

Martingale Characterizations Of Risk-Averse Stochastic Optimization Problems

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Abstract

This paper addresses risk awareness of stochastic optimization problems. Nested risk measures appear naturally in this context, as they allow beneficial reformulations for algorithmic treatments. The reformulations presented extend usual Hamilton–Jacobi–Bellman equations in dynamic optimization by involving risk awareness in the problem formulation.

Nested risk measures are built on risk measures, which originate by conditioning on the history of a stochastic process. We derive martingale properties of these risk measures and use them to prove continuity. It is demonstrated that stochastic optimization problems, which incorporate risk awareness via nesting risk measures, are continuous with respect to the natural distance governing these optimization problems, the nested distance.

Keywords: Risk measures, Stochastic optimization, Stochastic processes

Classification: 90C15, 60B05, 62P05

1 Introduction

Risk measures have been found useful in various disciplines of applied mathematics, particularly in mathematical finance and in stochastic optimization. Many applications involve them in various places to account for risk. It is hence natural to investigate risk measures in a multistage or dynamic optimization framework as well. One of the first occurrences of dynamic risk measures in the literature is Riedel [23], conditional risk measures are discussed in Ruszczyński and Shapiro [27] (consider also the references therein).

It seems that there is no general consensus on how to incorporate risk measures in a more general framework which involves time. One of the conceptual difficulties arising in a problem setting involving time is time consistency. In short, the decisions considered optimal at some stage of time should not be rejected from a later perspective.

Risk-averse multistage stochastic programs incorporate risk awareness in multistage decision making. These problems have been considered in Ruszczyński [26] and Dentcheva and Ruszczyński [6], while applications can be found in Philpott and de Matos [19], Philpott et al. [20] or Maggioni et al. [14], e.g., where stochastic dual dynamic programming methods are addressed, cf. also Römisch and Guigues [25], Girardeau et al. [9]. In economics, the spread between *risk-averse* and *risk-neutral* preferences is associated with a risk or insurance premium. For this, the prevailing idea of risk in these papers is the interpretation as insurance on a rolling horizon basis.

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This paper introduces conditional risk functionals based on the history of the governing stochastic process. These functionals are nested to obtain risk functionals accounting for the risk at each stage of the stochastic process. We elaborate their continuity properties and for important cases we compare them with simple risk measures spanning the entire horizon as a whole.

Building on the idea in Pflug [16] we introduce the nested distance via conditional probabilities. We relate these concepts by verifying that nested risk functionals are continuous with respect to the nested distance and provide an explicit expression of the modulus of continuity.

Martingales are present in stochastic optimization since its very beginning, cf. Rockafellar and Wets [24]. The approach taken here to verify the results is based on generalized martingales. They reflect the evolution of risk over time, as risk measures replace risk-neutral expectations. It is demonstrated that the nested distance, as well as nested risk measures, follow martingale characteristics in this generalized sense.

It is a consequence that risk-averse multistage stochastic programs are continuous with respect to the nested distance. The optimal solutions constitute a stochastic process, which again follows a martingale-like pattern. We finally give a verification theorem. This is a risk-averse generalization of Hamilton–Jacobi–Bellman equations, which are well-known from dynamic optimization.

Outline of the paper. Section 3 introduces nested risk functionals after an introductory discussion (Section 2). Section 4 addresses the main features of the nested distance which are important and relevant to cover the discussion on continuity of the multistage stochastic programs in Section 5. Risk martingales are introduced in Section 6. We conclude with the main result in Section 7.

2 Notation and preliminaries

We consider the Polish spaces (Ξ_t, d_t) , $t \in \{1, \dots, T\}$. We shall associate $t \in \{1, \dots, T\}$ with *stage* or *time* advancing in discrete steps from 1 to T , where $T \in \{1, 2, \dots\}$ is the time horizon (terminal time) or final stage. Each space Ξ_t , $t = 1, \dots, T$, contains the information revealed at time t . In what follows it will often be sufficient to consider the spaces $\Xi_t = \mathbb{R}^{m_t}$.

The product $\Xi := \Xi_{1:T} := \Xi_1 \times \dots \times \Xi_T$ is endowed with the metric d and

$$(\Xi, d) \tag{1}$$

is Polish as well (for example, choose $d(x, y) := \ell_p(x, y) := \left(\sum_{t=1}^T d_t(x_t, y_t)^p \right)^{1/p}$). We denote elements $x \in \Xi_{1:T}$ by $x_{1:T} := x = (x_1, \dots, x_T)$ and by pr_t the canonical (i.e., coordinate) projection $\text{pr}_t(x_{1:T}) := x_{1:t}$ onto the subspace $\Xi_{1:t} := \Xi_1 \times \dots \times \Xi_t$. To allow a compact notation we also introduce the empty tuple $x_{1:0} = ()$.

On the Borel sets $\mathcal{F}_T := \mathcal{B}(\Xi_{1:T})$ we consider the probability measure

$$P: \mathcal{F}_T \rightarrow [0, 1].$$

The probability measures restricted to the sub-sigma algebra $\mathcal{F}_t := \sigma(\text{pr}_t)$ are the image measures defined by

$$P_t(A) := P^{\text{pr}_t}(A) = P(A \times \Xi_{t+1} \times \dots \times \Xi_T),$$

where $A \in \mathcal{B}(\Xi_{1:t})$, the Borel sigma algebra on $\Xi_{1:t}$. The sequence $\mathcal{F} := \mathcal{F}_{0:T} := (\mathcal{F}_t)_{t=0}^T$ is the canonical (i.e., coordinate) filtration and $(\Xi_{1:T}, \mathcal{F}_{0:T}, P)$ is a filtered probability space (a.k.a. stochastic basis), where we include the trivial sigma algebra $\mathcal{F}_0 := \{\emptyset, \Xi_{1:T}\}$ for completeness and convenience.

The disintegration theorem (cf. Dellacherie and Meyer [5, III-70] or Ambrosio et al. [1, Section 5.3]) allows ‘disintegrating’ the probability measure with respect to the coordinates.

Theorem 1 (Disintegration theorem). *There is a regular kernel, i.e., a P_t -a.s. uniquely defined family of measures $P(\cdot | x_{1:t})$ so that*

- (i) $x_{1:t} \mapsto P(B | x_{1:t})$ is measurable for every $B \in \mathcal{B}(\Xi_{t+1} \times \dots \times \Xi_T)$ and
- (ii) $P(A \times B) = \int_A P(B | x_{1:t}) P_t(dx_{1:t})$, where $A \in \mathcal{B}(\Xi_1 \times \dots \times \Xi_t)$ and $B \in \mathcal{B}(\Xi_{t+1} \times \dots \times \Xi_T)$.

The conditional probability measures

$$P_{t+1}(\cdot | x_{1:t}) \text{ on } \mathcal{B}(\Xi_{t+1}) \quad (2)$$

are called (regular) kernels and the substring $x_{1:t}$ is also called a fiber.

By disintegrating the measures P_t and composing their kernels at subsequent stages we obtain the nested expressions

$$P_t(A_1 \times \dots \times A_t) = \int_{A_1} \int_{A_2} \dots \int_{A_t} P_t(dx_t | x_{1:t-1}) \dots P_2(dx_2 | x_{1:1}) P_1(dx_1) \quad (3)$$

and the conditional probability measures

$$P(A_{t+1} \times \dots \times A_T | x_{1:t}) = \int_{A_{t+1}} \dots \int_{A_T} P_T(dx_T | x_{1:T-1}) \dots P_{t+1}(dx_{t+1} | x_{1:t}). \quad (4)$$

Both expressions reveal the initial probability measure P , which can be seen by substituting $t = T$ in (3) or $t = 0$ in (4).

Remark 2. The kernels derived from the projected measures (2) are conditioned on the history $x_{1:t}$ and they do depend explicitly on the entire history up to t . In the Markovian case this dependence reduces (simplifies) to

$$P_{t+1}(\cdot | x_{1:t}) = P_{t+1}(\cdot | x_t).$$

An important algorithm in stochastic optimization is Stochastic Dual Dynamic Programming (SDDP). In this context the probabilities are typically assumed to be *stagewise independent*, i.e.,

$$P_{t+1}(\cdot | x_{1:t}) = P_{t+1}(\cdot)$$

(cf. Goulart and da Costa [10]).

3 Conditional and nested risk measures

To define conditional risk functionals we recall the definition of *law invariant, coherent risk functionals* $\mathcal{R}: L \rightarrow \mathbb{R}$ defined on some vector space L of \mathbb{R} -valued random variables first. They satisfy the following axioms introduced by Artzner et al. [2].

- A1 Monotonicity: $\mathcal{R}(Y_0) \leq \mathcal{R}(Y_1)$, provided that $Y_0 \leq Y_1$ almost surely;
- A2 Translation equivariance: $\mathcal{R}(Y + c) = \mathcal{R}(Y) + c$ for $c \in \mathbb{R}$;
- A3 Convexity: $\mathcal{R}((1 - \lambda)Y_0 + \lambda Y_1) \leq (1 - \lambda)\mathcal{R}(Y_0) + \lambda\mathcal{R}(Y_1)$;
- A4 Positively homogeneity: $\mathcal{R}(\lambda Y) = \lambda\mathcal{R}(Y)$;
- A5 Law invariance: $\mathcal{R}(Y) = \mathcal{R}(Y')$, whenever Y and Y' have the same law, i.e., $P(Y \leq y) = P(Y' \leq y)$ for all $y \in \mathbb{R}$.

We shall make frequently use of the following proposition, which is an immediate consequence of the monotonicity axiom [A1](#).

Proposition 3. *The essential infimum of a set of random variables apparently satisfies $\text{ess inf}_{\iota \in I} Y_\iota \leq Y_\iota$ for every $\iota \in I$. Hence, by the monotonicity axiom, [A1](#), $\mathcal{R}(\text{ess inf}_{\iota \in I} Y_\iota) \leq \mathcal{R}(Y_\iota)$ and subsequently*

$$\mathcal{R}\left(\text{ess inf}_{\iota \in I} Y_\iota\right) \leq \inf_{\iota \in I} \mathcal{R}(Y_\iota).$$

The Average Value-at-Risk at level $\alpha \in [0, 1)$ defined on $L^1(P)$ by

$$\text{AV@R}_\alpha(Y) := \inf_{q \in \mathbb{R}} \left\{ q + \frac{1}{1-\alpha} \mathbb{E}(Y - q)_+ \right\} \quad (5)$$

is the most prominent coherent risk functional satisfying the axioms [A1–A5](#) above. The Average Value-at-Risk at risk level $\alpha = 0$ is the expectation,

$$\text{AV@R}_0(Y) = \mathbb{E} Y$$

and, for $Y \in L^\infty$, the convenient setting

$$\text{AV@R}_1(Y) := \lim_{\alpha \nearrow 1} \text{AV@R}_\alpha(Y) = \text{ess sup } Y$$

continuously extends the Average Value-at-Risk to $\alpha = 1$.

The Average Value-at-Risk turns out to be of central importance, it can be interpreted as an extreme point in the set of risk functionals and, similarly to Choquet's representation, every risk functional is a convex combination of AV@Rs. The following general representation (Kusuoka's representation, cf. Kusuoka [[13](#)]) highlights this relation. The statement is a consequence of the Fenchel–Moreau theorem in convex analysis (cf. Föllmer and Schied [[8](#), Lemma 4.55] or Shapiro et al. [[31](#)], Shapiro [[28](#)], Pichler and Shapiro [[22](#)]).

Definition 4. A function $\sigma: [0, 1) \rightarrow \mathbb{R}$ is a *distortion function*, if $\sigma(\cdot)$ is non-decreasing, $\sigma(\cdot) \geq 0$ and $\int_0^1 \sigma(u) du = 1$.

Proposition 5 (Derived from Kusuoka's representation, cf. Pflug and Pichler [[17](#)]). *Every law invariant, coherent risk functional $\mathcal{R}: L \rightarrow \mathbb{R}$ has the representation*

$$\mathcal{R}(Y) = \sup_{\sigma \in \mathcal{S}} \mathcal{R}_\sigma(Y), \quad (6)$$

where \mathcal{S} is an appropriate collection of distortion functions and

$$\mathcal{R}_\sigma(Y) := \sup \left\{ \mathbb{E} Y \zeta \mid \begin{array}{l} \zeta \geq 0, \mathbb{E} \zeta = 1 \text{ and} \\ \text{AV@R}_\alpha(\zeta) \leq \frac{1}{1-\alpha} \int_\alpha^1 \sigma(u) du \text{ for all } \alpha \in (0, 1) \end{array} \right\}. \quad (7)$$

The vector space L can be assumed to be $L = \{Y: \mathcal{R}_\sigma(|Y|) < \infty\}$ (cf. Pichler [[21](#)]). In applications, as well in what follows it will be enough to consider the Lebesgue spaces L^p , $p \geq 1$, or L^∞ .

The representation of the distortion risk functional (6) implicitly involves the probability measure P via the expectation \mathbb{E} and the Average Value-at-Risk in (7). We want to make the probability measure P explicit by rewriting (6) as

$$\mathcal{R}(Y) = \mathcal{R}_{\mathcal{S}; P}(Y) := \sup \left\{ \mathbb{E}_P Y \zeta \mid \begin{array}{l} \zeta \geq 0, \mathbb{E}_P \zeta = 1 \text{ and} \\ \text{AV@R}_{\alpha; P}(\zeta) \leq \frac{1}{1-\alpha} \int_\alpha^1 \sigma(u) du, \alpha \in (0, 1) \\ \text{for some } \sigma(\cdot) \in \mathcal{S} \end{array} \right\}, \quad (8)$$

where the expectation in $\text{AV@R}_{\alpha; P}$ is with respect to the probability measure P as well, cf. (5).

Example 6. The Kusuoka representation of the Average Value-at-Risk according to Proposition 5 is given by $\mathcal{S} = \{\sigma_\alpha(\cdot)\}$, where the distortion function is $\sigma_\alpha(u) := \begin{cases} \frac{1}{1-\alpha} & \text{if } u \geq \alpha, \\ 0 & \text{else.} \end{cases}$

3.1 Conditional risk measures

To define conditional versions of risk measures on product spaces we employ the conditional measures available by the disintegration theorem, Theorem 1.

Definition 7. Let \mathcal{S}_{t+1} be a collection of distortion functions. The *conditional risk measure* or *risk measure conditioned on the fiber* $x_{1:t}$ of the regular kernels of the probability measure P is

$$\mathcal{R}_{\mathcal{S}_{t+1}}(Y \mid x_{1:t}) := \sup_{\sigma \in \mathcal{S}_{t+1}} \mathcal{R}_{\sigma; P(\cdot \mid x_{1:t})}(Y). \quad (9)$$

As a consequence of Theorem 1(i) and the representations (6) and (7), the mapping

$$\begin{aligned} \mathcal{R}_{\mathcal{S}_{t+1}}(Y \mid \cdot): \Xi_{1:t} &\rightarrow \mathbb{R} \\ x_{1:t} &\mapsto \mathcal{R}_{\mathcal{S}_{t+1}}(Y \mid x_{1:t}) \end{aligned} \quad (10)$$

is a random variable on $\Xi_{1:t}$, which is P_t a.s. well-defined and measurable with respect to \mathcal{F}_t . For $t = 0$, the conditional risk functional (9) is

$$\mathcal{R}_{\mathcal{S}_1}(Y \mid x_{1:0}) = \mathcal{R}_{\mathcal{S}_1}(Y) = \sup_{\sigma \in \mathcal{S}_1} \mathcal{R}_{\sigma; P}(Y) = \mathcal{R}_{\mathcal{S}_1}(Y),$$

a deterministic number.

3.2 Nested risk measures

The conditional risk measures (9) are well-defined on a fiber $x_{1:t}$. As each risk functional (10) is a random variable, they can be combined and considered in the following recursive, or nested way.

Definition 8 (Nested risk functional). Let $s, t \in \{1, \dots, T\}$ with $s < t$. The *nested risk functional* for a sequence $\mathcal{S}_{s+1:t} := \mathcal{S}_{s+1} \times \dots \times \mathcal{S}_t$ of collections of distortion functionals is

$$\mathcal{R}_{\mathcal{S}_{s+1:t}}(Y \mid x_{1:s}) := \mathcal{R}_{\mathcal{S}_{s+1}}\left(\dots \mathcal{R}_{\mathcal{S}_{t-1}}\left(\mathcal{R}_{\mathcal{S}_t}(Y \mid x_{1:t-1}) \mid x_{1:t-2}\right) \dots \mid x_{1:s}\right). \quad (11)$$

Remark 9. The nested risk functional $\mathcal{R}_{\mathcal{S}_{1:T}}(\cdot)$ maps real-valued random variables $Y: \Xi \rightarrow \mathbb{R}$ defined on Ξ to the real line. The nested risk functional satisfies generalizations of the axioms A1–A4, but it is *not* law invariant any longer, i.e., A5 is not necessarily satisfied.

The construction employed in Shapiro [29] to discuss rectangular sets is similar to nested risk measure given in Definition 8 above. Indeed, they can be recovered by choosing the feasible set as given in the general representation (8). A major difference is given by the fact that law invariant risk functionals have the Kusuoka representation (8), which is not the case for more general risk functionals.

Importantly, the nested risk measures are recursive as specified in the following proposition.

Proposition 10. *The nested risk functional $\mathcal{R}_{\mathcal{S}_{t+1:T}}$ is recursive, it holds that*

$$\mathcal{R}_{\mathcal{S}_{t+1:T}}(Y \mid x_{1:t}) = \mathcal{R}_{\mathcal{S}_{t+1:s}}(\mathcal{R}_{\mathcal{S}_{s+1:T}}(Y \mid x_{1:s}) \mid x_{1:t}) \quad (12)$$

whenever $0 \leq t < s < T$.

Proof. The assertion is an immediate consequence of the recursion (11) in Definition 8. \square

Example 11 (Conditional expectation). The risk-neutral special case is given by choosing the simplest distortion functions $\mathcal{S}_{t+1} = \{\mathbb{1}\}$, i.e., the distortions consisting only of the constant function $\sigma(\cdot) = \mathbb{1}(\cdot) = 1$. In this case the risk functional (9) is

$$\mathcal{R}_{\mathcal{S}_{t+1};P}(Y \mid x_{1:t}) = \mathbb{E}(Y \mid x_{1:t}),$$

i.e.,

$$\mathcal{R}_{\mathcal{S}_{t+1};P}(Y \mid \cdot) = \mathbb{E}^{\mathcal{F}_t}(Y)$$

(recall that $\mathbb{E}^{\mathcal{F}_t}(Y)$ is indeed an \mathcal{F}_t random variable). The recursion (12) reflects the tower property of the conditional expectation.

Definition 12 (Nested Average Value-at-Risk, cf. Pflug and Römisch [18]). The nested Average Value-at-Risk for $\alpha_{s+1:t} \in [0, 1]^{t-s}$ is a composition of AV@Rs at risk levels dependent on the state t . More explicitly, we set

$$\text{nAV@R}_{\alpha_{s+1:t}}(Y \mid x_{1:s}) := \text{AV@R}_{\alpha_{s+1};P(\cdot \mid x_{1:s})} \left(\dots \text{AV@R}_{\alpha_{t-1};P(\cdot \mid x_{1:t-2})} \left(\text{AV@R}_{\alpha_t;P(\cdot \mid x_{1:t-1})}(Y) \right) \right). \quad (13)$$

The nested Average Value-at-Risk can be bounded by the Average Value-at-Risk. Indeed, it follows from Xin and Shapiro [34, Proposition 4.2] that $\text{nAV@R}_{\alpha_{1:T};P}(Y) \leq \text{AV@R}_{\alpha}(Y)$ provided that the risk level α satisfies $\alpha \geq 1 - (1 - \alpha_1) \dots (1 - \alpha_T)$.

4 The distance adapted to nested risk measures

Generalizing the concept of distance from probability spaces to filtered probability spaces corresponds to generalizing the distance from random variables to stochastic processes. As a metric for probability measures we recall the Wasserstein distance first here, which we then generalize to a metric of stochastic processes.

4.1 Wasserstein metric

Consider the Polish space (Ξ, d) and probability measures

$$P, \tilde{P}: \mathcal{F} \rightarrow [0, 1]$$

on the Borel sigma algebra $\mathcal{F} := \mathcal{B}(\Xi)$.

Definition 13 (Wasserstein metric). Let P and \tilde{P} be probability measures on Ξ and $r \in [1, \infty)$. The Wasserstein metric of order r with respect to the cost function $c: \Xi \times \Xi \rightarrow \mathbb{R}$ is

$$w_r(P, \tilde{P}; c) := \inf_{\pi} (\mathbb{E}_{\pi} c^r)^{1/r} = \inf_{\pi} \left(\iint_{\Xi \times \Xi} c(x, y)^r \pi(\mathrm{d}x, \mathrm{d}y) \right)^{1/r}, \quad (14)$$

where the infimum in (14) is among all bivariate probability measures $\pi \in \mathcal{P}(\Xi \times \Xi)$ with marginals P and \tilde{P} , i.e.,

$$\pi(A \times \Xi) = P(A), \quad A \in \mathcal{B}(\Xi) \quad \text{and} \quad (15)$$

$$\pi(\Xi \times B) = \tilde{P}(B), \quad B \in \mathcal{B}(\Xi). \quad (16)$$

For the Wasserstein distance of order $r = 1$ we shall also write simply $w(P, \tilde{P})$.

Remark 14. The Wasserstein metric introduced in (14) is based on a cost functions $c(\cdot)$ (cf. also Villani [33]). This setting slightly generalizes the usual definition, which is based on the distance function d of the space (Ξ, d) in lieu of c . In what follows, this extension will be essential.

4.2 The nested distance

The Wasserstein metric w_r introduced in Definition 13 is of course well defined for measures P and \tilde{P} on the product space $(\Xi_{1:T}, d)$. The nested distance generalizes the Wasserstein metric by involving the filtration in addition. The filtration carries the information revealed over time. The filtration considered here is the coordinate filtration, and for this we may introduce the nested distance on coordinate basis as well, i.e., sequentially by defining the process stage by stage.

Definition 15 (Cost process, nested distance). Let P and \tilde{P} be probability measures on $\Xi_{1:T}$, let $r \in [1, \infty)$ and let $c: \Xi_{1:T} \times \Xi_{1:T} \rightarrow \mathbb{R}$ be a lower semi-continuous (lsc.) function.

(i) Cost process c_t for $t = T$ down to 0:

(a) The cost function c_T on $\Xi_{1:T} \times \Xi_{1:T}$ at terminal time T is

$$c_T(x_{1:T}, y_{1:T}) := c(x_{1:T}, y_{1:T}).$$

We shall refer to c_T also as the *terminal* cost function.

(b) The cost functions c_t for $t < T$ are defined in a backwards recursive way by

$$c_{t-1}(x_{1:T}, y_{1:T}) := w_r(P_t(\cdot \mid x_{1:t-1}), \tilde{P}_t(\cdot \mid y_{1:t-1}); c_t), \quad t = T, \dots, 1, \quad (17)$$

where w_r is the Wasserstein metric of order r .

(c) The *cost-process* is the stochastic process $c = (c_t)_{t=0}^T$.

(ii) The nested distance: let $c = (c_t)_{t=0}^T$ be the cost process with terminal cost

$$c_T(\cdot) = d(\cdot), \quad (18)$$

the distance of the space $\Xi_{1:T}$ (cf. (1)). The *nested distance* of order $r \geq 1$ of the measures P and \tilde{P} is

$$\text{dl}_r(P, \tilde{P}) := c_0. \quad (19)$$

Remark 16. The function c_t is defined for $(x_{1:T}, y_{1:T}) \in \Xi_{1:T} \times \Xi_{1:T}$, but its definition in (17) notably involves only the truncated states $(x_{1:t}, y_{1:t}) \in \Xi_{1:t} \times \Xi_{1:t}$. The cost function c_t thus is unambiguously defined for $(x_{1:t}, y_{1:t})$, irrespective of future realization $(x_{t+1:T}, y_{t+1:T})$. It follows that c_t is $\mathcal{F}_t \otimes \mathcal{F}_t$ measurable and the cost process $(c_t)_{t=0}^T$ is adapted to the filtration $\mathcal{F} \otimes \mathcal{F}$.

In particular, c_0 is independent of the formal argument $(x_{1:T}, y_{1:T})$ (the string $x_{1:0}$ is empty for $t = 0$ in (17)) so that c_0 is a number ($c_0 = \text{dl}_r(P, \tilde{P}) \in \mathbb{R}$) and the nested distance is well-defined by (19).

Remark 17. It is a consequence of Hölder's inequality that $w_r(P, \tilde{P}) \leq w_{r'}(P, \tilde{P})$ whenever $r \leq r'$. By monotonicity of (17) we thus get that

$$dl_r(P, \tilde{P}) \leq dl_{r'}(P, \tilde{P}) \quad (r \leq r'). \quad (20)$$

Remark 18 (Relation to Wasserstein metric). For $T = 1$ we have $\Xi_{1:T} = \Xi_1$ and there are no intermediary stages present. In this case, the nested distance reduces to the usual Wasserstein metric and it holds that

$$dl_r(P, \tilde{P}) = w_r(P, \tilde{P}; d) \quad (T = 1).$$

Remark 19. As for the Wasserstein distance we also write $dl(P, \tilde{P})$ if the order is $r = 1$ (cf. Remark 14).

An important case in practice is the cost functions, where costs occur sequentially at every stage and total costs are accumulated over time. The cost process reflects this additive property, as the following proposition outlines.

Proposition 20 (Additive cost functions). *Suppose the terminal cost function is of particular form*

$$c_T(x, y) = \ell_r(x, y) = \left(\sum_{t=1}^T d_t(x_t, y_t)^r \right)^{1/r}, \quad (21)$$

where d_t , $t = 1, \dots, T$ are functions on $\Xi_t \times \Xi_t$ (distance functions, e.g.). Then the process

$$\tilde{c}_t := \left(c_t^r - \sum_{j=1}^{t-1} d_j^r \right)^{1/r} \quad (22)$$

satisfies the recursive equations

$$\tilde{c}_{t-1}^r = d_{t-1}^r + w_r(P_t(\cdot | x_{1:t-1}), \tilde{P}_t(\cdot | y_{1:t-1}); \tilde{c}_t)^r \quad (23)$$

with $\tilde{c}_T = d_T$.

Further, the nested distance is

$$dl_r(P, \tilde{P}) = \tilde{c}_0.$$

Remark 21. The recursive equation (23) is actually the initial attempt in defining a distance on the nested spaces $\Xi_t \times \mathcal{P}(\Xi_{t-1})$ for the particular case $r = 1$, where $\mathcal{P}(\Xi_{t-1})$ is the set of probability measures on Ξ_{t-1} . We refer to Pflug [16] for the initial and complete discussion on nested spaces and nested distances.

Proof. From (17) we have that

$$c_{t-1}(x_{1:T}, y_{1:T})^r = w_r(P_t(\cdot | x_{1:t-1}), \tilde{P}_t(\cdot | y_{1:t-1}); c_t)^r.$$

As d_j are \mathcal{F}_{t-1} -measurable for every $j < t$ it follows further that

$$c_{t-1}(x_{1:T}, y_{1:T})^r = \sum_{j=1}^{t-1} d_j^r + w_r \left(P_t(\cdot | x_{1:t-1}), \tilde{P}_t(\cdot | y_{1:t-1}); \left(c_t^r - \sum_{j=1}^{t-1} d_j^r \right)^{1/r} \right)^r$$

and hence

$$\begin{aligned}
\tilde{c}_{t-1}^r(x_{1:T}, y_{1:T})^r &= c_{t-1}(x_{1:T}, y_{1:T})^r - \sum_{j=1}^{t-2} d_j^r \\
&= d_{t-1}^r + w_r \left(P_t(\cdot | x_{1:t-1}), \tilde{P}_t(\cdot | y_{1:t-1}); \left(c_t^r - \sum_{j=1}^{t-1} d_j^r \right)^{1/r} \right) \\
&= d_{t-1}^r + w_r (P_t(\cdot | x_{1:t-1}), \tilde{P}_t(\cdot | y_{1:t-1}); \tilde{c}_t)^r,
\end{aligned}$$

which is the assertion. \square

Remark 22. It is evident that the assertion of the previous statement holds as well in case of cost functions which are nonanticipative and of the form $c_T(x, y) = \left(\sum_{t=1}^T d_t(x_{1:t}, y_{1:t})^r \right)^{1/r}$.

4.3 Characterization as a martingale

For the measure P we have given the nested expressions (3) and (4) based on kernels explicitly. In the same way one may glue together the kernels which are optimal in (17) to compute the nested distance and cost process. To this end denote the optimal kernels on $\Xi_t \times \Xi_t$ obtained in (17) by $\pi_t(\cdot \times \cdot | x_{1:t}, y_{1:t})$. A well-known result of Brenier [3, 4] (see also McCann [15]) asserts that the Wasserstein problem (14) attains the infimum at a unique bivariate measure π for the quadratic cost function $c(x, y) = \|x - y\|^2$, if both measures P and \tilde{P} have finite variance and do not give mass to small sets (cf. Villani [33, Theorem 2.12]); the measures $\pi_t(\cdot \times \cdot | x_{1:t}, y_{1:t})$ thus exist.

The global measure governing all kernels then is

$$\begin{aligned}
\pi(A \times B) := & \iint_{A_1 \times B_1} \left(\iint_{A_2 \times B_2} \dots \left(\iint_{A_T \times B_T} \pi_T(dx_T, dy_T | x_{1:T-1}, y_{1:T-1}) \right) \right. \\
& \left. \dots \pi_2(dx_2, dy_2 | x_1, y_1) \right) \pi_1(dx_1, dy_1),
\end{aligned} \tag{24}$$

where $A = A_1 \times \dots \times A_T$ and $B = B_1 \times \dots \times B_T$. The measure π is a bivariate measure on the entire space $\Xi_{1:T} \times \Xi_{1:T}$.

We have the following alternative characterization of the governing bivariate measure (24).

Proposition 23. *The conditional marginals of the measure π defined in (24) satisfy*

$$\pi(A \times \Xi | x_{1:t}, y_{1:t}) = P(A | x_{1:t}), \quad A \in \mathcal{F}_T \quad \text{and} \tag{25}$$

$$\pi(\Xi \times B | x_{1:t}, y_{1:t}) = \tilde{P}(B | y_{1:t}), \quad B \in \mathcal{F}_T, \tag{26}$$

for every $t \in \{0, \dots, T-1\}$.

Proof. The most inner integral in (24) satisfies

$$\iint_{A_T \times \Xi_T} \pi(dx_T, dy_T | x_{1:T-1}, y_{1:T-1}) = \pi(A_T \times \Xi_T | x_{1:T-1}, y_{1:T-1}) = P(A_T | x_{1:T-1})$$

by construction of the measure $\pi(\cdot, \cdot | x_{1:T-1}, y_{1:T-1})$. This is (25) for the terminal time $t = T-1$.

Suppose now, by backwards inductions, that the marginal (25) is valid for $t + 1$. Then

$$\begin{aligned}
& \pi(A_{t+1:T} \times \Xi_{t+1:T} | x_{1:t}, y_{1:t}) \\
&= \iint_{A_{t+1} \times \Xi_{t+1}} \cdots \iint_{A_T \times \Xi_T} \pi(dx_T, dy_T | x_{1:T-1}, y_{1:T-1}) \cdots \pi(dx_{t+1}, dy_{t+1} | x_{1:t}, y_{1:t}) \\
&= \iint_{A_{t+1} \times \Xi_{t+1}} \pi(A_{t+2:T} \times \Xi_{t+2:T} | x_{1:t+1}, y_{1:t+1}) \pi(dx_{t+1}, dy_{t+1} | x_{1:t}, y_{1:t}) \\
&= \iint_{A_{t+1} \times \Xi_{t+1}} P(A_{t+2:T} | x_{1:t+1}) \pi(dx_{t+1}, dy_{t+1} | x_{1:t}, y_{1:t}) \\
&= \int_{A_{t+1}} P(A_{t+2:T} | x_{1:t+1}) P(dx_{t+1} | x_{1:t}) \\
&= P(A_{t+1:T} | x_{1:t}),
\end{aligned}$$

where we have used the decomposition (24), the induction hypothesis, the decomposition (4) and the setting $A_{t+1:T} := A_{t+1} \times A_{t+2:T}$. We conclude that identity (25) is valid for all t .

The remaining identity (26) follows analogously. \square

The process $(c_t)_{t=0}^T$ given in Definition 15 is constructed by recursively averaging with respect to the conditional measures of π given in (24). We thus have the following characterization as a martingale.

Theorem 24 (Martingale characterization). *Let $\pi(\cdot, \cdot)$ be the measure defined in (24) and $r \geq 1$. Then the cost process $c = (c_t)_{t=1}^T$ is a martingale with respect to π and the canonical filtration, i.e.,*

$$c_t^r = \mathbb{E}_\pi(c_{t+1}^r | \mathcal{F}_t \otimes \mathcal{F}_t).$$

Proof. By definition of the process c_t in (17) we have that

$$c_{t-1}(x_{1:T}, y_{1:T})^r = \iint_{\Xi_t \times \Xi_t} c_t(x_{1:T}, y_{1:T})^r \pi(dx_t, dy_t | x_{1:t-1}, y_{1:t-1}),$$

where $\pi(\cdot, \cdot | x_{1:t-1}, y_{1:t-1})$ is the measure with marginals $P(\cdot | x_{1:t-1})$ and $\tilde{P}(\cdot | y_{1:t-1})$, resp., for which the Wasserstein distance attains the infimum in (17). This is the conditional martingale property for the fibers $(x_{1:t-1}, y_{1:t-1})$. The assertion follows as the measure π in (24) combines these optimal, conditional measures. \square

Corollary 25 (Alternative characterization). *The nested distance is given by*

$$dl_r(P, \tilde{P}) = \inf_{\pi} (\mathbb{E}_{\pi} d^r)^{1/r} = \inf_{\pi} \left(\iint_{\Xi \times \Xi} d(x, y)^r \pi(dx, dy) \right)^{1/r},$$

where the infimum is among all probability measures $\pi \in \mathcal{P}(\Xi \times \Xi)$ satisfying the conditional marginal constraints (25)–(26). The infimum is attained for the measure π defined in (24).

Proof. Let $\pi(\cdot | \cdot)$ satisfy the marginals (25)–(26). Then every conditional measure $\pi(\cdot, \cdot | x_{1:t-1}, y_{1:t-1})$ satisfies the constraints (15)–(16) to compute the Wasserstein distance. It follows that $dl_r(P, \tilde{P})^r \leq \mathbb{E}_{\pi} d^r$.

The measure π defined in (24) satisfies the constraints (25)–(26) as well. However, we have from Theorem 24 that c_t^r is a martingale. The assertion follows from the power property of the conditional expectation, as $c_T^r = d^r$ and

$$\begin{aligned} \text{dl}_r(P, \tilde{P})^r &= c_0^r = \mathbb{E}_\pi (\dots \mathbb{E}_\pi (c_{t+1}^r | \mathcal{F}_t \otimes \mathcal{F}_t) \cdots | \mathcal{F}_1 \otimes \mathcal{F}_1) \\ &= \mathbb{E}_\pi (\dots \mathbb{E}_\pi (\dots \mathbb{E}_\pi (d^r | \mathcal{F}_T \otimes \mathcal{F}_T) \cdots | \mathcal{F}_t \otimes \mathcal{F}_t) \cdots | \mathcal{F}_1 \otimes \mathcal{F}_1) \\ &= \mathbb{E}_\pi d^r; \end{aligned}$$

hence the result. \square

For additive cost functions the distance of the individual stages have to be taken care of. The following corollary describes the process in analogy to Proposition 20 above.

Corollary 26 (Additive cost functions). *Let $\pi(\cdot, \cdot)$ be the optimal measure (24) and c_T the additive cost function (21) for $r \geq 1$. Then the process*

$$\tilde{c}_t^r + \sum_{j=1}^{t-1} d_j^r$$

is a martingale with respect to the measure π (cf. (22)).

Proof. This is immediate as $\tilde{c}_t^r = c_t^r - \sum_{j=1}^{t-1} d_j^r$ by definition of the process (22) and as c_t is a martingale by Theorem 24. \square

5 Continuity properties

The risk functionals defined in (6) above are continuous with respect to the Wasserstein distance. We generalize the results here and verify that nested risk functionals are continuous with respect to the nested distance. This section elaborates the modulus of continuity.

Proposition 27 (Continuity of risk functionals). *Let \mathcal{R}_S be a general risk functional according (8). Suppose that the random variables $Y, \tilde{Y}: \Xi \rightarrow \mathbb{R}$ satisfy*

$$Y(x) - \tilde{Y}(y) \leq L \cdot d(x, y)^\beta \tag{27}$$

for some $\beta \leq 1$. Then

$$\begin{aligned} \mathcal{R}_{S;P}(Y) - \mathcal{R}_{S;\tilde{P}}(\tilde{Y}) &\leq L \cdot \sup_{\sigma \in S} \|\sigma\|_q \cdot w_{\beta r}(P, \tilde{P})^\beta \\ &\leq L \cdot \sup_{\sigma \in S} \|\sigma\|_q \cdot w_r(P, \tilde{P})^\beta, \end{aligned}$$

where $q \in (1, \infty]$ is the Hölder conjugate exponent of r (the order of the Wasserstein metric) for which $\frac{1}{q} + \frac{1}{r} = 1$.

Proof. Let $\zeta \geq 0$ with $\mathbb{E} \zeta = 1$ be chosen so that the supremum in (8) is attained up to $\varepsilon > 0$, i.e., $\mathbb{E} Y \zeta > \mathcal{R}_{S;P}(Y) - \varepsilon$. Let π have marginals P and \tilde{P} . Note that $\mathbb{E}_\pi \zeta = \mathbb{E}_P \zeta = 1$, so that

$$\mathcal{R}_{S;\tilde{P}}(\tilde{Y}) = \mathcal{R}_{S;\pi}(\tilde{Y}) \geq \mathbb{E}_\pi \tilde{Y} \zeta.$$

It follows from Hölder's inequality that

$$\begin{aligned}
\mathcal{R}_{S;P}(Y) - \varepsilon - \mathcal{R}_{S;\tilde{P}}(\tilde{Y}) &\leq \iint_{\Xi \times \Xi} (Y(x) - \tilde{Y}(y)) \zeta(x) \pi(\mathrm{d}x, \mathrm{d}y) \\
&\leq L \iint_{\Xi \times \Xi} d(x, y)^\beta \zeta(x) \pi(\mathrm{d}x, \mathrm{d}y) \\
&\leq L \left(\iint_{\Xi \times \Xi} d(x, y)^{\beta r} \pi(\mathrm{d}x, \mathrm{d}y) \right)^{1/r} (\mathbb{E} \zeta^q)^{1/q}. \tag{28}
\end{aligned}$$

Now note that $(\mathbb{E} \zeta^q)^{1/q} = \|\sigma\|_q$ where $\sigma(\cdot) := F_\zeta^{-1}(\cdot) \in \mathcal{S}$ is the generalized inverse distribution function. We obtain the desired result by taking the infimum in (28) over all possible measures with marginals P and \tilde{P} and after letting $\varepsilon \rightarrow 0$.

For the remaining inequality observe that

$$\left(\mathbb{E}_\pi d^{\beta r} \right)^{1/\beta r} = \|d\|_{\beta r} \leq \|d\|_r = (\mathbb{E}_\pi d^r)^{1/r}$$

by Hölder's inequality, so that

$$(28) \leq L \left(\iint_{\Xi \times \Xi} d(x, y)^r \pi(\mathrm{d}x, \mathrm{d}y) \right)^{\beta/r} \cdot \sup_{\sigma \in \mathcal{S}} \|\sigma\|_q.$$

This is the assertion. □

Corollary 28 (Continuity of the Average Value-at-Risk). *Suppose that $Y(x) - \tilde{Y}(y) \leq L \cdot d(x, y)$. Then*

$$\text{AV@R}_{\alpha;P}(Y) - \text{AV@R}_{\alpha;\tilde{P}}(\tilde{Y}) \leq \frac{L}{1-\alpha} w(P, \tilde{P}; d).$$

Proof. This is a special case of Proposition 27 for $r = 1$ and $q = \infty$ (cf. Example 6). □

Theorem 29 (Continuity of nested risk functionals). *Suppose that the random variables $Y: \Xi \rightarrow \mathbb{R}$ is Hölder continuous with constant L and exponent $\beta \leq 1$,*

$$|Y(x) - Y(y)| \leq L \cdot d(x, y)^\beta.$$

Then the nested risk functional $\mathcal{R}_{S_{1:T}}(Y)$ is continuous with respect to the nested distance, it holds that

$$\left| \mathcal{R}_{S_{1:T};P}(Y) - \mathcal{R}_{S_{1:T};\tilde{P}}(Y) \right| \leq \sup_{\sigma \in \mathcal{S}_t, t=1, \dots, T} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot L \cdot \mathbf{d}_r(P, \tilde{P})^\beta.$$

Proof. We infer from Proposition 27 with $\tilde{Y} = Y$ that

$$\begin{aligned}
\mathcal{R}_{S_T;P(\cdot|\cdot_{1:T-1})}(Y) - \mathcal{R}_{S_T;\tilde{P}(\cdot|\cdot_{1:T-1})}(Y) \\
\leq L \cdot \sup_{\sigma_T \in \mathcal{S}_T} \|\sigma_T\|_q \cdot w_r(P(\cdot|\cdot_{1:T-1}), \tilde{P}(\cdot|\cdot_{1:T-1}); c_T)^\beta, \tag{29}
\end{aligned}$$

where the terminal cost function is the distance as in the definition of the nested distance (cf. (18)),

$$c_T = d.$$

Define the random variables

$$Y_{T-1}(x_{1:T-1}) := \mathcal{R}_{S_T; P(\cdot | x_{1:T-1})}(Y) \quad \text{and} \quad \tilde{Y}_{T-1}(y_{1:T-1}) := \mathcal{R}_{S_T; \tilde{P}(\cdot | y_{1:T-1})}(Y),$$

so that we have

$$Y_{T-1}(x_{1:T-1}) - \tilde{Y}_{T-1}(y_{1:T-1}) \leq L \cdot \sup_{\sigma_T \in S_T} \|\sigma_T\|_q \cdot c_{T-1}(x_{1:T}, y_{1:T})^\beta$$

by (29) and the definition of the process c_t in (17). The random variables Y_{T-1} and \tilde{Y}_{T-1} thus satisfy the condition (27) with respect to the cost function c_{T-1} . So we may again apply Proposition 27 to the measures $P(\cdot | x_{1:T-2})$ and $\tilde{P}(\cdot | y_{1:T-2})$ and repeating this procedure for $t = T - 2$ down to $t = 0$ gives

$$\mathcal{R}_{S_{1:T}; P}(Y) - \mathcal{R}_{S_{1:T}; \tilde{P}}(Y) \leq \sup_{\sigma_t \in S_t, t=1, \dots, T} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot L \cdot c_0^\beta,$$

with terminal cost function $c_T = d$. We have that $c_0 = \text{dl}_r(P, \tilde{P})$ and thus

$$\mathcal{R}_{S_{1:T}; P}(Y) - \mathcal{R}_{S_{1:T}; \tilde{P}}(Y) \leq \sup_{\sigma_t \in S_t, t=1, \dots, T} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot L \cdot \text{dl}_r(P, \tilde{P})^\beta.$$

The result follows finally by exchanging the probability measures P and \tilde{P} . \square

Corollary 30 (Continuity of the nested Average Value-at-Risk). *Suppose that Y is Lipschitz continuous with constant L . Then the nested Average Value-at-Risk, nAV@R , is continuous with respect to the nested distance dl . More precisely, it holds that*

$$|\text{nAV@R}_{\alpha_{1:T}; P}(Y) - \text{nAV@R}_{\alpha_{1:T}; \tilde{P}}(Y)| \leq \frac{L}{1 - \alpha} \text{dl}_r(P, \tilde{P})$$

for every $r \geq 1$, where $\alpha \geq 1 - (1 - \alpha_1) \cdot \dots \cdot (1 - \alpha_T)$ (cf. (13)).

Proof. The statement for $r = 1$ is immediate by the definition of the nested Average Value-at-Risk, Corollary 28 and Theorem 29. The statement for general $r \geq 1$ follows from (20). \square

6 Dynamic equations and the martingale property

In what follows we consider multistage optimization problems with cost function

$$Q: \mathcal{Z}_{0:T} \times \Xi_{1:T} \rightarrow \mathbb{R},$$

where a sequence of subsequent decisions $z_t \in \mathcal{Z}_t$, $t = 0, \dots, T$, is chosen from $\mathcal{Z}_{0:T} = \mathcal{Z}_0 \times \dots \times \mathcal{Z}_T$. To account for risk-averse decision making under uncertainty we involve risk functionals at each stage.

Definition 31 (Policy). The random variable $z_t: \Xi \rightarrow \mathcal{Z}_t$ is a random *policy* or *decision* at time t , $t = 0, \dots, T$. The decision z_t is *nonanticipative* (or *adapted*) if $z_t: \Xi \rightarrow \mathcal{Z}_t$ is \mathcal{F}_t -measurable for every $t = 0, \dots, T$, abbreviated by $z_t \triangleleft \mathcal{F}_t$. The function $z: \Xi \rightarrow \mathcal{Z}_{1:T}$ with $z(x)_t := z_t(x)$ is nonanticipative (adapted; in short, $z \triangleleft \mathcal{F}$), if each component z_t is nonanticipative for every $t = 0, \dots, T$.

Remark 32. It is a consequence of the Doob–Dynkin lemma that z_t is nonanticipative if it depends solely on the information available at time $t \in \{0, \dots, T\}$, i.e., if $z_t(x_{1:T}) = \tilde{z}_t(x_{1:t})$ for some measurable function $\tilde{z}_t: \Xi_{1:t} \rightarrow \mathcal{Z}_t$ (cf. Kallenberg [11, Lemma 1.13] or Shiryaev [32, Theorem II.4.3]). As the filtration $\mathcal{F} = (\mathcal{F}_t)_{t=0}^T$ is the coordinate filtration it follows that every nonanticipative random decision $z \triangleleft \mathcal{F}$ can be written explicitly as

$$z_{0:T}(x_{1:T}) = \begin{pmatrix} z_0 \\ z_1(x_1) \\ z_2(x_1, x_2) \\ \vdots \\ z_T(x_1, \dots, x_T) \end{pmatrix}$$

for adequate, measurable functions $z_t: \Xi_{0:t} \rightarrow \mathcal{Z}_t$.

Definition 33 (Multistage optimization). Let $Q: \mathcal{Z}_{0:T} \times \Xi_{1:T} \rightarrow \mathbb{R} \cup \{\infty\}$ be a lsc. cost function. The risk-averse multistage optimization problem is

$$\inf_{z_{0:T} \triangleleft \mathcal{F}_{0:T}} \mathcal{R}_{S_{1:T}} \left(Q(z_{0:T}(\cdot); \cdot) \right), \quad (30)$$

where the infimum is among all adapted policies $z \triangleleft \mathcal{F}$. We emphasize and indicated the random component in (30) by ‘ \cdot ’.

Remark 34. To avoid confusions or ambiguities regarding the arguments of the function Q we separate the arguments $z \in \mathcal{Z}_{0:T}$ and $x \in \Xi_{1:T}$ explicitly and write $Q(z; x)$. This will turn out helpful in what follows, for example in expressions as $Q(z_{0:t-1}, z_{t:T}; x_{1:t}, x_{t+1:T})$.

Remark 35. Constraints of the form $z_{0:t}(x_{1:t}) \in \mathcal{Z}_t(x_{1:t}) \subseteq \mathcal{Z}_t$ for some multifunction $\mathcal{Z}_t(\cdot)$ appear naturally in applications involving optimization under uncertainty. They are easily incorporated in the problem formulation (30) just by employing the function $Q(z_{0:T}, x_{1:T}) \cdot \mathbb{1}_{\mathcal{Z}_t(x_{1:t})}(z_{0:T})$ instead of Q . This setting is not advisable for real world implementations, but convenient for the conceptual treatment envisaged here.

The multistage problem (30) thus consists in finding optimal functions $z_0, z_1(\cdot), \dots, z_T(\cdot)$ (only z_0 is deterministic) and therefore can be considered as *optimization on function spaces*.

6.1 The essential infimum

We shall make use of the following interchangeability principle, cf. also Shapiro [30]. For $z \in \mathcal{Z}$ fixed, the mapping $x \mapsto Q(z, x)$ is a random variable for which we write $Q(z, \cdot)$. In what follows we discuss the expression $\inf_z Q(z, \cdot)$ and its measurability. We refer to Karatzas and Shreve [12, Appendix A] for a formal definition of the essential infimum $\text{ess inf}_{z \in \mathcal{Z}} Q(z, \cdot)$, which is a measurable random variable as well.

Proposition 36. *Let \mathcal{Z} be a vector space and consider all policies with values $z(\cdot) \in \mathcal{Z}$. Then there exists a sequence $z_n(\cdot)$ of simple functions so that*

$$\lim_{n \rightarrow \infty} Q(z_n(\cdot), \cdot) = \text{ess inf}_{z(\cdot) \in \mathcal{Z}} Q(z(\cdot), \cdot) \quad \text{almost surely} \quad (31)$$

and $Q(z_n(\cdot), \cdot)$ is nonincreasing.

Proof. Denote the set of simple functions $z(\cdot) = \sum_{i=1}^k a_i \mathbb{1}_{A_i}(\cdot)$ by s . For $z(\cdot)$ and $z'(\cdot)$ simple functions define

$$z''(x) := \begin{cases} z(x) & \text{if } Q(z(x), x) \leq Q(z'(x), x), \\ z'(x) & \text{else,} \end{cases} \quad (32)$$

which is a simple function again and measurable. (The maximization (32) actually defines a directed set or preorder on s .) It holds that $Q(z''(\cdot), \cdot) \leq Q(z'(\cdot), \cdot)$ and $Q(z''(\cdot), \cdot) \leq Q(z(\cdot), \cdot)$ and the set $\{Q(z(\cdot), \cdot) : z \in s\}$ thus is closed under pairwise minimization. It follows from Karatzas and Shreve [12, Theorem A.3] that there is a sequence $z_n(\cdot)$ of simple functions so that

$$\operatorname{ess\,inf}_{z(\cdot)} Q(z(\cdot), \cdot) = \lim_{n \rightarrow \infty} Q(z_n(\cdot), \cdot) \quad \text{almost everywhere}$$

and thus the assertion. \square

Corollary 37. *Let s be a set of policies containing all simple functions and suppose that*

$$x \mapsto Q(z, x) \quad (33)$$

is upper semi-continuous for every $z \in \mathcal{Z}$. Then there exists a sequence $z_n(\cdot)$ of policies so that

$$\lim_{n \rightarrow \infty} Q(z_n(\cdot), \cdot) = \inf_{z \in \mathcal{Z}} Q(z, \cdot) \quad \text{almost everywhere.}$$

Proof. The set s contains the constant functions and thus

$$\inf_{z \in \mathcal{Z}} Q(z, x) = \inf_{z(\cdot) \in \mathcal{Z}} Q(z(x), x) \quad \text{for every } x.$$

We have that $\{x : \inf_{z \in \mathcal{Z}} Q(z, x) < \alpha\} = \bigcup_{z \in \mathcal{R}} \{x : Q(z, x) < \alpha\}$ for every $\alpha \in \mathbb{R}$ so that the additional assumptions ensure that $x \mapsto \inf_{z \in \mathcal{Z}} Q(z, x)$ is measurable. The assertion thus follows as

$$\inf_{z(\cdot) \in \mathcal{Z}} Q(z(\cdot), \cdot) = \operatorname{ess\,inf}_{z(\cdot) \in \mathcal{Z}} Q(z(\cdot), \cdot) = \lim_{n \rightarrow \infty} Q(z_n(\cdot), \cdot),$$

where $z_n(\cdot)$ is the sequence found in Proposition 36. \square

Convention 38. In what follows we shall always understand the measurable version when writing $\inf_{z \in \mathcal{Z}} Q(z, \cdot)$, i.e., we set

$$\inf_{z \in \mathcal{Z}} Q(z, \cdot) := \operatorname{ess\,inf}_{z(\cdot) \in \mathcal{Z}} Q(z(\cdot), \cdot). \quad (34)$$

The preceding Corollary 37 provides general conditions so that the convention is void and automatically valid in these cases.

Proposition 39 (Risk functional at the essential infimum, cf. Shapiro et al. [31, Proposition 6.60]). *Suppose that \mathcal{R} is continuous at $\inf_z Q(z, \cdot)$ with respect to convergence in L^p . Then it holds that*

$$\inf_{z(\cdot) \in \mathcal{Z}} \mathcal{R}(Q(z(\cdot), \cdot)) = \mathcal{R}\left(\inf_{z \in \mathcal{Z}} Q(z, \cdot)\right).$$

Proof. The result is a consequence Lebesgue's dominated convergence theorem in view of our setting (34) and the representation as nonincreasing limit given in (31). \square

6.2 Martingale characterization

Section 4.3, in particular Theorem 24, characterize the nested distance as a martingale process. This concept extends to the value process of the stochastic optimization problem when generalizing the concept of martingales. We incorporate risk awareness in the definition of the martingale term first and characterize the optimal solution of the multistage stochastic optimization problem as a martingale with respect to the risk functionals involved.

Definition 40 (Risk martingale). The stochastic process $v = (v_t)_{t=0}^T$ is a *submartingale* (*supermartingale*, resp.) with respect to the risk functionals $\mathcal{R}_{\mathcal{S}_t}$ (an \mathcal{R} -submartingale, for short), if

$$v_t \leq \mathcal{R}_{\mathcal{S}_{t+1}}(v_{t+1}) \text{ a.s.} \quad (v_t \geq \mathcal{R}_{\mathcal{S}_{t+1}}(v_{t+1}) \text{ a.s., resp.}) \quad (35)$$

for very $t \in \{0, 1, \dots, T\}$. The process v_t is an \mathcal{R} -martingale, if (35) holds with equality.

For the expectation, $\mathcal{R} = \mathbb{E}$, the notion of an \mathcal{R} -martingale (sub-, supermartingale, resp.) coincides with the usual term martingale (sub-, supermartingale, resp.).

Remark 41. A process $v = (v_t)_{t=0}^T$, which is an \mathcal{R} -submartingale, satisfies in addition

$$v_s \leq \mathcal{R}_{\mathcal{S}_{s+1:t}}(v_t), \quad 0 \leq s < t < T.$$

This follows as the risk functionals $\mathcal{R}_{\mathcal{S}_t}$ are monotone (Axiom A1) and from the recursive definition of the nested risk functional given in Definition 8.

Theorem 42. Let $z = (z_t)_{t=0}^T$ be an adapted policy. Then the process

$$v_t(x_{1:t}) := \mathcal{R}_{\mathcal{S}_{t+1:T}} \left(Q(z_{0:t}(x_{1:t}), z_{t+1:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \mid x_{1:t} \right) \quad (36)$$

is an \mathcal{R} -martingale with terminal value

$$v_T = Q(z_{0:T}(\cdot); \cdot). \quad (37)$$

Proof. Choosing $t = T$ in the defining equation (36) gives $v_T(x_{1:T}) = Q(z_{0:T}(x_{1:T}); x_{1:T})$ and thus (37).

Apply $\mathcal{R}_{\mathcal{S}_T}$ and it follows from (37) that

$$\begin{aligned} \mathcal{R}_{\mathcal{S}_T} \left(v_T \mid x_{1:T-1} \right) &= \mathcal{R}_{\mathcal{S}_T} \left(Q(z_{0:T}(x_{1:T}); x_{1:T}) \mid x_{1:T-1} \right) \\ &= \mathcal{R}_{\mathcal{S}_T} \left(Q(z_{0:T-1}(x_{1:T-1}), z_{T:T}(x_{1:T-1}, \cdot); x_{1:T-1}, \cdot) \mid x_{1:T-1} \right) \\ &= v_{T-1}(x_{1:T-1}), \end{aligned}$$

as z is adapted. This is the desired martingale property for $t = T - 1$.

Apply next $\mathcal{R}_{\mathcal{S}_{T-1}}$ to the latter equation and observe that

$$v_{T-2}(x_{1:T-2}) = \mathcal{R}_{\mathcal{S}_{T-1:T}} \left(Q(z_{0:T-1}(x_{1:T-1}); x_{1:T-1}) \mid x_{1:T-2} \right) = \mathcal{R}_{\mathcal{S}_{T-1:T}} \left(v_{T-1}(x_{1:T-1}) \mid x_{1:T-2} \right),$$

which is the assertion for $t = T - 1$. The general assertion is immediate by repeatedly applying the risk functional corresponding to the individual stage. \square

As a consequence we have the following immediate property of an optimal policy.

Corollary 43. Let $z_{0:T}^* : \Xi \rightarrow \mathcal{Z}_{0:T}$ be an optimal policy in the multistage stochastic optimization problem (30) and $v^* = (v_t^*)_{t=0}^T$ the value process (36) associated with the policy $z_{0:T}^*$. Then v^* is an \mathcal{R} -martingale and the starting value v_0^* is the solution of the optimization problem (30).

6.3 The value process is a martingale

Associated with the optimal solution of the reference problem (30) is an optimal policy. We shall characterize the evolution of this process now by highlighting their martingale properties.

Definition 44 (The value process). Let $z_{0:T} : \Xi \rightarrow \mathcal{Z}_{0:T}$ be a policy. The *value process* associated with the policy $z_{0:T}$ is $v(z) := (v_t(z))_{t=0}^T$. The marginal functions $v_t(z) := v_t(z_{0:t-1} | \cdot) : \Xi_{0:t} \rightarrow \mathbb{R}$ are defined by

$$v_t(z_{0:t-1} | x_{1:t}) := \inf_{z_{t:T} \triangleleft \mathcal{F}_{t:T}} \mathcal{R}_{S_{t+1:T}} \left(Q(z_{0:t-1}(x_{1:t}), z_{t:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) | x_{1:t} \right), \quad t = 0, \dots, T, \quad (38)$$

where the infimum in (38) is among all adapted processes $z_{t:T}(x_{1:T}) = \begin{pmatrix} z_t(x_1, \dots, x_t) \\ \vdots \\ z_T(x_1, \dots, x_t, \dots, x_T) \end{pmatrix}$.

Remark 45. The value process at initial time $t = 0$ is

$$v_0^* := \inf_{z_{0:T} \triangleleft \mathcal{F}_{0:T}} \mathcal{R}_{S_{1:T}} \left(Q(z_{0:T}(\cdot); \cdot) \right),$$

this value coincides with the risk-averse multistage stochastic program (30) given in Definition 33. The quantity v_0^* is a deterministic number and not random.

In addition, we have for $t = T$ that

$$v_T(z_{0:T-1} | x_{1:T}) = \inf_{z_T \triangleleft \mathcal{F}_T} Q(z_{0:T-1}(x_{1:T}), z_T(x_{1:T}); x_{1:T}),$$

so that the terminal value function does not involve a risk measure any longer and the terminal optimization problem is deterministic, i.e., not random either.

Theorem 46 (Submartingale characterization of the value process). *The value process is an \mathcal{R} -submartingale for any given policy $z_{0:T}$.*

Proof. We have that

$$\begin{aligned} \mathcal{R}_{S_{t+1}}(v_{t+1}) &= \mathcal{R}_{S_{t+1}} \left(\inf_{z_{t+1:T} \triangleleft \mathcal{F}_{t+1:T}} \mathcal{R}_{S_{t+2:T}} \left(Q(z_{0:t}(x_{1:t+1}), z_{t+1:T}(x_{1:t+1}, \cdot); x_{1:t+1}, \cdot) | x_{1:t+1} \right) \right) \\ &= \inf_{z_{t+1:T} \triangleleft \mathcal{F}_{t+1:T}} \mathcal{R}_{S_{t+1}} \left(\mathcal{R}_{S_{t+2:T}} \left(Q(z_{0:t}(x_{1:t+1}), z_{t+1:T}(x_{1:t+1}, \cdot); x_{1:t+1}, \cdot) | x_{1:t+1} \right) \right) \quad (39) \\ &= \inf_{z_{t+1:T} \triangleleft \mathcal{F}_{t+1:T}} \mathcal{R}_{S_{t+1:T}} \left(Q(z_{0:t}(x_{1:t+1}), z_{t+1:T}(x_{1:t+1}, \cdot); x_{1:t+1}, \cdot) | x_{1:t+1} \right), \quad (40) \end{aligned}$$

where we have employed (34) in (39).

The result follows now, as that value process (38) is the infimum among all $z_{t:T} \triangleleft \mathcal{F}_{t:T}$, while the infimum in (40) is among $z_{t:T} \triangleleft \mathcal{F}_{t+1:T}$, which is one dimension less. \square

Dynamic optimization employs verification theorems which give sufficient conditions for a solution to the optimal control problem, cf. Fleming and Soner [7, Theorems 5.1 and 5.2]. The following theorem provides the corresponding statement for the risk-averse multistage stochastic problem.

Theorem 47 (Martingale characterization, dynamic equations, verification theorem). *For the value process it holds that*

$$v_0^* = \inf_{z_{0:T} \triangleleft \mathcal{F}_{0:T}} \mathcal{R}_{S_{1:T}}(v_t(z_{0:t})), \quad t \in \{0, 1, \dots, T\}.$$

More generally, for $s < t$ we have the recursive equations

$$v_s(z_{0:s-1}) = \inf_{z_{s:t} \triangleleft \mathcal{F}_{s:t}} \mathcal{R}_{S_{s+1:t}}(v_t(z_{1:s}, z_{s+1:t})). \quad (41)$$

Proof. Applying the conditional risk functional $\mathcal{R}_{S_t}(\cdot \mid x_{1:t-1})$ to (38) gives

$$\begin{aligned} \mathcal{R}_{S_t}(v_t(z_{0:t-1} \mid x_{1:t}) \mid x_{1:t-1}) &= \mathcal{R}_{S_t} \left(\inf_{z_{t:T} \triangleleft \mathcal{F}_{t:T}} \mathcal{R}_{S_{t+1:T}} \left(Q(z_{0:t-1}(x_{1:t}), z_{t:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \mid x_{1:t} \right) \mid x_{1:t-1} \right) \\ &= \inf_{z_{t:T} \triangleleft \mathcal{F}_{t:T}} \mathcal{R}_{S_t} \left(\mathcal{R}_{S_{t+1:T}} \left(Q(z_{0:t-1}(x_{1:t}), z_{t:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \mid x_{1:t} \right) \mid x_{1:t-1} \right) \\ &= \inf_{z_{t:T} \triangleleft \mathcal{F}_{t:T}} \mathcal{R}_{S_{t:T}} \left(Q(z_{0:t-1}(x_{1:t}), z_{t:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \mid x_{1:t-1} \right), \end{aligned} \quad (42)$$

where we have used the monotonicity axiom, A1 and Proposition 3 to obtain “ \leq ” in (42). The converse inequality “ \geq ” involves the Lebesgue Dominated Convergence Theorem and is a consequence of Proposition 39.

At this stage take the infimum with respect to $z_{t-1} \triangleleft \mathcal{F}_{t-1}$ and thus

$$\begin{aligned} \inf_{z_{t-1} \triangleleft \mathcal{F}_{t-1}} \mathcal{R}_{S_t}(v_t(z_{0:t-1} \mid x_{1:t}) \mid x_{1:t-1}) &= \inf_{z_{t-1:T} \triangleleft \mathcal{F}_{t-1:T}} \mathcal{R}_{S_{t:T}} \left(Q(z_{0:t-1}(x_{1:t}), z_{t:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \mid x_{1:t-1} \right) \\ &= v_{t-1}(z_{0:t-2} \mid x_{1:t-1}), \end{aligned}$$

which is the martingale property of the value process $v(z)$. The remaining equation (41) follows in line with Remark 41.

The converse inequalities follow from the submartingale characterization, Theorem 46. \square

The dynamic equations derived in this section can be employed to characterize optimal solution of the multistage stochastic optimization problem. The conceptual advantage lies in the fact that each stage can be considered for its own. For this the dynamic equations can be employed in algorithms to improve suboptimal policies at each stage individually.

7 Continuity of risk-averse multistage programs

The value of the risk-averse multistage stochastic optimization problem (30) depends on the probability measure P . We shall make this explicit by writing

$$v_P := \inf_{z_{0:T} \triangleleft \mathcal{F}_{0:T}} \mathcal{R}_{S_{1:T}; P} \left(Q(z_{0:T}(\cdot); \cdot) \right). \quad (43)$$

It is known that the *risk-neutral* version of the multistage problem (43) is continuous with respect to changing the probability measure.

The following main result elaborates continuity of the *risk-averse* problem with respect to the nested distance and gives the modulus of continuity explicitly.

Theorem 48 (Continuity of the risk-averse MSO problem). *Suppose that*

$$x \mapsto Q(z; x), \quad z \in \mathcal{Z},$$

is uniformly Lipschitz, i.e.,

$$|Q(z; x) - Q(z; y)| \leq L \cdot d(x, y) \quad \text{for all } x, y \in \Xi_{1:T} \text{ and } z \in \mathcal{Z} \quad (44)$$

and

$$z \mapsto Q(z; x) \quad (x \in \Xi_{1:T})$$

is convex for every x fixed. Then the risk-averse optimization problem (30) is continuous with respect to changing the probability measure. More specifically, we have that

$$|v_P - v_{\tilde{P}}| \leq \sup_{\sigma \in \mathcal{S}_t, t=1, \dots, T} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot L \cdot \text{dl}_r(P, \tilde{P}),$$

where the exponents r and q are Hölder conjugates, $\frac{1}{r} + \frac{1}{q} = 1$.

Remark 49. The assumption on Lipschitz continuity of the function Q notably insures the Convention 38 as Q is particularly usc., cf. (33).

Proof of Theorem 48. To compare with the second problem $v_{\tilde{P}}$ define the new policy

$$\tilde{z}_t(y_{1:t}) := \mathbb{E}_\pi(z_t(x) \mid \text{pr}_t(x, y) = y_{1:t}),$$

where $\text{pr}_t(x_{1:t}, y_{1:t}) := y_{1:t}$ is the projection onto the second marginal and consider the specific random variables

$$Y_t(x_{1:t}) := \mathcal{R}_{\mathcal{S}_{t+1:T}; P(\cdot | x_{1:t})} \left(Q(z_{0:t}, z_{t+1:T}(x_{1:t}, \cdot); x_{1:t}, \cdot) \right) \quad (45)$$

and

$$\tilde{Y}_t(y_{1:t}) := \mathcal{R}_{\mathcal{S}_{t+1:T}; \tilde{P}(\cdot | y_{1:t})} \left(Q(z_{0:t}, \tilde{z}_{t+1:T}(y_{1:t}, \cdot); y_{1:t}, \cdot) \right), \quad (46)$$

where $z_{0:t} \in \mathcal{Z}$ is fixed and ‘ \cdot ’ indicates the random component.

For $\varepsilon > 0$ pick a policy $z = (z_{0:t}(x_{1:t}))_{t=1}^T$ so that

$$v_P > \mathcal{R}_{\mathcal{S}_{1:T}; P} (Q(z_{0:T}(\cdot); \cdot)) - \varepsilon. \quad (47)$$

Further, let the measure $\pi(\cdot, \cdot)$ have conditional marginals $P(\cdot)$ and $\tilde{P}(\cdot)$ with respect to the nested distance, cf. (24).

In line with Definition 15 we set $c_T := d$ and proceed by backwards induction from $t = T$ down to $t = 0$.

Base case: Note that $Y_T(x_{1:T}) = Q(z_{0:T}; x_{1:T})$ and $\tilde{Y}_T(y_{1:T}) = Q(z_{0:T}; y_{1:T})$. By Lipschitz continuity (44) it holds that $\tilde{Y}_T(y_{1:T}) - Y_T(x_{1:T}) \leq L \cdot d(x_{1:T}, y_{1:T})$. This is the statement

$$\tilde{Y}_t(y_{1:t}) - Y_t(x_{1:t}) \leq L \cdot \sup_{\sigma \in \mathcal{S}_{t+1:T}} \|\sigma_{t+1}\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot c_t(x_{1:t}, y_{1:t}) \quad (48)$$

for the case $t = T$ (and by setting the empty product to $\prod_{t \in \emptyset} f_t := 1$).

Inductive step: In what follows we shall employ the statement (48) as induction hypothesis and deduce the statement for $t - 1$ instead of t . From Jensen’s inequality we infer that

$$Q(\tilde{z}(y); y) = Q(\mathbb{E}_\pi(z(x) \mid \text{pr}_t(x, y) = y); y) \leq \mathbb{E}_\pi(Q(z(x); y) \mid \text{pr}(x, y) = y). \quad (49)$$

To be more specific we emphasize that z is a vector of functions, $z = (z_t)_{t=0}^T$ and further, each z_t is a function of the variables x_1, \dots, x_t , $z_t = z_t(x_{1:t})$. Jensen's inequality applies to each function z_t and each argument x_t separately, so that the inequality (49) is actually the result of applying Jensen's inequality t times repeatedly at each stage t .

Now let ζ be chosen so that $\mathbb{E} \tilde{Y}_t \zeta > \mathcal{R}_{\mathcal{S}_{t:T}; \tilde{P}(\cdot | y_{1:t-1})}(\tilde{Y}_t) - \varepsilon'$ and $\text{AV@R}_\alpha(\zeta) \leq \frac{1}{1-\alpha} \int_\alpha^1 \sigma(u) du$ ($\alpha \in (0, 1)$) for some $\sigma(\cdot) \in \mathcal{S}_t$. As the risk functional is recursive we deduce from (45) and (46) that

$$\begin{aligned} \tilde{Y}_{t-1} - Y_{t-1} - \varepsilon' &= \mathcal{R}_{\mathcal{S}_{t:T}; \tilde{P}(\cdot | y_{1:t-1})}(\tilde{Y}_t) - \varepsilon' - \mathcal{R}_{\mathcal{S}_{t:T}; P(\cdot | x_{1:t-1})}(Y_t) \\ &\leq \mathbb{E}_\pi Q(\tilde{z}(y); y) \zeta(y) - \mathbb{E}_\pi Q(z(x); x) \zeta(y) \\ &\leq \mathbb{E}_\pi \mathbb{E}_\pi (Q(z(x); y) \mid \text{pr}(x, y) = y) \zeta(y) - \mathbb{E}_\pi Q(z(x); x) \zeta(y), \end{aligned}$$

where we have used (49). By the tower property of the conditional expectation, Lipschitz continuity (44) and Hölder's inequality it follows further that

$$\begin{aligned} \tilde{Y}_{t-1} - Y_{t-1} - \varepsilon' &\leq \mathbb{E}_\pi Q(z(x); y) \zeta(y) - \mathbb{E}_\pi Q(z(x); x) \zeta(y) \\ &\leq \mathbb{E}_\pi \zeta(y) c_t(x_{1:t}, y_{1:t}) \\ &\leq L \sup_{\sigma \in \mathcal{S}_{t:T}} \|\sigma_t\|_q \cdot \dots \cdot \|\sigma_T\|_q w_r(P(\cdot | x_{1:t-1}), \tilde{P}(\cdot | y_{1:t-1}); c_t) \\ &= L \sup_{\sigma \in \mathcal{S}_{t:T}} \|\sigma_t\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot c_t(x_{1:t-1}, y_{1:t-1}), \end{aligned}$$

as π has conditional marginals $\tilde{P}(\cdot | y_{1:t-1})$ and $P(\cdot | x_{1:t-1})$. By letting $\varepsilon' \rightarrow 0$ we get the assertion (48) for $t - 1$. By repeatedly applying the previous reasoning we thus get that

$$\tilde{Y}_0 - Y_0 \leq L \sup_{\sigma \in \mathcal{S}_{1:T}} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot \text{dl}_r(P, \tilde{P}). \quad (50)$$

Now note that $v_P > Y_0 - \varepsilon$ by (47) and we thus have found a policy \tilde{z} so that $v_{\tilde{P}} \leq \tilde{Y}_0$. It follows with (50) that

$$v_{\tilde{P}} - v_P \leq \tilde{Y}_0 - (Y_0 - \varepsilon) \leq L \sup_{\sigma \in \mathcal{S}_{1:T}} \|\sigma_1\|_q \cdot \dots \cdot \|\sigma_T\|_q \cdot \text{dl}_r(P, \tilde{P}) + \varepsilon.$$

The result finally follows by letting $\varepsilon \rightarrow 0$ and by interchanging the role of P and \tilde{P} . □

8 Summary

This paper addresses risk-averse stochastic optimization problems. To define the risk functionals based on partial observations we introduce conditional risk measures first. They are defined on fibers and can be composed to nested risk measures. We demonstrate that these nested risk measures are continuous and we establish the modulus of continuity. As a consequence, the optimization problems are continuous as well, these problems inherit the modulus of continuity from the risk functionals.

All results come along with characterizations as generalized martingales. It is demonstrated that the underlying distance is a usual martingale with respect to the natural filtration. The value functions are shown to follow a generalized, risk-averse martingale pattern as well.

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