# Multivariate Distribution Model for Financial Risks Management

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**Abstract.** The method for constructing joint distribution copula based models is proposed. The copula model parameters are estimated by the method of maximum likelihood which turned out to be effective according to the mean squared error criterion. The model for estimating risks of various types is proposed and constructed using copula function approach. Higher quality of the tail risk measures was achieved for the data samples selected that leads to the necessity of improvement formal description for the central part of observations. This is true when the model is used based on combined marginal distributions along with normal and generalized Pareto distributions. Results of computational experiments are also provided.

**Keywords:** Financial Risks, Copula, Marginal Distribution, Risks Analysis, Risks Measures, Value-at-Risk

## **1** Introduction

Risk estimation and management can be performed by changing the structure of financial instruments portfolio that requires adequate mathematical model constructing for multivariable processes. One of the possible approaches is based on separate modeling of marginal distributions and dependency structure between corresponding variables by using of special copula functions [1 - 4]. The adequacy criterion for such models is based on the quality of risk measures estimation [5 - 9]

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that provide a quantitative characteristics of risk degree, and it is convenience for applying such traditional procedures for risk management as scenario analysis [10, 11]. The complexity of constructing joint distribution for several variables that belong to different types of risks imposes restrictions on the procedures of analytical expressions development for estimating appropriate risk measure. It should be stressed that using of Monte Carlo procedures to obtain quality estimates of risk measures [12, 13] requires application of effective methods for generating appropriate pseudorandom numbers (PRN) by making use of such special functions as copula [13–16].

# 2 Problem statement

In the frames of this study it is planned to solve the following problems: (1) to formulate systemic approach to risks analysis for multidimensional portfolio of financial instruments (PFI); (2) to determine possible applications of widely used risk measures when the portfolio under risk is composed; (3) to develop theoretically substantiated risk measures for the PFI risk measures on the basis of probabilistic and statistical modeling using the outlier theory and copulas; (4) to determine the possibility and effectiveness of practical application of proposed methods using actual statistical data.

# 3 Mathematical model of risks distribution

To perform statistical analysis of financial risk level the models are used on the basis of combined marginal distributions and copulas. The joint distribution for possible risks is considered in the following form:

 $H(x_1, ..., x_n) = P[X_1 \le x_1, ..., X_n \le x_n] = C(F_1(x_1), ..., F_n(x_n))$ 

Here,  $F_1, ..., F_n$ , are marginal distribution functions for separate risks; C is n-copula, that characterizes the dependency structure between the analyzed risks. Application of the model is oriented to elliptical copulas, Archimedes copulas, and extreme value copulas. The right tail of marginal distribution is formally described by the distribution of a threshold exceeding values in the form of generalized Pareto distribution (GPD) as follows:

$$GPD_{\xi,\beta}(x) = \begin{cases} 1 - (1 + \xi \ x / \beta)^{-1/\xi}, & \xi \neq 0\\ 1 - \exp(-x / \beta), & \xi = 0, \end{cases}$$

where,  $\beta > 0$ , and,  $x \ge 0$ , with  $\xi > 0$ , and,  $0 \le x \le \frac{-\beta}{\xi}$ , with  $\xi < 0$ ;  $\xi$  is parameter of the distribution form;  $\beta$  is scale parameter. The central data observations are modeled with normal distribution.

# 4 Scenario analysis

One of the well-known approaches to formal risks description is based upon computing PFI cost for the risk positions under various market conditions. This approach is called scenario analysis. The scenario analysis approach has the following two basic types: shock testing and sensitivity analysis.

The shock testing is dedicated to analysis of consequences provoked by the sudden and substantial variations of market processes. It is especially useful in the cases when the market movements are resulted in qualitative changes of some market conditions. Actually the shock testing is based on the process of creating and analysis of such situations that could be helpful in identifying extreme portfolio risks and implementing the scenarios in a simulation model for estimating the qualitative and quantitative risk measures.

The scenario development is performed on the basis of observations received during former crisis events, the observations coming from other spheres of activities, or realistic, potentially extreme scenarios. Systemic approaches to creating scenarios of this type do not exist at the moment. It is necessary to create a viable model of risks to analyze results of the scenarios implementation that are characterized by substantial quantitative and qualitative deviations of the processes under study from normal mode of functioning. The model on the basis of existing link functions and combined marginal distributions provides a possibility for variations in the scenarios regarding the parameters of separate risk distributions and utilize the types of marginal distributions that differ from the types of distributions characteristic for the basic accepted model. It is also possible to incorporate extreme modifications into the process under study through the values of tail distribution parameters and vary the dependency characteristics between the risks. However, the shock testing does not provide the possibility for estimating probabilities of the scenarios that result in high losses. So, its application in risk management systems is restricted by taking into account the risks for each potential risky event separately.

The sensitivity analysis provides a possibility for estimating the risk characteristic of a portfolio in non-critical market conditions when actual modifications are not substantial. This type of analysis is used in conditions of availability of a large number of realistic scenarios that do not predict/suppose extreme events. Using of the models based upon link functions and combined marginal distributions provides a possibility for changing not only for distribution parameters but also the non-tail parts of marginal distributions can be changed leaving unchanged the model for extreme values that is constructed on the data generated within longer periods of time and very often is weakly depending upon the distribution for central values.

# 5 The risk measures

The quantitative estimate of risk with using specific measures is one of the basic tasks of risk management due to necessity of comparing the risks provoked by different decision alternatives. On one side, the risk can be considered as the losses or shortened income that is determined by the future value of a certain position. On the other side, the risk is provoked by the high volatility of value for some positions.

The examples of risks dynamical models are shown for credit and outflow risks in [17, 18] with risk degree and level estimation through the probability and size of the losses. The risk measures that are based upon understanding the risk as a future cost of an instrument are called coherent. The necessary requirements to the coherent risk measure,  $\rho$ , were formulated in the form of axioms in [19]. These requirements are of financial nature and are derived from the rules of accepting risks proposed by investment managers as well as practically induced regulations. The non-subadditive coherent risk measures should satisfy the axioms 1 - 3 given below.

Let X and Y are random variables that reflect future values for portfolio composed of market positions. The portfolio components which exhibit higher future values than positions of another portfolio is considered as a less risky one.

Axiom 1. Monotonicity: if  $X \ge Y$ , then,  $\rho(X) \le \rho(Y)$ . Investing into the stock instrument with known return level, r, for which the future value is deterministic and known, decreases the risk by the invested value.

Axiom 2. Invariance to bias: for the constant a,  $\rho(X + a \cdot (1 + r)) = \rho(X) - a$ , the risk is growing with growth of risky position value. For large value positions the simple growth of risk (according to the growth of invested capital) is added to the growth of liquidity risk. The large position cannot be realized at the market as easy as the small one. However, the risk measure for small position should be proportional to the value of the position.

**Axiom 3.** The positive homogeneity:  $\rho(bX) = b\rho(X)$  for any positive number, b.

The measures of risk deviation,  $\delta$ , should be defined by some other axioms [20]. The non-subadditive measures of risk deviation satisfy the axioms 4 - 6. The characteristic of deviations for non-deterministic risk X should always be positive and equal to zero in a case when X is a constant.

**Axiom 4.**  $\forall X, \quad \delta(X) \ge 0$ .

Investing into stock instrument the future value of which is known thanks to the level of return, r, that is not volatile should not influence the characteristic of portfolio deviations.

**Axiom 5.**  $\delta(X + a \cdot (1 + r)) = \delta(X)$ .

For the market of complete liquidity increase of the investment volume results in proportional growth of cost deviation.

**Axiom 6.**  $\forall X, \forall \lambda > 0 \quad \delta(\lambda X) = \lambda \delta(X)$ .

This subadditivity axiom is based upon suggestion that diversification of the stock instruments results in decreasing of risks. Besides, the subadditive risk measures are convenient for the use at the enterprise level, as far as they guaranty that integrated risk will not exceed the sum of risks characteristic for separate divisions, portfolios or positions.

Axiom 7. The subadditivity of coherent risk measures is determined as follows:  $\rho(X + Y) \le \rho(X) + \rho(Y)$ . **Axiom 8.** The subadditivity of measures for risk deviation is determined as follows:  $\delta(X + Y) \le \delta(X) + \delta(Y)$ .

Holding of axiom 7 is necessary condition for the coherent risk measures (subadditive coherent risk measures); and axiom 8 for the deviation measures of risks (subadditive measures of risk deviation).

The axioms 1 - 3 and subadditivity axiom could be used to derive an expression for the coherent risk measures of the following type [19]:

$$\rho(X) = \sup_{\mu \in \Psi} E_{\mu} \left[ \frac{-X}{1+r} \right], \tag{1}$$

where,  $\Psi$ , is a family of probabilistic measures. Thus, any coherent measure of risk is defined by the mathematical expectation of maximum loss for a certain set of scenarios. As it can be seen from (1) extension of the scenario set leads to enhancement of the risk measure too. The fact that the risk measures depend on the set of scenarios proves practical validity of the scenario analysis approach in the risk management systems. The coherent risk measures for which the following inequality holds:  $\rho(X) > E[-X]$ , when X is not a constant, and,  $\rho(X) < E[-X]$ , when X is a constant are called restricted expectations. Between coherent risk measures that are restricted by expectation and risk deviation measures there exists one-to-one correspondence of the form:

$$\delta(X) = \rho(X - EX),$$
  

$$\rho(X) = E[-X] + \delta(X).$$

If these conditions hold, then  $\delta$  is a measure of risk deviation linked to the coherent measure of risk,  $\rho$ , and,  $\rho$ , is a measure of risk linked to  $\delta$ . The couple of the two measures together create a risk profile.

#### 6 Estimation of basic risk measures

The measure of Value-at-Risk (VaR). VaR, or value at risk is a response to extreme financial events and catastrophes. The initial purpose was to create quantitative measure of risk based on available statistical techniques. Thus, VaR provides for a probabilistic measure of potential loss pointing out to the threshold that could be exceeded with expected loss in normal market conditions on definite time horizon and given confidence level.

**Definition 1**. For a given value of  $\alpha \in (0,1]$  and random variable, X, the  $\alpha$  – quantile is defined by the expression:

$$q_{\alpha} = \{ x \in R \mid P[X \le x] \ge \alpha \}.$$

The values of quantiles with sufficient set of values for  $\alpha$  characterize sufficiently form and scatterplot of probabilities distribution.

**Definition 2.** The measure of risk VaR at confidence level  $\alpha$  for random variable (process) X (returns and losses where losses are negatively defined values) is formally defined as follows:

$$VaR_{\alpha}(X) = q_{\alpha}(-X)$$

If losses have distribution function, F, then  $VaR_{\alpha} = F^{-1}(\alpha)$ , where  $F^{-1}$  is inverse for F.

As a quantitative characteristic of short-term market risk, VaR is often used in management systems for analysis of credit and operational risks. This measure is one of the most important components of general methodology of quantitative risk estimation that is used in practice. The VaR-based instruments are used for solving the problems of investments and estimating the compromise between returns and risk. The measure satisfies the axioms 1 - 3 for the coherent risk measures but does not satisfy the axiom of subadditivity.

However, analytical estimation of VaR is not always convenient due to relative complexity of finding appropriate solution. This is true for the case considered in this study where joint distribution for risks with combined marginal distributions is used. That is why an empirical estimate will be used based on application of the Monte Carlo technique. If a sample of random values  $\{X_{i;i}\}$  is considered of power, n, that

is arranged in the way that  $X_{1:n} \le ... \le X_{n:n}$ , then empirical estimate of risk measure VaR is determined as follows:

$$VaR_{\alpha}(X) = X_{\max(i \in N | i \le n\alpha): n}.$$

The computations necessary for estimating this value include the following steps:

- 1. Selection of confidence level,  $\alpha$ .
- 2. Generating sufficiently large number of pseudorandom vectors according to previously estimated marginal distributions for the central and tail parts as well as link function for the joint distribution of losses constructed.
- Using the price models compute simulated samples of losses and returns for all instruments.
- 4. Computing VaR as the lowest losses for  $1-\alpha$  percentage of the worst cases.

The risk measure ES (Expected Shortfall). An alternative to VaR coherent risk measure is Expected Shortfall.

**Definition 3.** The measure of risk ES at confidence level  $\alpha$  for random variable *X* is determined as follows:

$$ES_{\alpha}(X) = -\frac{1}{1-\alpha} \Big( E\Big[ X \mid_{\{-X \ge q_{\alpha}(-X)\}} \Big] + q_{\alpha}(-X) \Big\{ \alpha - P[-X < q_{\alpha}(-X)] \Big\} \Big).$$

ES is mathematical expectation of losses on a given time horizon and confidence level; at the same time VaR represents minimum loss:

$$ES_{\alpha} = -E[X \mid X \leq -VaR_{\alpha}].$$

The measure ES is also called conditional VaR. An empirical estimate of ES is determined by the expression:

$$ES_{\alpha}(X) = -\frac{\sum_{k=1}^{\max(i \in N \mid i \le n \alpha)} X_{k:n}}{\max(i \in N \mid i \le n \alpha)}$$

**Markowitz' measures of risk.** First, Markowitz proposed to use as a measure of risk the standard deviation,  $\sigma(X) = |X - EX|$ , and then semi-deviation  $\sigma_{-}(X) = |[X - EX]_{-}|$  [5]. Both risk measures satisfy all four axioms for risk

deviation measures. In risk management systems it is more convenient to use the measure in the form of semi-deviation,  $\sigma_{-}(X)$ . For a sample,  $X_i$ , of power, n, the empirical estimates for risk deviation measures are the following:

– for standard deviation:

$$\hat{\sigma}(X) = \sqrt{\frac{1}{n} \sum_{k=1}^{n} \left( X_i - \frac{1}{n} \sum_{k=1}^{n} X_i \right)^2};$$

– and for semi-deviation:

$$\hat{\sigma}_{-}(X) = \sqrt{\frac{1}{n} \sum_{k=1}^{n} \left[ \left[ X_i - \frac{1}{n} \sum_{k=1}^{n} X_i \right]_{-} \right]^2}$$

The empirical estimates of the risk measure VaR or coherent risk measure ES and risk deviation measures together provide an estimate of risks portfolio.

Generating risk measures from copulas. When copulas are used in the risk management systems the Monte Carlo techniques should be used for estimating risk measures, risk deviation measures, and scenario analysis. It means generating pseudorandom dependent values with given marginal distributions and copula. There exists general scheme for sample generating from joint distribution, the structure of dependence in which is determined by copula. Using Sklyar theorem, the procedure of generating random sample,  $X_1,...,X_n$ , with marginal distribution functions,  $F_1,...,F_n$  and copula, C, is described as given below [21].

1. Generate random numbers  $u_1, ..., u_n$  with scalar uniform distributions with parameters, [0, 1], and joint distribution over copula C.

2. Compute the values of  $x_i$  using the transformation:

$$x_i = F_i^{-1}(u_i),$$

where,  $F_i^{-1}$ , is inverse for  $F_i$ .

Depending on the features of the copula family implementation of the step 1 can be simplified.

### Elliptical copulas.

**Definition 4.** The characteristic function  $\varphi: \mathbb{R}^n \to C$  for a random *n*-vector *X* is defined as follows:

$$\varphi_X(t) = E[e^{itX}],$$

where, tX, is coordinate-to-coordinate product of the vectors t and X. The characteristic function represents probabilistic distribution as a single-valued meaning.

**Definition 5** [22]. The probabilistic distribution for *n*-vector *X* is called elliptical if for a certain vector,  $\mu \in \mathbb{R}^n$ , positively defined matrix,  $\Sigma \in \mathbb{R}^{n \times n}$ , and some

function  $\phi:[0,\infty) \to R$ , the characteristic function of vector,  $X - \mu$ , is defined as follows:  $\phi_{X-\mu}(t) = \phi\left(t^T \Sigma t\right)$ .

The elliptical functions got their name due to the elliptical form of invariability lines for densities of the distributions. Any combination of elliptical distributions also creates elliptical distribution. The marginal distributions of joint elliptical distribution also belong to elliptical distributions. The elliptical copulas are defined through joint elliptical distributions, H, and inverse functions,  $F_i^{(-1)}(u_i)$ , as follows:

$$C(u_1,...,u_n) = H(F_1^{(-1)}(u_1),...,F_n^{(-1)}(u_n)).$$

Generating random numbers with the dependency structure defined by the elliptical copula is actually sampling from respective elliptical distribution. The algorithm given below generates PRN from Gaussian copula, corresponding to normal distribution, in the form of *n*-vector,  $(X_1,...,X_n)$ , with correlation matrix,  $\rho$ , as follows:

- 1. Generate *n* independent random numbers,  $u_1, ..., u_n$ , according to normal distribution.
- 2. Represent,  $\rho$ , in the form of decomposition,  $\rho = A \cdot A^T$ , where A is a low triangle matrix.
- 3. Compute,  $\vec{y} = A \cdot \vec{u}$ .
- 4. Using the scalar normal distribution function,  $\Phi$ , compute,  $x_i = \Phi(y_i)$ .

To generate PRN from more complex elliptical copulas any zero mean elliptically distributed vector, X, is represented via centered normally distributed vector, N, with the same correlation matrix, and independent on, N, random value, r: X = rN. For example, for Student *t*-distribution with  $\nu$  degrees of freedom, r has  $\chi^2$ -distribution with,  $\nu$ , degrees of freedom.

**The Archimedean copulas**. The copulas that can be represented in the form:  $C(u_1,...,u_n) = \varphi^{[-1]}(\varphi(u_1) + ... + \varphi(u_n))$ , are called Archimedean. The unified algorithm for generating the copulas is constructed on the following basis:

 $C(u_1,...,u_n) = C_n(u_n \mid u_1,...,u_{n-1})...C_2(u_2 \mid u_1)C_1(u_1),$ 

where,  $C_2$ , is copula for the first k components; and  $C_1(u_1) = u_1$ . To accomplish sampling from this joint distribution it is necessary to perform the following steps:

- Generate, n, independent PRN,  $v_1, ..., v_n$ ;

then, sequentially compute,  $u_1 = v_1$ ,  $u_2 = C_2^{-1}(v_2 | u_1)$ , ...,

 $u_n = C_n^{-1}(v_n \mid u_1, ..., u_{n-1}).$ 

For Archimedean copulas the computational algorithm can be simplified to the following form:

$$u_{2} = \varphi_{2}^{[-1]} \left( \varphi_{2} \left( \varphi_{1}^{[-1]} \left( \frac{\varphi_{1}(u_{1})}{v_{2}} \right) \right) - \varphi_{2}(u_{1}) \right).$$

**The extreme value copulas.** Let,  $(X_{i1}, ..., X_{im})$ , i = 1, 2, ..., are independent identically distributed *m*-vectors with distribution function, *F*, and let

$$M_{ij} = \max_{1 \le i \le n} X_{ij}, \quad j = 1, ..., m$$

represents maximum for each component. The multidimensional distribution for extreme values is a limit of the random vectors,  $\left(\frac{(M_{1n}-a_{1n})}{b_{1n}}, \dots, \frac{(M_{mn}-a_{mn})}{b_{mn}}\right)$ . If the

limit distribution exists, then each component of the distribution represents onedimensional distribution of extreme values that can be represented in the form:

$$C(H(z_1;\gamma_1),...,H(z_m;\gamma_m)),$$

where,  $H(z_j; \gamma_j)$ , is generalized distribution of extreme values, and, *C*, is a copula. It is natural that marginal distributions are related to the extreme value distributions, but more interesting is the case of the following copula:

$$G(z) = \lim_{n \to \infty} F^n(a_{1n} + b_{1n}z_1, ..., a_{mn} + b_{mn}z_m) = C(H(z_1; \gamma_1), ..., H(z_m; \gamma_m)).$$
(2)

Let,  $r_j$ , is strictly increasing transform for  $X_{ij}$ ; denote the variables and maximum after the transform as,  $X_{ij}^*$ , and  $M_{in}^*$ , respectively. Suppose also that expression,  $\frac{M_{in}^* - a_{jn}^*}{b_{jn}^*}$ , tends to the extreme value distribution with,  $n \to \infty$ , for all

j; and, let

$$G(z) = \lim_{n \to \infty} P(M_{1n}^* \le a_{1n}^* + b_{1n}^* z_1^*, ..., M_{mn}^* \le a_{mn}^* + b_{mn}^* z_m^*)$$
$$= C^* \Big( H(z_1^*; \gamma_1^*), ..., H(z_m^*; \gamma_m^*) \Big)$$

It can be shown that,  $C = C^*$ , and,

$$C(u_1^t, ..., u_m^t) = C^t(u_1, ..., u_m),$$
(3)

for all t > 0. The copulas that satisfy this condition are called extreme value copulas.

This set of copulas includes, for example, the family of two-dimensional Gumbel copulas of the form:

$$C_{\theta}(u,v) = \exp(-\left[\left(-\ln u\right)^{\theta} + \left(-\ln v\right)^{\theta}\right]^{\frac{1}{\theta}}),$$

where,  $\theta \ge 1$ . Sampling of PRN from extreme copulas is based on the use of universal generating procedures from known probabilistic distributions. To solve the problem it is proposed to apply multidimensional generalization of one-dimensional method of so called slice sampling [14].

The method of slice sampling. Suppose it is necessary to perform sampling from probabilistic distribution on some subset from,  $R^n$ , that is determined by some probability density function (PDF) proportional to some function, f(x). Such task can be accomplished by the uniform sampling from n+1-dimensional area under graph of the function, f(x). Formally this idea can be implemented by introducing

extra variable, y, and determining joint distribution for, x, and y, that is uniform over the area under the boundary described by the function f(x):

$$U = \{(x, y) : 0 < y < f(x)\}.$$

Thus, the joint distribution for (x, y) is determined as follows:

$$p(x, y) = \begin{cases} 1/\int f(x) \, dx, & 0 < y < f(x) \\ 0, & y \le 0, & y \ge f(x) \end{cases}.$$

Generating of independent values uniformly distributed over the set, U, is not simple task. That is why Markov chain is generated converging to this distribution. One of possible solutions is based upon Gibbs sampling, i.e. sampling from conditional distribution, P(y|x). The distribution is uniform over the interval, (0, f(x)), and from conditional distribution, P(x|y), that is uniform over the interval,  $S = \{x : y < f(x)\}$ . The resulting distribution is called a "slice" (or sector) defined over the variable, y. Rather complicated task can be generating of independent, uniformly distributed value from, S. It can be replaced by some innovation for, x, that doesn't violate uniform distribution over S.

Introduce the following notations: let f(x) is the function proportional to the density distribution we are looking for;  $x_0$  is current (initial) state;  $x_1$  is a new state. Now, the method for sampling in one-dimensional case can be formulated as given below [14].

1. Generate real value, y, uniformly from  $(0, f(x_0))$ , and this way get horizontal "slice",  $S = \{x : y < f(x)\}$ . Note, that  $x_0$  always belongs to, S.

2. Define the interval, I = (L, R), around  $x_0$  that contains at least large part of the slice.

3. Generate new point,  $x_1$ , that belongs to the interval defined, i.e. to the intersection,  $S \cap I$ .

On the first initialization step a value for an extra variable should be selected that is characteristic for the particular case of slice sampling. There is no need to save the value for further iterations. To avoid the problem touching upon precision of representing floating point numbers it is recommended to use the following function:  $g(x) = \log(f(x))$ , instead of f(x). In this case the extra variable is defined as follows:  $z = \log(y) = g(x_0) - e$ , where, e, is exponentially distributed value with unity mathematical expectation; and the slice is defined as  $S = \{x : z < g(x)\}$ . The second and the third steps of sampling in one-dimensional case could be implemented in various ways but the result should be in the form of Markov chain that does not affect the distribution defined by the function, f(x).

On the second step, the respective numerical interval is determined. It is desirable that the interval should include as larger part of the slice as possible. This is necessary for distinguishing the new point (number) from the previous one as much as possible. Though it is also necessary to avoid such intervals that would exceed the slice

substantially; the matter is that this results in less effective generating of the new point. The interval can be found in several possible ways given below.

It is ideally to have,  $L = \inf(S)$ , and,  $R = \sup(S)$ . It means that interval, I, 1. is equal to the least possible interval that contains the whole set, S. However, this situation is difficult to implement in practice.

If the values of, x, are restricted, then interval, I, may be defined over the 2. whole admissible range. However, this is not effective approach when the slice size is substantially less than the range.

3. Define size of the slice as,  $\omega$ , and randomly select initial interval,  $\omega_0$ , that would include the point,  $x_0$ , with possibility of expanding the interval. For example, it could be doubled to one side or expanded to other side until both ends of the interval will be outside of the slice.

In multidimensional case there are two different ways of sample generating.

The first one supposes application of one-dimensional approach to each variable from the multidimensional distribution. The second approach is based on application of slice sampling to generating multidimensional distribution by forming uniform sample from the area under the graph of PDF for this distribution.

The second approach was used in performing computational experiments because it is more natural and can be substantiated analogously to one-dimensional case. Here sample generating is performed from uniform distribution in the interval from zero to the value of density function at current point. The uniform value defined by slice is determined by the vertical dimension. It is obvious that uniform sample generating in a case of multidimensional slice is more difficult. In this case the interval, I = (L, R), is replaced by the hyper-interval:  $H = \left\{ x : L_i < x_i < R_i \quad \forall i = 1, ..., n \right\}.$ 

Here,  $L_i$  and,  $R_i$ , determine the length of the hyper-interval along the axis,  $x_i$ .

The simplest way of determining, H, is its placement randomly over dimensions in such a way that its value would be uniform over all possible values of, H, that contain initial point,  $x_0$ . Other procedures for determining the interval don't exhibit such simple generalization. For example, the procedure of expanding the interval until all boundaries will exceed the slice limits would not be effective because ndimensional interval has  $2^n$  peaks. That is why in the computational experiments the approach was used based on the interval expanding until the point taken uniformly from the interval belongs to the slice.

#### 7 Estimation of basic risk measures

The model has been constructed for analysis of exchange rates for the currencies: Swiss franc, British pound, Japanese yen, and US dollar to euro. To estimate model parameters the data for daily exchange rates were taken from 03.1998 to 01.2006. After preliminary data processing 1643 observations were selected for the model building. Using maximum likelihood technique parameters for one-dimensional marginal distributions estimated for the exchange rates selected, as for respective copula parameters (Table 1).

Copula	Parameter	Value	MSE
Gumbel	θ	1.6720	0.0158
Normal	$\rho_1$	0.5637	0.0118
	$\rho_2$	0.3318	0.0136
	ρ <sub>3</sub>	0.5943	0.0120
	$\rho_4$	0.8241	0.0054
	ρ <sub>5</sub>	0.8593	0.005
	ρ <sub>6</sub>	0.8037	0.0061
Frank	β	4.5874	0.0911

Table 1. Estimates for copula parameters

An empirical measure of VaR with quantile 0.03, i.e. for 50 observations that exceed the threshold selected amounted to 3.4967. For quantile 0.01, there are 16 observations that exceed the threshold 3.5345. There are enough observations for the quintile 0.03 to use in practice for calculating empirical estimate, but for quantile 0.01 the data sample is too short. That is why it is necessary to construct risk distribution model and estimate the VaR value using this model (Table 2).

Table 2 shows that VaR measures computed using the models on the basis of combined marginal distributions that were used to form joint distribution with Gumbel and Frank copulas exhibit relative error to the empirical value for quintile 0.03 of the following values: 0.203%, and 0.022%, respectively.

Copula	Quantile	Sample size		
		100	1000	10000
Gumbel	0.03	3.414	3.477	3.489
	0.01	3.437	3.566	3.601
Normal	0.03	3.500	3.538	3.534
	0.01	3.617	3.688	3.660
Frank	0.03	3.498	3.479	3.495
	0.01	3.513	3.563	3.589

Table 2. VaR estimates using the models

The model with normal copulas showed the error of about 1%. On the basis of the results obtained it can be concluded that all three models constructed are adequate but the model on the basis of Frank copula showed the best result. Thus, the VaR measure for, 0.01, quantile will be the value of 3.5892. The same three models were used to compute the coherent risk measures ES. An empirical value for measure ES with quantile 0.03 is 3.6074. Table 3 shows that the most adequate model for estimating the measure ES is also the model based on Frank copula: an estimate for ES with

<b>Table 3.</b> Estimates of risk measure ES				
Copula	Quintile	Sample size		
		100	1000	10000
Gumbel	0.03	3.512	3.575	3.583
	0.01	3.534	3.647	3.673
Normal	0.03	3.610	3.647	3.643
	0.01	3.712	3.773	3.752
Frank	0.03	3.585	3.566	3.583
	0.01	3.602	3.655	3.682

quintile 0.01 is 3.6822. All three models showed worse results for the risk deviation measure proposed by Markowitz than two tail measures considered above (Table 4).

Table 4. Estimates of	f Markowitz measure
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Estimation method	$\sigma_+$
Gumbel copula	0.1330
Normal copula	0.1462
Frank copula	0.1370
Empirical estimate	0.1733

Thus, quality of the models constructed shows that they can be used for active management of risks by changing the portfolio structure aiming to optimization of the risk measure selected.

# 8 Conclusion

The method for constructing joint distribution, copula based models, is proposed. The copula model parameters are estimated with the method of maximum likelihood which turned out to be effective according to the mean squared error criterion. Three types of copulas were studied and three models constructed. All three models turned out to be adequate and practically useful. Used approach to evaluate the risk measures based on sampling measures provided high precision of estimates when non-extreme quantiles were estimated. At the same time the quality of risk deviation measures that create a part of a risk profile requires model refinement in the future.

High quality of the tail measure estimates supports the idea that the model based upon combined marginal distributions, with the use of normal and generalized Pareto distribution, requires description improvement for the central part of observations. The future studies will take into consideration the possibility for applying to modeling central observations some other types of distribution than normal one. The purpose is to further improve quality of the model in the form of combined marginal distribution, and to refine final results of risk estimation.

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