading portfolio, excluding those specifically defined as *non-material* risks.

In the practice, the default of each issuer is modeled by a Merton type model. The *credit worthiness* dynamics of the obligor n ($n = 1, \dots, N$) is defined by the model

$$\Delta X_n = \sum_{k=1}^K \beta_{k,n} \cdot W_k + \sigma \cdot \varepsilon_n \quad (14)$$

where the coefficients $\beta_{k,n}$ are the *factor loadings* and describe the systematic risk, driven by a set of risk factors $k = 1, \dots, K$, the term ε_n is the specific (obligor) factor, with W_k correlated standardized Normal random variables, with $E[W_k \varepsilon_n] = 0 \forall n, k$.

In this standardized framework, the default happens if the credit worthless is below a given threshold, $D_n = \{X_n \leq \Phi_n^{-1}(PD_n)\}$, where Φ_n and PD_n indicate the cumulative distribution of ΔX_n and the default probability of the n -th issuer in the portfolio. The factor loading coefficients are estimated by a statistical regression step, by combining the time series of the list of risk factors with the issuer equity prices or spreads. In large banks, we generally have hundreds of issuers (N) and some dozens of risk factors (K). The distribution of the portfolio loss is given by 2, that we recall here:

$$Loss_{DRC} = \sum_{n=1}^N EAD_n \cdot \mathbf{1}_{D_n} \cdot LGD_n.$$

This quantity may not be available in closed form, thus most banks develop Montecarlo simulation algorithms. Indeed, also the Basel regulation explicitly refers to the simulation approach. Due to the extremely high confidence level $\alpha = 99.9\%$, to achieve a sufficient accuracy the number of simulation scenarios must be very high, usually in the range $[100k, 1M]$. At the end of the simulation cycle, the quantile is estimated (empirically or using some smoother estimator, in line with the usual approach adopted for computing the VaR by historical simulation).

With respect to the work by [Faugeras and Pagès, 2024], where the authors deal with the losses coming from the claims of the clients (i.e., the domain of the distribution lies in one-side in the real axis), we recall that for the VaR of the financial portfolio, a bank can experience both profits and losses, while for the DRC we just have losses, so making the context identical to that of the insurance sector.

4.2.3 Market Risk in the Insurance field: Monte Carlo approach

While in the banking sector the liabilities side of the balance sheet typically mirrors the asset side—comprising the same types of instruments such as bonds, equities, derivatives, and loans with opposite signs—in the insurance business the situation is structurally different. The asset side, usually consisting of bonds and investment funds, is primarily designed to hedge the liabilities, namely the claims arising from policyholders' underwriting activity.

Since the products sold to clients span very heterogeneous categories (life, longevity, accident, financial, and others) and often include numerous complex clauses linked to a wide range of contingent events, it becomes extremely challenging to construct a fully granular model incorporating all the specific inputs. For this reason, most insurance companies adopting a probabilistic approach to risk modeling rely instead on a compact set of risk factors, which are expected to adequately capture the distribution of profits and losses across the entire balance sheet.

To do that, the insurances use an *economic scenario generator* (ESG) that adopts the Monte Carlo simulation to project thousands of paths of the selected risk factors. For each path, all the positions of the insurance are evaluated to obtain the profile of profits and losses. Finally, the desired risk measures (percentiles, expected shortfall, etc) are estimated by analyzing the simulated results. For a recent extended review see [EIOPA, 2025]. Briefly, the Monte Carlo approach is not a choice in the insurance sector, but it represents a mandatory tool, due to the high number of risk factors and to the complexity of the products managed in insurance. Of course, also in the practical applications several improvements have been implemented to reduce the weaknesses of the basic Monte Carlo approach, such as the simulation error. The improvements are related to both the simulation step (by low discrepancy algorithm, quasi Monte Carlo methods, etc) and the estimation step, by smoothing the straight empirical estimator with more robust estimators, e.g, the Harrel-Davis estimator. In our application, we collected samples of 100K or 200K simulations related to the market risk of a major European insurance company and compared the 3-points results, mainly m_1 and m_2 , with respect to the 0.995 1 year Solvency VaR .

4.3 The Dataset

The data on which we apply the approach presented in this paper includes 1 year of daily $PnLs$ time series of a representative portfolio of a large European bank, encompassing both its banking and trading books. As explained in the previous section, the $PnLs$ have been obtained by full revaluation of the positions composing the portfolio with historical simulations of the risk factors. The time window includes 254 business dates (July 3rd 2023- July 1st 2024).

The portfolio consists of more that 100k positions, spanning from bonds to equities to derivatives. A couple of business dates where some outliers were detected (due to failures in the bank software systems) have been removed. It is worth to note that the process to get the portfolio VaR is typically a bottom-up

process. In other words, Formula 13 is just a synthetic definition, as it is practically calculated as

$$PnL_{[t, t_0+h]} = \sum_{m=1}^M f_m(t_0 + h, \mathbf{Z} + \mathbf{Y}_{t+h}, \mathbf{P}_n), \quad (15)$$

being n the index of each position in the portfolio and \mathbf{P}_n the information (maturity, coupon, currency, etc) related to any position. The VaR in historical simulation (Case 1) is then usually reported at a portfolio level, but it can be easily broken down for any purpose at any more granular level, according to the different analysis that are required.

As concerns the DRC (Case 2), we have for the same bank a set of some end-of-month calculation, referred to March 2024, June 2024, December 2024 and to February 2025. The outcomes are based on 200k simulations. The portfolio consists of many thousands positions, belonging to $O(10^3)$ issuers that could default. Finally, for the insurance portfolio (Case 3), we analyzed an array of 100k PnL generated by a multi dimensional Gaussian copula based on about 150 risk factors.

To test the methodology and the numerical procedures in a more comprehensive context, we exploit the methods for the three fields defined in the previous section.

- Case 1: Market Risk in the Banking sector. In this case the methodology for the PnL distribution is based on the historical simulation with 250 scenarios. Regulatory VaR (Basel regulation): 1 day, 99% confidence level. Risk sources: equity, interest rates, forex, spread, commodities.
- Case 2: Credit Risk in the Banking sector. Then we refer to binary events driven by PD (Default Probability), that determine the losses by the LGD (loss given default) parameter applied to the defaulted position. LGD could be stochastic, e.g. a Beta random variable in the $[0, 100\%]$ range. Methodology for the PnL distribution: Montecarlo simulation, multivariate gaussian copula, Merton-type default model. Regulatory VaR (Basel regulation): 1 year, 99.9% confidence level. Risk source: default.
- Case 3: Market Risk in the Insurance sector. Here the methodology for the PnL distribution is based on Montecarlo simulation, multivariate gaussian copula, and Merton-type default model. Regulatory VaR (Solvency regulation): 1 year, 99.5% confidence level. Risk sources: equity, interest rates, forex, spread, credit migration.

4.4 Results

In this subsection we will show that overall, the results demonstrate that our methodology based on the magnitude propensity remains theoretically sound, numerically stable, and flexible enough to account for diverse risk profiles, while preserving interpretability across market, credit, and insurance domains.

4.4.1 Case 1: Market Risk in the Banking Sector

In the initial stage of analysis, we consider the analytical solution obtained by quantization using the fixed-point method, allowing the quantities m_1 and m_2 to vary freely, i.e. subject only to the natural constraint $0 < m_1 < m_2 < \text{worst case}$. Figure 2 illustrates the daily evolution of m_1 and m_2 , and compares them with the conventional risk measures— VaR at the 99% confidence level and ES at the 97.5% level, the latter chosen to ensure comparability with the 99% VaR under a Gaussian framework. The unconstrained three-point distribution provides a clear and interpretable structure for quantifying risk, yielding results of a conservative nature that capture both moderate and extreme losses while avoiding excessive overestimation.

The interpretation of m_1 as a measure of moderate risk appears straightforward; however, the parameter m_2 often exhibits overly conservative behavior, frequently taking values below the corresponding VaR. This outcome is expected, as the VaR constraint becomes more binding in market environments characterized by moderate tail risk and smoother loss distributions. Consequently, the resulting estimates tend to be slightly conservative, consistent with the regulatory framework that defines VaR at the 99% confidence level over a one-day horizon. To maintain coherence with this framework—where VaR is associated with extreme risk levels—we subsequently impose a constraint requiring m_2 to exceed the VaR threshold.

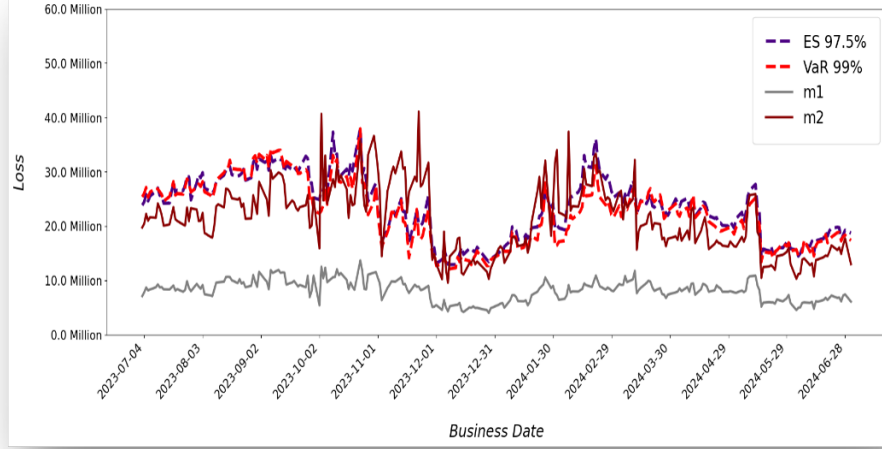


Figure 2: Time series plot of VaR 99%, ES 97.5%, m_1 and m_2 (without constraints on m_2).

If we add a constraint on the m_2 , the methodology provides the most satisfactory results, see Figure 3.

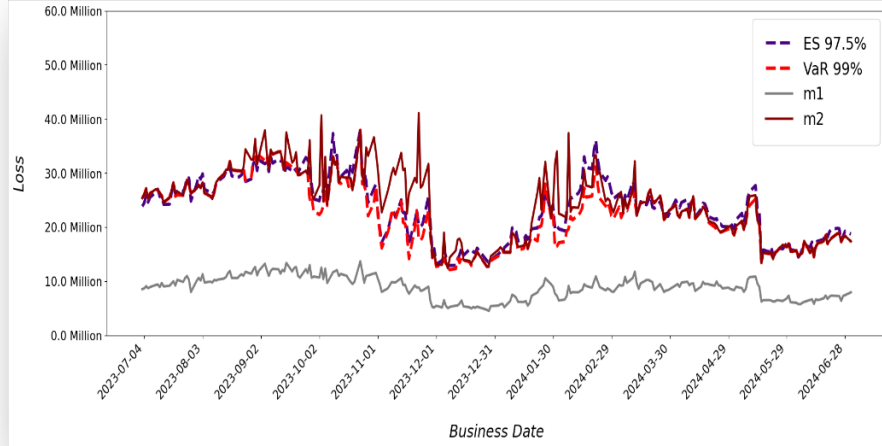


Figure 3: Time series plot of VaR 99%, ES 97.5%, m_1 and m_2 (with constraint $m_2 > VaR$).

The magnitude parameters m_1 and m_2 are intuitive, capturing both moderate and extreme losses, with stable dynamics over time relative to VaR and ES. Meanwhile, p_1 and p_2 remain smooth, with p_2 constrained

by $m_2 > VaR$ and p_1 free to vary within the $(0, 1)$ range.

As concerns the volatility of m_2 , this naturally follows from the variability of the VaR computed by the historical simulation method. Indeed, when the scenarios contributing to the empirical percentile fall outside the rolling window used for VaR, jumps may occur independently of the portfolio composition. This phenomenon, known as the *reshuffling effect*, could be mitigated or controlled through smoothing techniques, in analogy with those employed in filtered historical simulation, see e.g. [Barone-Adesi et al., 1999].

Furthermore, Table 2 shows that employing different optimization methods—including those based on Differential Evolution (briefly recalled in Appendix D) and on the Optimal Transport approach introduced earlier—yields remarkably consistent results. Computational times are extremely short (on the order of one second), thereby demonstrating the robustness and efficiency of the proposed methodology.

Parameter Metric	Without Constraint Fixed Point Eq. Semi-Analytical	Without Constraint Quantization Diff. Evol.	VaR 99% Constraint Quantization Diff. Evol.	Without Constraint Opt. Transport Sinkhorn-Knopp	VaR 99% Constraint Opt. Transport Sinkhorn-Knopp
m_1	8,206,458	8,210,913	8,878,795	8,465,072	8,834,289
m_2	21,452,567	21,453,539	24,397,039	21,971,367	23,925,852
p_0	67.96%	67.99%	69.34%	68.51%	69.25%
p_1	26.77%	26.74%	27.26%	26.94%	27.26%
p_2	5.27%	5.27%	3.40%	4.54%	3.48%
VaR 99%			22,531,887		
ES 97.5%			23,636,174		
Worst-case			34,616,367		

Table 2: Market risk — Fixed Point Eq. (Semi-analytical using (10),(11)), Quantization with Differential Evolution and Optimal Transport (Sinkhorn-Knopp) methods. Comparison with and without the VaR 99% constraint on m_2 .

4.4.2 Case 2: Credit Risk in the Banking Sector

Within the Credit Risk framework—modeled through a Merton-type setting with a multivariate Gaussian copula—the VaR constraint on m_2 plays a less active role, as we can immediately see from Table 3. Due to portfolio concentration effects, the presence of dominant exposures shifts the loss distribution to the right, leading to a naturally high value of m_2 , which already captures the fat-tail behavior of the underlying risk factors. Moreover, due to typical combinatorial effects associated with default events, significantly different scenarios may lead to very similar loss values, thereby making risk measures structurally unstable and leading to even more fragile risk decomposition techniques, as illustrated in the Appendix B. Consequently, imposing a constraint on the value of m_2 (for instance, requiring m_2 to exceed the VaR) may be of limited relevance, since the tail of the distribution is already heavily skewed toward extreme values, forcing m_2 into a region that the VaR is not able to reach, even with the typical credit risk regulatory requirement of a 99.9% confidence level. In other words, the extreme values of the distribution are so pronounced that the VaR no longer provides an adequate representation of tail risk, whereas m_2 effectively captures the true risk embedded in the distribution—yielding substantially higher values and thereby making any additional VaR-based constraint redundant.

A comparison between m_2 and the worst-case scenario, together with the graphical inspection of the losses histogram (see Figure 4), clearly points out how m_2 , acting as a barycenter of the extreme-loss scenarios, is positioned well above the VaR level, but remains consistently below the worst-case value. In fact, from Figure

Parameter Metric	Without Constraint Fixed Point Eq. Semi-Analytical	Without Constraint Quantization Diff. Evol.	VaR 99.9% Constraint Quantization Diff. Evol.	Without Constraint Opt. Transport Sinkhorn-Knopp	VaR 99.9% Constraint Opt. Transport Sinkhorn-Knopp
m_1	66,944,517	66,944,702	66,944,602	66,944,517	66,944,517
m_2	455,901,571	455,900,234	455,901,519	455,901,571	455,901,571
p_0	98.84%	98.84%	98.84%	98.84%	98.84%
p_1	1.13%	1.13%	1.13%	1.13%	1.13%
p_2	0.03%	0.03%	0.03%	0.03%	0.03%
VaR			125,882,878		
ES			256,621,107		
Worst-case			980,312,060		

Table 3: Credit risk comparison — Fixed Point Eq. (Semi-analytical using (10),(11)), Quantization with Differential Evolution and Optimal Transport (Sinkhorn-Knopp) methods. Comparison with and without the VaR 99.9% constraint on m_2 .

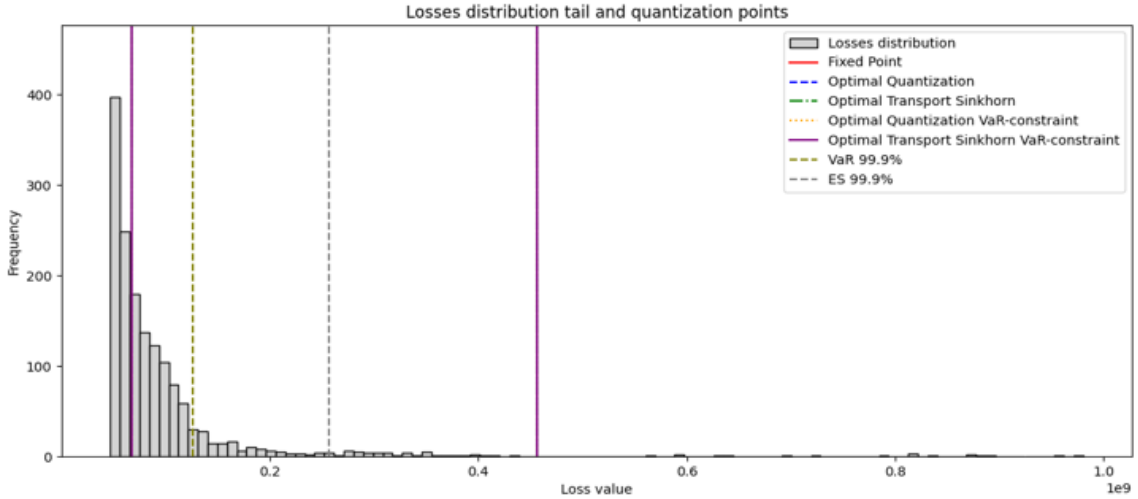


Figure 4: Credit Risk in the Banking Sector (VaR 99.9%, ES 99.9 %): tail of the Loss distribution.

4 we realize a sparse distribution on the right tail, leading to extreme losses with relatively non negligible probabilities. This is well captured by a relatively high value for m_2 with respect to the VaR and ES that underestimate the risk.

Due to the very high magnitude and low frequency of extreme losses, the associated parameter p_2 has a very small value. It is worth recalling, from an intuitive standpoint and in analogy with the theory of optimal transport, that the parameters (m, p) jointly minimize the “effort” required to transfer probability masses and values from a continuous distribution toward its discrete representation. This observation underscores the inherent tendency of VaR and ES to underestimate the probability of extreme losses, as is often the case in credit risk modeling.

In conclusion, the proposed methodology once again provides a more accurate and comprehensive representation of both moderate and extreme risks than traditional measures based on VaR and ES, particularly

within the complex framework of credit risk.

4.4.3 Case 3: Market Risk in the Insurance Sector

In the Insurance Market Risk case, the one-year horizon and 99.5% confidence level adopted under Solvency II standards confirm the adaptability of the proposed approach to different regulatory settings, see Table 4.

Parameter Metric	Without Constraint Fixed Point Eq. Semi-Analytical	Without Constraint Quantization Diff. Evol.	VaR 99.5% Constraint Quantization Diff. Evol.	Without Constraint Opt. Transport Sinkhorn-Knopp	VaR 99.5% Constraint Opt. Transport Sinkhorn-Knopp
m_1	292,726,533	292,251,249	292,250,135	292,726,533	292,726,533
m_2	1,326,307,230	1,326,305,007	1,326,304,534	1,326,307,230	1,326,307,230
p_0	90.03%	90.00%	90.00%	90.03%	90.03%
p_1	9.21%	9.24%	9.24%	9.21%	9.21%
p_2	0.76%	0.76%	0.76%	0.76%	0.76%
VaR			1,103,006,334		
ES			1,524,216,012		
Worst-case			2,791,623,043		

Table 4: Insurance market risk — Fixed Point Eq. (Semi-analytical using (10),(11)), Quantization with Differential Evolution and Optimal Transport (Sinkhorn-Knopp) methods. Comparison with and without the VaR 99.5% constraint on m_2 .

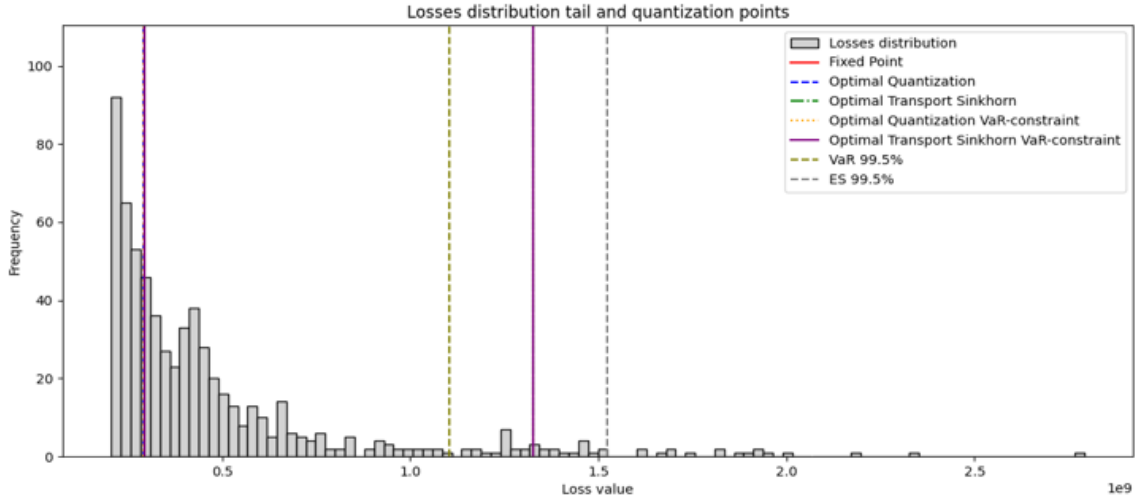


Figure 5: Insurance market risk (VaR 99.5%, ES 99.5%): tail of the Loss distribution.

In Figure 5 we show the tail of the loss distribution, from which we deduce a considerable amount of losses beyond the VaR, although the difference between the parameter m_2 and the VaR level is much less dramatic than in the previous case of Credit Risk. This behavior is consistent with the nature of the portfolio and the types of risks considered. Indeed, market risk within the insurance context also encompasses spread risk and migration risk (see Section 4.2.3). As a result, the PnL distribution may include potential jump

events, yet with tail effects that are less pronounced than those typically observed in pure credit risk settings. Consequently, m_2 assumes a meaningful position from an informative standpoint, identifying potentially significant losses, while p_2 —associated with extreme risk events—reflects a degree of conservatism (or prudence) broadly in line with the regulatory confidence level.

5 Conclusion

One of the long-standing challenges in risk management is the search for a risk measure that simultaneously satisfies the goals of objectivity, intuitiveness, and theoretical soundness. In this respect, the long and ongoing regulatory transition from Value-at-Risk (VaR) to Expected Shortfall (ES) clearly illustrates the difficulty of introducing new measures that can meet the expectations and requirements of all potential users.

In our work, we build on a frequency–severity framework, a well-established approach in the insurance domain that has recently received a rigorous theoretical foundation from [Faugeras and Pagès, 2024], who formalized the search for optimality within this setting. Our main contribution consists in extending this framework to the three-point case, which is particularly relevant from a practical perspective, as it captures the essential structure of risk profiles through three representative outcomes: “no loss,” “small loss” (with medium-to-high probability), and “extreme loss” (with low probability). This setting is especially meaningful in the insurance field, where such discrete representations of risk are commonly employed.

The magnitude–propensity framework was tested across three distinct and relevant use cases, with the aim of assessing both the behavior of the numerical procedures and the practical implementation of the results. In the unconstrained case, the optimization is free to explore the parameter space for the optimal configuration (m_1, m_2, p_1, p_2) . We then introduced a constraint on m_2 to align with regulatory requirements—specifically, the prescribed confidence level and time horizon associated with the Value-at-Risk (VaR) measure—thus enforcing $m_2 \geq VaR$.

In the unconstrained setting, the numerical procedures exhibit mutual convergence and match the analytical fixed-point solution derived from the theoretical framework. This confirms that (i) the numerical algorithms are robust and reliable, and (ii) the theoretical formulation—yielding substantial computational savings—is correct. In the constrained case, where a theoretical benchmark is not yet available, the numerical methods still converge to consistent results. From a practical perspective, running multiple procedures in parallel may serve as a valuable cross-validation strategy to ensure stability and reliability of the optimal solution.

Finally, we note that the computational performance of the fixed-point approach is excellent, with execution times close to zero, confirming its efficiency and suitability for large-scale or real-time applications.

In the risk management process, an essential step is capital allocation, that is, the ex-ante assignment of the risk budget to individual business units and the ongoing monitoring to verify the risk reward performance of each business owner. This task requires rigorous methods of risk decomposition, particularly in the case of financial conglomerates or banking groups. A seminal reference in this area includes [Garman, 1997], see also [Tasche, 2002] and [Tasche, 2004] for a formal rigorous framework.

As an initial approach, we consider it useful to extend the existing literature by treating the discrete variables m_1 and m_2 as if they represented the true loss distribution, allowing the application of the above mentioned techniques. Roughly speaking, one has to calculate the expected loss of each business unit conditioned on the total loss of the parent portfolio, i.e., VaR in the classical approach, while m_2 in our setting. Under this perspective, risk decomposition reduces to the computation of a conditional expectation with respect to the discrete values assumed by the distribution.

We believe that our approach could be evaluated from a managerial perspective, not only as a tool for risk measurement and reporting, but also as a useful instrument to support the quantification of strategic parameters for financial institutions, such as risk appetite and risk tolerance.

A Proofs

A.1 Proof of Lemma 3.2

1. Fix x such that, w.l.o.g., $h_0(x) < \min(h_1(x); h_2(x))$ and define $d := \min(h_1(x); h_2(x)) - h_0(x) > 0$. Take $0 < \epsilon_1 < d, 0 < \epsilon_2 < d - \epsilon_1$ (that is, $\epsilon_1 + \epsilon_2 < d$). Since h_1, h_2 are continuous, also $h_1 \wedge h_2$ is continuous, then there exists $\delta > 0$ such that $\forall v \in \mathbb{R}^2$ such that $\|v\| < \delta$ we have

$$\begin{aligned} h_0(x + v) &\leq h_0(x) + \epsilon_1, \\ h_1(x + v) \wedge h_2(x + v) &\geq h_1(x) \wedge h_2(x) - \epsilon_2. \end{aligned}$$

Now, from $h_0(x) = h_1(x) \wedge h_2(x) - d$ we get $h_0(x) + \epsilon_1 = h_1(x) \wedge h_2(x) + \epsilon_1 - d$, so that

$$h_0(x + v) \leq h_1(x) \wedge h_2(x) + \epsilon_1 - d < h_1(x) \wedge h_2(x) - \epsilon_2 \leq h_1(x + v) \wedge h_2(x + v).$$

Therefore $H(x + v) = h_0(x + v) \quad \forall \|v\| < \delta$, so that H is differentiable, with derivative given by $\nabla H^B(x)(v) = \nabla H(x)v = \nabla h_0(x)v$.

2. Take w.l.o.g. $h_0(x) = h_1(x) < h_2(x)$, so that $H(x) = h_0(x) \wedge h_1(x)$. By repeating the same reasoning of the previous point, there exists $\delta > 0$ such that for all $\|v\| < \delta$

$$H(x + v) = h_0(x + v) \wedge h_1(x + v).$$

Now, from Lemma 3.3 in [Faugeras and Pagès, 2024] we have that $h_0 \wedge h_1$ is B-differentiable, so that H is also B-differentiable and $\nabla^B H(x)(v) = \nabla h_1(x)v$.

3. We have $H(x) = h_0(x) = h_1(x) = h_2(x)$, so that

$$\begin{aligned} H(x + v) - H(x) &= \min(h_0(x + v) - h_0(x); h_1(x + v) - h_1(x); h_2(x + v) - h_2(x)) \\ &= \min(h_0(x + v) - h_0(x); \min(h_1(x + v) - h_1(x); h_2(x + v) - h_2(x))). \end{aligned}$$

Now, using the inequality $|\min(a; b) - \min(c; d)| \leq \max(|a - c|; |b - d|)$ with

$$\begin{aligned} a &= h_0(x + v) - h_0(x), \\ b &= \min(h_1(x + v) - h_1(x); h_2(x + v) - h_2(x)), \\ c &= \nabla h_0(x)v, \\ d &= \nabla^B(h_1 \wedge h_2)(x)(v), \end{aligned}$$

we get

$$\begin{aligned} &|H(x + v) - H(x) - \min(\nabla h_0(x)v; \nabla^B(h_1 \wedge h_2)(x)(v))| \\ &\leq \max(|h_0(x + v) - h_0(x) - \nabla h_0(x)v|; |\min(h_1(x + v) - h_1(x); h_2(x + v) - h_2(x)) - \nabla^B(h_1 \wedge h_2)(x)(v)|) \\ &= \max(|h_0(x + v) - h_0(x) - \nabla h_0(x)v|; |h_1 \wedge h_2(x + v) - h_1 \wedge h_2(x) - \nabla^B(h_1 \wedge h_2)(x)(v)|) \\ &= \max(\|v\|o(1); |h_1 \wedge h_2(x + v) - h_1 \wedge h_2(x) - \min(\nabla h_1(x)v; \nabla h_2(x)v)|) \\ &= \|v\| \max(o(1); |o(1)|), \end{aligned}$$

therefore

$$\begin{aligned} \nabla^B H(x)(v) &= \min(\nabla h_0(x)v; \min(\nabla h_1(x)v; \nabla h_2(x)v)) \\ &= \min(\nabla h_0(x)v; \nabla h_1(x)v; \nabla h_2(x)v), \end{aligned}$$

and the proof is complete.

A.2 Proof of Theorem 3.3

The proof is organized in two steps.

A.2.1 Step 1: B-derivatives of the function H

Let $H(x_1, x_2) = X^2 \wedge (X - x_1)^2 \wedge (X - x_2)^2$ with $0 < x_1 < x_2$. Fix X and set $h_0(x_1, x_2) = X^2$; $h_1(x_1, x_2) = (X - x_1)^2$; $h_2(x_1, x_2) = (X - x_2)^2$. We have then the following seven possibilities.

1. $h_0(x) = h_1(x) = h_2(x)$ for $x = (x_1, x_2) \in \{(0, 0), (0, 2X), (2X, 0), (2X, 2X)\}$. In this case we have $H(x_1, x_2) = X^2$ and the B-derivatives are given respectively by:

$$\begin{aligned}\nabla^B H(0, 0)(v) &= \min(0; -2Xv_1; -2Xv_2), \\ \nabla^B H(0, 2X)(v) &= \min(0; -2Xv_1; 2Xv_2), \\ \nabla^B H(2X, 0)(v) &= \min(0; 2Xv_1; -2Xv_2), \\ \nabla^B H(2X, 2X)(v) &= \min(0; 2Xv_1; 2Xv_2).\end{aligned}$$

2. $h_0(x) = h_1(x) < h_2(x)$ for $x = (x_1, x_2)$ such that $(x_1 = 0 \text{ and } x_2 > 2X)$ or $(x_1 = 2X \text{ and } x_2 > 2X)$. In this case we still have $H(x_1, x_2) = X^2$ and the B-derivative is given by:

$$\begin{aligned}\nabla^B H(0, x_2)(v) &= \min(0; -2Xv_1), \\ \nabla^B H(2X, x_2)(v) &= \min(0; 2Xv_1),\end{aligned}$$

provided that $x_2 > 2X$.

3. $h_0(x) = h_2(x) < h_1(x)$ for $(x_1 > 2X \text{ and } x_2 = 0)$ or $(x_1 > 2X \text{ and } x_2 = 2X)$. In this case we still have $H(x_1, x_2) = X^2$ and the B-derivative is given by:

$$\begin{aligned}\nabla^B H(x_1, 0)(v) &= \min(0; -2Xv_2), \\ \nabla^B H(x_1, 2X)(v) &= \min(0; 2Xv_2),\end{aligned}$$

provided that $x_1 > 2X$.

4. $h_1(x) = h_2(x) < h_0(x)$ for $(x_1 + x_2 = 2X \text{ and } 0 < x_1 < x_2 < 2X)$. This time we have $H(x_1, x_2) = (X - x_1)^2 = (X - x_2)^2$ and the B-derivative is given by:

$$\nabla^B H(x_1, 2X - x_1)(v) = \min(-2(X - x_1)v_1; -2(X - x_2)v_2),$$

provided that $x_1 + x_2 = 2X$ and $0 < x_1 < x_2 < 2X$.

5. $h_0(x) < h_1(x) \wedge h_2(x)$ for $x_1 > 2X$. We have $H(x_1, x_2) = X^2$ and the B-derivative is zero.

6. $h_1(x) < h_0(x) < h_2(x)$ for $(x_1 < 2X \text{ and } x_1 + x_2 > 2X)$. This time we have $H(x_1, x_2) = (X - x_1)^2$ and the B-derivative is given by:

$$\nabla^B H(x_1, x_2)(v) = -2(X - x_1)v_1,$$

provided that $x_1 < 2X$ and $x_1 + x_2 > 2X$.

7. $h_2(x) < h_0(x) < h_1(x)$ for $(x_1 + x_2 < 2X)$. This time we have $H(x_1, x_2) = (X - x_2)^2$ and the B-derivative is given by:

$$\nabla^B H(x_1, x_2)(v) = -2(X - x_2)v_2,$$

provided that $x_1 + x_2 < 2X$.

A.2.2 Step 2: B-differentiability of the objective function D and critical points for the distortion

Set

$$\nabla^B D(x)(v) := \mathbb{E}[\nabla^B H(x)(v)], \quad (16)$$

and consider now the different possibilities listed in the previous subsection.

1. The point $x = (0, 0)$ is critical for the distortion function if $\nabla^B D(0, 0)(v) \geq 0 \forall v \in \mathbb{R}^2$. Now,

$$\begin{aligned} \nabla^B D(0, 0)(v) &= \mathbb{E}[\min(0; -2Xv_1; -2Xv_2)] \\ &= \mathbb{E}[\min(-2Xv_1; -2Xv_2)] < 0 \end{aligned}$$

for $v_1 > 0, v_2 > 0$, therefore the point $x = (0, 0)$ is not critical for the distortion. The same holds true for $(0, 2X)$, since $\nabla^B D(0, 2X)(v) = \mathbb{E}[\min(0; -2Xv_1; -2Xv_2)] < 0$ for $v_1 > 0, v_2 < 0$. Using the same argument one can exclude also $(2X, 0), (2X, 2X)$.

2. Consider now $x_2 > 2X$, we get

$$\begin{aligned} \nabla^B D(0, x_2)(v) &= \mathbb{E}[\min(0; -2Xv_1)] < 0 \quad \text{for } v_1 > 0, \\ \nabla^B D(2X, x_2)(v) &= \mathbb{E}[\min(0; 2Xv_1)] < 0 \quad \text{for } v_1 < 0, \end{aligned}$$

so that the $(0, x_2), (2X, x_2)$ are not critical for $x_2 > 2X$.

3. Let $x_1 > 2X$,

$$\begin{aligned} \nabla^B D(x_1, 0)(v) &= \mathbb{E}[\min(0; -2Xv_2)] < 0 \quad \text{for } v_2 > 0, \\ \nabla^B D(x_1, 2X)(v) &= \mathbb{E}[\min(0; 2Xv_2)] < 0 \quad \text{for } v_2 < 0, \end{aligned}$$

therefore $(x_1, 0), (x_1, 2X)$ are not critical for $x_1 > 2X$.

4. Consider $x_1 + x_2 = 2X$ and $0 < x_1 < x_2 < 2X$:

$$\nabla^B D(x_1, 2X - x_1)(v) = \mathbb{E}[\min(-2(X - x_1)v_1; -2(X - x_2)v_2)] < 0 \quad \text{for } v_1 > 0, v_2 < 0,$$

therefore $(x_1, 2X - x_1)$ is not critical for $x_1 + x_2 = 2X$ and $0 < x_1 < x_2 < 2X$.

5. We have $\nabla^B D(x_1, x_2)(v) = 0$ for $x_1 > 2X$.

6. Take $x_1 < 2X$ and $x_1 + x_2 > 2X$:

$$\begin{aligned} \nabla^B D(x_1, x_2)(v) &= \mathbb{E}[-2(X - x_1)v_1 \mathbf{1}_{x_1 < 2X} \mathbf{1}_{x_1 + x_2 > 2X}] \\ &\quad + \mathbb{E}[\min(0; 2Xv_1) \mathbf{1}_{x_1 = 2X} \mathbf{1}_{x_2 > 2X}] \\ &\quad + \mathbb{E}[\min(-2(X - x_1)v_1; -2(X - x_1)v_2) \mathbf{1}_{x_1 < 2X} \mathbf{1}_{x_1 + x_2 = 2X}], \end{aligned}$$

so that (x_1, x_2) may be a critical point.

7. Finally, let consider $x_1 + x_2 < 2X$:

$$\nabla^B D(x_1, x_2)(v) = \mathbb{E}[-2(X - x_2)v_2 \mathbf{1}_{x_1 + x_2 < 2X}],$$

then (x_1, x_2) may be a critical point.

In conclusion, it turns out that the function $v \rightarrow \nabla^B D(x)(v)$ is positively homogeneous for $x \in \mathbb{R}_+^2$, hence D is B-differentiable. Moreover, collecting the different cases one gets (9), and the proof of the theorem is complete.

B Intrinsic Instability of the Quantile in Credit Risk Modeling

Extreme quantiles such as the 99.9% VaR play a central role in the computation of the *Default Risk Charge* (DRC). However, when applied to concentrated credit portfolios, these percentiles may exhibit a structural instability that does not originate from Monte Carlo error, but from the intrinsic combinatorial geometry of the loss distribution. Before analysing this phenomenon, we introduce the simplified analytical setup used to isolate and quantify such behavior.

Analytical setup. Consider a reduced set of the N largest obligors (covering the majority of portfolio EAD) and model defaults via a one-factor Gaussian copula. For obligor $n = 1, \dots, N$, we define

$$Z_n = \sqrt{\rho_n} U + \sqrt{1 - \rho_n} \varepsilon_n,$$

where U, ε_n are independent standard Gaussian random variables, and a default occurs whenever Z_n falls below the threshold given by the corresponding default probability, i.e. $b_n := \Phi^{-1}(\text{PD}_n)$. The portfolio loss in 2 is therefore

$$\text{Loss}_{\text{DRC}} = \sum_{n=1}^N \text{EAD}_n \cdot \mathbf{1}_{\{Z_n \leq b_n\}} \cdot \text{LGD}_n,$$

with deterministic LGD in this experiment.

Since each obligor may be in default or not, the loss distribution is supported on the 2^N possible default/non-default combinations. For each configuration $r \subset \{1, \dots, N\}$, the corresponding loss is

$$\text{Loss}_r = \sum_{n \in r} \text{EAD}_n \cdot \mathbf{1}_{\{Z_n \leq b_n\}} \cdot \text{LGD}_n,$$

while its probability under the copula reduces to a one-dimensional integral:

$$P_r = \int_{-\infty}^{+\infty} \left[\prod_{n \in r} \Phi\left(\frac{b_n - \sqrt{\rho_n} u}{\sqrt{1 - \rho_n}}\right) \right] \left[\prod_{n \notin r} \left(1 - \Phi\left(\frac{b_n - \sqrt{\rho_n} u}{\sqrt{1 - \rho_n}}\right)\right) \right] \phi(u) du, \quad (17)$$

evaluated via Gaussian quadrature. Sorting all (Loss_r, P_r) pairs yields the exact loss distribution and, in particular, the exact 99.9% Default Risk Charge DRC is defined as

$$\text{DRC} = \min \{ \text{Loss}_r : P_{\text{cum}}(r) \geq 0.999 \},$$

where $P_{\text{cum}}(r) = \sum_{|m| \leq |r|} P_m$, where the sum is extended to incompatible events.

Focusing on the percentile instability. To analyse how the percentile behaves in the tail, we examine the symmetric window

$$\{k - 500, \dots, k + 500\},$$

where k is the index of the order statistic corresponding to the 99.9% quantile. This window contains the ranked scenarios whose cumulative probabilities lie immediately below and above 0.999—precisely where small perturbations may change the identity of the selected quantile.

Figure 6 displays, for each scenario in the quantile window, (i) the default pattern across the $N = 31$ obligors (orange dots), (ii) the cumulative probability α_r , and (iii) the corresponding loss Loss_r . The figure highlights a crucial structural phenomenon: *despite the fact that default patterns differ significantly across scenarios, the resulting losses in this tail region are numerically almost identical*. In other words, the instability is not related to the magnitude of the loss, but rather to the specific configuration of the defaulting obligors. This arises because many combinations of $\text{EAD}_n \text{LGD}_n$ sum to nearly the same value, an effect amplified by the combinatorial explosion of the support.

optimal transport problem reads

$$\min_{\Pi \in \mathbb{R}_+^{n \times m}} \left\{ \langle C, \Pi \rangle + \varepsilon \sum_{i=1}^n \sum_{j=1}^m \Pi_{ij} (\log \Pi_{ij} - 1) \right\} \quad \text{s.t.} \quad \Pi \mathbf{1}_m = \mathbf{p}, \quad \Pi^\top \mathbf{1}_n = \mathbf{q}, \quad (18)$$

where $\varepsilon > 0$ is the regularization parameter, $\langle C, \Pi \rangle = \sum_{ij} c_{ij} \Pi_{ij}$, and $\mathbf{1}_k$ denotes the k -dimensional vector of ones.

Introducing the kernel matrix

$$K_{ij} = \exp\left(-\frac{c_{ij}}{\varepsilon}\right), \quad i = 1, \dots, n, \quad j = 1, \dots, m,$$

it is well known that any solution of (18) can be written in the scaling form

$$\Pi^\star = \text{diag}(\mathbf{u}) K \text{diag}(\mathbf{v}), \quad (19)$$

for some positive vectors $\mathbf{u} \in \mathbb{R}_{++}^n$ and $\mathbf{v} \in \mathbb{R}_{++}^m$. The Sinkhorn–Knopp algorithm is an iterative matrix scaling procedure that computes (\mathbf{u}, \mathbf{v}) such that the marginal constraints in (18) are satisfied.

Let $\mathbf{u}^{(0)} \in \mathbb{R}_{++}^n$ and $\mathbf{v}^{(0)} \in \mathbb{R}_{++}^m$ be initial scaling vectors, typically chosen as vectors of ones. At iteration $\ell = 0, 1, 2, \dots$, the algorithm performs the updates

$$u_i^{(\ell+1)} = \frac{p_i}{\sum_{j=1}^m K_{ij} v_j^{(\ell)}}, \quad i = 1, \dots, n, \quad (20)$$

$$v_j^{(\ell+1)} = \frac{q_j}{\sum_{i=1}^n K_{ij} u_i^{(\ell+1)}}, \quad j = 1, \dots, m \quad (21)$$

The sequence $\{(\mathbf{u}^{(\ell)}, \mathbf{v}^{(\ell)})\}_{\ell \geq 0}$ generated by (20)–(21) defines a sequence of transport plans

$$\Pi^{(\ell)} = \text{diag}(\mathbf{u}^{(\ell)}) K \text{diag}(\mathbf{v}^{(\ell)}),$$

whose row and column sums converge to \mathbf{p} and \mathbf{q} , respectively, under mild assumptions on K (e.g. all entries strictly positive). In practice, the iterations are stopped when the violation of the marginals is below a prescribed tolerance $\tau > 0$, i.e. when

$$\left\| \Pi^{(\ell)} \mathbf{1}_m - \mathbf{p} \right\|_1 + \left\| (\Pi^{(\ell)})^\top \mathbf{1}_n - \mathbf{q} \right\|_1 \leq \tau. \quad (22)$$

For numerical stability, in particular when ε is small and the entries of K may underflow, the updates (20)–(21) can be implemented in the log-domain. Defining $\alpha^{(\ell)} = \log \mathbf{u}^{(\ell)}$ and $\beta^{(\ell)} = \log \mathbf{v}^{(\ell)}$, one replaces the matrix–vector products by log-sum-exp operations, which reduces the risk of numerical underflow/overflow while leaving the fixed point (19) unchanged.

Once convergence has been reached (according to (22)), the transport plan Π^\star in (19) is used to compute the (regularized) optimal transport cost

$$\mathcal{W}_\varepsilon(\mathbf{p}, \mathbf{q}) = \langle C, \Pi^\star \rangle + \varepsilon \sum_{i,j} \Pi_{ij}^\star (\log \Pi_{ij}^\star - 1),$$

which is then employed in the subsequent stages of the analysis.

D The Differential Evolution (DE) Algorithm

In this appendix, we provide some details on the numerical procedure we adopted for the optimization problem.

The Differential Evolution (DE) algorithm, originally proposed by [Storn and Price, 1997], is a stochastic and population-based global optimization technique belonging to the family of evolutionary algorithms. It is particularly effective for solving non-convex, non-differentiable, and multi-modal optimization problems. Unlike gradient-based methods, DE relies exclusively on function evaluations, which makes it robust even when the objective function $f(\mathbf{x})$ is noisy or discontinuous.

The algorithm evolves a population of candidate solutions across generations through a sequence of variation mechanisms — *mutation*, *crossover*, and *selection*. At each iteration, these operations allow the population to progressively explore the search space and concentrate around high-performing regions, eventually converging to a near-optimal solution for the problem under consideration.

The optimization problem is defined as

$$\min_{\mathbf{x} \in \Omega} f(\mathbf{x}), \quad (23)$$

where $\Omega \subseteq \mathbb{R}^d$ denotes the feasible domain of the d -dimensional decision vector \mathbf{x} . The approach consists in evolving a set of candidate solutions over successive generations.

At generation g , the algorithm maintains a population of N_P candidate solutions,

$$\mathbf{x}_{i,g} = (x_{i,g}^{(1)}, x_{i,g}^{(2)}, \dots, x_{i,g}^{(d)}), \quad i = 1, 2, \dots, N_P, \quad (24)$$

which evolve iteratively through mutation, crossover, and selection.

Mutation. For each target vector $\mathbf{x}_{i,g}$, a mutant vector $\mathbf{v}_{i,g}$ is generated according to the classical DE/rand/1/bin scheme,

$$\mathbf{v}_{i,g} = \mathbf{x}_{r_1,g} + F(\mathbf{x}_{r_2,g} - \mathbf{x}_{r_3,g}), \quad (25)$$

where $r_1, r_2, r_3 \in \{1, \dots, N_P\}$ are distinct indices different from i , and $F \in [0, 2]$ is the mutation factor controlling the amplification of the differential variation. Equation (25) therefore perturbs a base vector $\mathbf{x}_{r_1,g}$ by adding a scaled difference of two other individuals, thereby promoting exploration of the search space.

Crossover. To increase population diversity, a trial vector $\mathbf{u}_{i,g}$ is formed by combining components of the mutant and target vectors:

$$u_{i,g}^{(j)} = \begin{cases} v_{i,g}^{(j)}, & \text{if } \text{rand}_j(0, 1) \leq C_R \text{ or } j = j_{\text{rand}}, \\ x_{i,g}^{(j)}, & \text{otherwise,} \end{cases} \quad j = 1, \dots, d, \quad (26)$$

where $C_R \in [0, 1]$ is the crossover probability and j_{rand} is a randomly chosen index ensuring that at least one component originates from the mutant vector. Hence, (26) mixes the information contained in $\mathbf{x}_{i,g}$ and $\mathbf{v}_{i,g}$, with the parameter C_R regulating the expected proportion of mutated components.

Selection. Selection is then performed to determine whether the trial vector replaces the target vector in the next generation:

$$\mathbf{x}_{i,g+1} = \begin{cases} \mathbf{u}_{i,g}, & \text{if } f(\mathbf{u}_{i,g}) \leq f(\mathbf{x}_{i,g}), \\ \mathbf{x}_{i,g}, & \text{otherwise.} \end{cases} \quad (27)$$

According to (27), the better (or equal) solution survives, which guarantees that the population's overall fitness does not deteriorate over successive generations. The evolutionary process continues until a stopping

criterion is met, such as a maximum number of generations G_{\max} , or when the relative improvement in the best fitness value falls below a threshold $\varepsilon > 0$, i.e.,

$$\frac{|f_{\text{best},g} - f_{\text{best},g-1}|}{|f_{\text{best},g-1}|} < \varepsilon. \quad (28)$$

Condition (28) prevents unnecessary iterations once the algorithm has essentially converged.

To apply the approach to empirical data in our context, the algorithm is used to solve the nonlinear systems defining the quantities (m_x, p_x) and (m_1, m_2, p_1, p_2) for the 2-point and 3-point cases, respectively. In a historical simulation setting, the vector of the S P&Ls is interpreted as the empirical distribution of the underlying risk factor. Each point in the vector is therefore assigned the probability mass

$$p_s = \frac{1}{S}, \quad s = 1, \dots, S, \quad (29)$$

with $S = 250$ or $S = 500$ in most of our applications.

Once the P&Ls have been reordered, we denote by $\text{PnL}(s)$ the empirical cumulative distribution, which can be written as

$$F(\text{PnL}(s)) = \frac{s}{S}, \quad s = 1, \dots, S, \quad (30)$$

while for any generic point $z \in \mathbb{R}$ the empirical distribution function is

$$F(z) = \frac{\#\{s : \text{PnL}_s \leq z\}}{S}. \quad (31)$$

Equations (29)–(31) provide the empirical probabilities and cumulative distribution that constitute the input to the objective function optimized by the DE algorithm in (23).

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