The Scenario Approach: a Tool at the Service of Data-Driven Decision Making

M.C. Campi*, A. Carè

Dipartimento di Ingegneria dell'Informazione - Università di Brescia, via Branze 38, 25123 Brescia, Italia.

S. Garatti

Dipartimento di Elettronica, Informazione e Bioingegneria - Politecnico di Milano, piazza Leonardo da Vinci 32, 20133 Milano, Italia.

Abstract

In the eye of many control scientists, the theory of the scenario approach is a tool for determining the sample size in certain randomized control-design methods, where an uncertain variable is replaced by a random sample of scenarios. This point of view is rooted in the history of the scenario approach and stands on a long track record of successful applications. However, in the last two decades the theory of the scenario approach has gone beyond its original motivations and applications, and has unveiled some fundamental relationships between the complexity of a design and its generalization capabilities. The new knowledge brought by the theory provides a solid ground for a framework where data can be exploited in a flexible and wise manner throughout a large variety of engineering activities. By this article we aim at providing an access point to a set of state-of-the-art results in the theory of the scenario approach that can be valuable to target important challenges in modern control-design and decision-making at large. In the first part of the article, we introduce a set-up for decision-making where the role of prior knowledge and user preferences can, and should, be distinguished from the role of data. Then, we show that the theory of the scenario approach offers a platform for conjugating heuristic approaches, which

*Corresponding author

in complex contexts are unavoidably based on incomplete and possibly imprecise information, with a solid theory for certifying the validity of the output of the decision process.

Keywords: data-driven methods, scenario approach, control, generalization.

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1. Introduction

1.1. The era of data

Opportunities and challenges intertwine in the era of data. On the one hand, pervasive networks of smart sensors collect, process and store measurements in extremely fast, reliable and cheap ways. On the other hand, control scientists are haunted by the question of how one can take advantage of the increasing availability of data, either to improve existing control systems or to bring order to newly born systems of systems.

The weaving of opportunities with old and new challenges is apparent in many fields; we mention below but a few, in full awareness that many others could have deserved to be included in the list.

- **Automotive systems control.** The availability and smart processing of real-time traffic measurements has made it possible to deploy automatic and effective coordinated ramp metering strategies for freeways, Papageorgiou & Papamichail (2014); Seo et al. (2017). Complex systems of systems arise not only from the conscious striving to go beyond ordinary solutions at the frontier of automotive technology (e.g., in the control of autonomous vehicles and of the systems arising from their interconnection Guanetti et al. (2018)), but also as technology-driven side-effects: for example, it has become urgent to deal with the disrupting effects of having many (ordinary) human drivers in (ordinary) cars, all relying simultaneously on the same (ordinary) mapping and route planning software, Macfarlane (2019).

- **Power generation and dispatchment.** At the dawn of the Second Industrial Revolution, connecting steam turbines to dynamos made large

scale power generation possible. Since then, energy has been generated and rationally distributed by leveraging well-understood physical laws, but the increasing penetration of renewable energy poses new and significant challenges to the planning and operation of modern power grids. In particular, power generation now depends on uncertain, hardly predictable, phenomena, ranging from weather conditions to individual and social human behaviors. The underpinning for modern decisions and control schemes in this field has necessarily to include, besides physics, economic and social sciences, as well as conceptual tools to integrate historical and real-time data in the decision process, Li et al. (2020).

- **Real-time control of biological systems.** Feedback, as a mechanism to ensure homoeostasis, has long been recognized as a central feature of life, Cannon (1939); Hoagland & Dodson (1995), and control scientists have been increasingly at work to design feedback loops to re-establish homoeostasis in medical patients. A remarkable example is the development of an artificial pancreas for keeping blood glucose into a healthy range. While the first closed-loop algorithms computed insulin infusion rates with poor information about the process to be controlled, modern control algorithms aim at combining a better understanding of the glucose metabolism with a refined exploitation of historical and real-time data. An account of this evolution can be found in Quiroz (2019), where one of the main future challenges in the design of glucose control algorithms is identified as the ability to address the large diversity of metabolic conditions that a diabetic patient can face in his/her daily life.

- **Medical computer-aided diagnosis.** A wise combination of domain knowledge, which remains fundamental to direct design efforts, and of data availability has fostered major progresses in medical computer-aided diagnosis. We just recall here the case of the automated classification of images of skin lesions as benign or malignant, which might soon become a widely available tool at the disposal of smartphone customer users, see

Esteva et al. (2017).

- **Machine-Learning-in-the-loop technologies.** Traditionally, an automatic classifier returns a result to a human being who is then responsible to make an informed decision (in the skin lesions example, this human being is a medical doctor or directly the patient). More and more often, however, classification algorithms, or other machine learning algorithms that are trained on data in a black-box manner, can be found in automated control loops, where they play the role of soft sensors, similarly to traditional state estimators. Examples range again from automotive systems (where classifiers are used *inter alia* for obstacle detection and avoidance, Gruyer et al. (2017); Devi et al. (2020)) to medical applications such as the artificial pancreas, Cappon et al. (2019). The more black-box learning algorithms enter safety-critical decision loops, the more urgent the need for a solid, scientific understanding of their limits and potentials.

Overall, it is clear that the increasing availability of computational power and of distributed large-scale optimization techniques enables the deployment of innovative data-driven decisions and control schemes, but it is also clear that this extraordinary potential cannot be fully expressed until it gets backed by a solid theoretical understanding.

1.2. Indirect vs. direct methods

In traditional model-based control, the "driving power" of data is often employed by engineers in the modeling of the reality to be controlled. In a typical workflow, first principles dictate the model class into which a reasonable representative of the reality should be sought, and experimental data are used to find the best parameter values to be plugged into the model equations. At the end of this modeling effort, the engineer has obtained a model that describes with a sufficient accuracy the way in which reality is expected to behave in response to various signals. At the final stage of the workflow, the engineer is called to

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use the model to make decisions, set-up policies and control mechanisms so as to optimize the objective values. This is what is called an *indirect approach*.

The increasing intricacy of many modern problems challenges this traditional design workflow: in fact, the input-output behavior of many modern systems is hardly captured by the class of models that can be found in the traditional system modeling toolboxes. Even in those lucky cases where, *in the eyes* of the data-analyst, available models look like acceptable descriptors of the reality, it is often the case that, *in the hands* of the control engineer, they lead to unsatisfactory control performances. The reason is that the metric according to which a model is said to capture "well" the hidden nature of the complex system is usually not tailored to the needs of the control engineer, whose goal is optimizing specific objective functions. This explains why the recent history of feedback control has witnessed a surge of the so-called *direct methods* (see e.g. Åström & Hägglund (1995); Safonov & Tsao (1995); Hjalmarsson et al. (1998); Guardabassi & Savaresi (2000); Gerencsér et al. (2002); Hjalmarsson (2002); Campi et al. (2002); Van Heusden et al. (2011); Bazanella et al. (2011); Moore (2012); Hou & Wang (2013); Formentin et al. (2014); Hori et al. (2016); Sutter et al. (2017); Karimi & Kammer (2017); Apkarian & Noll (2018); Novara & Milanese (2019); Formentin et al. (2019); Chiluka et al. (2021)). In direct methods, one uses the available information that is carried by experiments to directly assess how the controllers in a given class behave in various conditions, and this forms the ground on which a controller is selected; this is opposed to the traditional indirect methods where significant effort is spent in preliminarily approximating reality by means of a model.

The development of direct methods for decisions and control is part of a broader scientific trend, which is largely technology-driven and is not limited to the control community (the reader is referred to the box "Direct, goal-oriented approaches: a cultural, technology-driven trend" for more discussion on this point).

Direct, goal-oriented approaches: a cultural, technology-driven trend.

According to a highly influential paper, Breiman et al. (2001), two cultures abide in statistical sciences. The classic one aims at using data for modeling the data generation mechanism: if successful, this culture generates powerful tools not only to control but also to describe reality (control engineering practice reflects this culture when, e.g., the transfer function of a second-order linear system is estimated from noisy data); on the other hand, as the complexity of the data generation mechanism increases, modeling becomes more and more an ambitious task, prone to detrimental oversimplifications. Hence, the other culture aims at using data to inform problem-oriented procedures, and to issue certificates on the statistical effectiveness of such procedures: this task can be accomplished under much milder and realistic assumptions even in the presence of very complex data generating mechanisms.

In the context of statistical learning, Vapnik, see Vapnik (2013) - Section 1.9, formulated the following principle to be applied in the presence of restricted information: "When solving a given problem, try to avoid solving a more general problem as an intermediate step". In the wake of this advice, many recent advances in machine learning (in the fields of supervised classification and regression, Krizhevsky et al. (2012); clustering, Ghasedi Dizaji et al. (2017); reinforcement learning, Silver et al. (2016); etc.) share an agnostic approach with respect to the "true" or "best" description of the system at hand and focus on directly optimizing a cost function that maps the possible options that are available to the decision-maker into a value that quantifies the level of satisfaction with the selected option.

Generally speaking, direct approaches become more urgent as technology enables one to address problems of increasing complexity, which is the trend that applied science is nowadays experiencing at an increasingly fast

pace. The reader is also referred to Norvig (2017) for a thought-provoking discussion on matters relating to the topic of this box.

1.3. Content of this article

In this article, we present a framework for direct data-driven decision-making called the *scenario approach* that benefits from a series of technical results that have been established in the last two decades.

In the next Section 2, the scenario approach is introduced at a rather informal level as a general methodology to govern, in a well-grounded manner, the interplay between prior knowledge and data in decision-making. Some space (the whole Section 2.4) is devoted to introducing the probabilistic point of view that allows one to assess the quality of a decision not only with respect to the data collected before the decision is made, but also with respect to the infinitely larger set of the unseen cases (those that can occur at the time the decision is applied). The section ends with a non-technical preview (based on a simple example in Section 2.5) of the kind of statistical evaluations that are possible thanks to the theory of the scenario approach. This prepares the ground for the following, more technical, Section 3. Section 3 provides an easily accessible, but technical, gallery of results: moving from simple decision schemes based on convex worst-case optimization, the reader is gradually introduced to the state-of-the-art of the scenario theory, which encompasses non-convex optimization and a general decision-making framework. The article is closed by the concluding Section 4.

2. A set-up for direct data-driven decision-making: the scenario approach

2.1. When is a decision good?

Many decision and control problems can be abstracted as the problem of choosing an object x from within a decision set $\mathcal{X}$. For example, in control-design, x is a vector of parameters representing a controller in a given class; in

financial portfolio optimization, x is a vector where each element is the amount of money to be invested in a given asset; in binary classification, x encodes, according to some predefined rule, a function that maps any possible relevant object into a label belonging to $\{0, 1\}$, etc.

We will denote by x^* the specific object that represents the final output of the decision process. The quality of x^* can be judged in relation to intrinsic and extrinsic criteria as described in what follows.

1) INTRINSIC QUALITY

We call "intrinsic" a quality of x that depends only on x itself, in the light of cemented knowledge and preferences that are available to the decision-maker. Here are some examples.

Example (Filter Design). *If x is a digital filter, a good x may be expected to have a cut-off frequency in a certain range, to be physically realizable, possibly simple and cheap to implement, etc.*

Example (Prediction interval). *If x is an interval used to predict an unknown variable, its width should be small for the prediction to be informative.*

Example (Home temperature control system). *If x is a temperature controller, it should be designed so that, in nominal conditions, the control operates fast enough, overshooting and oscillations are limited, energy consumption is minimized, etc.*

2) EXTRINSIC QUALITY

The "extrinsic" quality refers to the performance of x in relation to various operating *situations* that may occur when x is applied. A bit more formally, we can think of the occurrence of a *situation* as an assignment of values to a vector of variables that we denote with the symbol δ, and, hence, extrinsic quality refers to the couple (x, δ): for every δ, x attains a performance as measured by a suitable indicator and the extrinsic quality refers to the variability of performances achieved by x as δ takes value in its range of variability.

Example (Filter Design). *If the aim of the filter x is to work in a mobile device as an audio channel equalizer and the frequency response of the channel is δ, then the extrinsic quality of x refers to the performance of x in relation to a variety of channels δ, and "high extrinsic quality" may refer to the ability of x to perform well over a large portion of δ's that may be encountered in the lifespan of the device.*

Example (Prediction interval). *Suppose that the interval x is used to predict how effective a medical therapy is. Since the effectiveness depends on the patient δ to whom the therapy is administered, the extrinsic quality may refer to how large the portion of potential patients for which x provides a correct prediction is.*

Example (Home temperature control system). *The performance of a building temperature controller x can be affected by the weather conditions and the minute actions of people in the building (which are partly unpredictable and will certainly differ from hour to hour and from day to day). Here, δ can be identified with a vector that includes quantities such as the external temperature and other weather conditions, the average amount of people in the building in a given time horizon, the number and the size of windows that happen to be open, etc., and the extrinsic quality may refer to the capability of the controller to keep the temperature within admissible limits for various δ.*

As an additional example, we anticipate that in Section 2.5 we shall be dealing with a simple machine learning problem where the prediction of an inaccessible variable is given by an interval that depends on a variable that can be measured. Similarly to the above "prediction interval" example, in that context the width of the interval is the intrinsic quality, while the extrinsic quality refers to the portion of cases that are mispredicted by the interval model.

2.2. The limits of knowledge

In traditional decision processes, the intrinsic quality and the extrinsic quality of a candidate solution are often judged at the decision-making stage on the

ground of available models.

Example (Traditional control design). *Let x be a controller to be applied to a linear plant whose poles (expressed in a vector δ) are somewhat uncertain. If the variability of δ is known, imposing suitable phase and gain margins may ensure that the design will work well for all the relevant values of δ.*

On the other hand, when we move from simple to complex application domains, we often experience that δ refers to articulated and elusive portions of the real world for which it is extremely difficult to obtain a satisfactory and complete model and, hence, a model-based assessment of the extrinsic quality becomes impractical. Under these circumstances, one may advocate the use of first-hand data, i.e., observed instances of δ, along a *direct decision-making approach*. This is the condition in which the scenario approach finds its natural application, as explained in the next section (see also the box "When should we consider using the scenario approach?" for a quick summary of the main ideas and applications domains to which the scenario approach has been applied).

2.3. The principles of the scenario approach

We shall denote the empirical instances of δ that are available at the decision-making stage by $\delta^{(1)}, \ldots, \delta^{(N)}$, and refer to them as "*scenarios*". When the variability of δ is difficult to describe by means of a model, the fact that δ impacts on the extrinsic performance suggests a different way of proceeding and, at an informal level, the scenario approach prescribes to choose the candidate solution x that

(i) works well for the scenarios $\delta^{(1)}, \ldots, \delta^{(N)}$

and, subject to (i),

(ii) optimizes the intrinsic quality.

In this way, the intrinsic quality is pursued directly while the scenarios $\delta^{(1)}, \ldots, \delta^{(N)}$ are used as an *empirical substitute* of the infinite amount of situations δ that
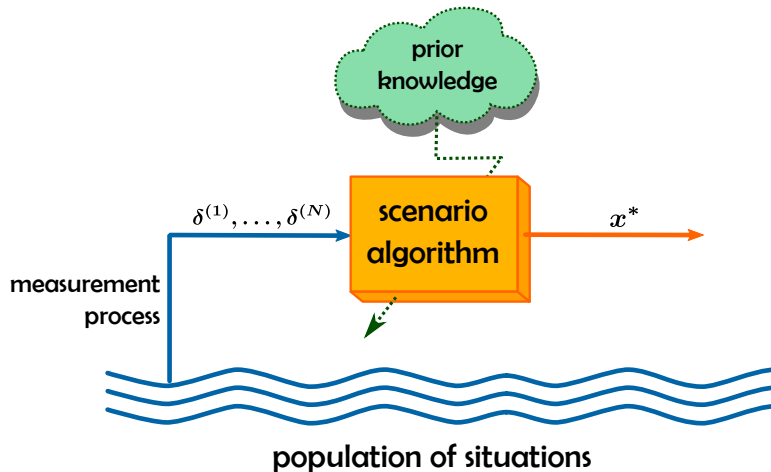
Figure 1: A scenario algorithm is a way to generate decisions able to cope with uncertainty. The picture shows in yellow the scenario box. It is constructed on the ground of prior knowledge about the problem at hand and is fed by a sample of situations (scenarios) elicited from a (typically infinite) population of situations.

could occur in a future use of the decision to heuristically secure the extrinsic quality.

In applications, (i) and (ii) must be formulated quantitatively and choosing suitable indicators is highly problem-dependent. Moreover, while the application may suggest the meaning of the expression "x works well for δ", still the decision-maker retains the right of deciding whether to enforce that the solution works well for all the N scenarios $\delta^{(i)}$ or rather for just some of them. There is also much flexibility in constructing the indicator of the intrinsic quality and the domain in which the solution is sought: they are typically based on prior knowledge and background preferences, but, as we shall see in Section 2.5, they can also be influenced by informal reasoning, second-hand information, guesses and even conjectures about the variability of δ. The workflow for the scenario approach is illustrated in Figure 1.

Importantly, making a wise decision based on observations requires tools to gain confidence in that the decision is good both intrinsically and extrinsically. While the intrinsic quality is directly measured (and optimized) because it only depends on x, replacing the large set of the unseen situations δ's with a

sample of scenarios is heuristic in nature and, when applied naively, may suffer from excessive empiricism: the final decision x^* performs well with respect to the instances $\delta^{(1)}, \ldots, \delta^{(N)}$ but it may fall short for other instances of δ that eventuates *after* the decision x^* has been implemented. The aim of the theory of the scenario approach is precisely that of taking control on this *generalization* issue.[1] This is important because only a trustful evaluation of the actual extrinsic quality along with the evidence about the intrinsic quality gives the decision-maker a handle to judge the solution. Based on a fair judgment, the decision-maker can:

(i) decide whether to "buy" the solution or to discard it;

(ii) decide how to subsequently use the solution, if the latter is part of a bigger decision process;

(iii) if the solution is not satisfactory, go back to the original choices in the problem formulation and re-calibrate them to construct solutions which are better aligned with the desires (for example, the definition of the intrinsic criterion can be modified, the size of the sample $\delta^{(1)}, \ldots, \delta^{(N)}$ can be re-tuned, etc.);

(iv) re-design the decision set $\mathcal{X}$ itself from which x^* is selected.

The main mathematical tool the scenario approach is based on is probability theory and the concepts introduced at a high-level in this section will be better formalized in the light of probability theory in the next section. Before proceeding, however, we feel advisable to highlight here two cornerstones of the theory as it will emerge in the remainder of this article.

(a) While one way of quantifying the extrinsic quality is by validation, this requires using data points for testing rather than designing. One key mes-

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[1] This theory is presented in Section 3. When the problem at hand belongs to the field of machine learning, the scenario theory of Section 3 is in the same vein as, and contributes to, the generalization theory of machine learning. In this case, the ability of a solution to perform well on the sample while having poor generalization capabilities has also been called *overfitting*. More generally, the scenario theory applies across various fields in decision making.

sage of the scenario approach is that data can be used to simultaneously make a design while also evaluating its extrinsic quality. This is made possible by an exploitation of the information contained in the data beyond what traditional approaches can do.

(b) In modern control and decision problems that deal with complex systems, besides scenarios one wants to exploit prior knowledge that comes from various sources, often including some that, while not completely trustworthy, can still be of help to obtain a satisfactory solution. However, if prior knowledge turns out to be lacking or even defective its use may downgrade the quality of the solution. Importantly, in the scenario approach the correctness of the evaluation of the extrinsic quality (which is not directly observable) remains intact independently of the correctness of the prior. This fact (which may came to the reader's surprise) is established in subsequent sections and is named "separation principle". The separation principle implies that the decision-maker can always correctly judge the impact of the prior information on the solution in terms of intrinsic (which is directly measurable) and extrinsic quality and decide to discard a prior if it is deemed unreliable. In this respect, it can be said that the scenario approach meets the pressing need in modern data-driven decision-making for a sound integration of domain knowledge and priors having uneven levels of trustworthiness with first-hand information given by data.

When should we consider using the scenario approach?

The scenario approach employs the data to directly target a design problem, without going through an intermediate step aiming at finding a description of the mechanism that generates δ. The scenario approach is of interest when:

  (I) the value taken by δ can significantly impact on the performance;

  (II) δ is made up of interconnected variables which take value according

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to mechanisms that are difficult, too expensive or even impossible to describe satisfactorily.

Because of (I), neglecting the variability of δ is not an option as this would lead to poor decisions. At the same time, (II) makes it impractical to find a description of the mechanism underlying the generation of δ. Hence, a good decision can be pursued heuristically driven by a direct use of the data. In this context, the scenario approach provides a theory to quantify the extrinsic quality of a decision and drives the user towards a wise selection.

Conditions (I) and (II) apply in innumerable applications, for example where the instances of δ are:

- pairs of input and output signals collected from an automotive system in different operational and environmental conditions, e.g., from a car to be controlled with different tires and in different road conditions, Rallo et al. (2016); Formentin et al. (2018);

- dynamical characteristics of an electro-mechanical system for the force control in brake-by-wire actuation, Riva et al. (2020);

- the power generation and load values in a power system with intermittent and distributed sources, Modarresi et al. (2019); Geng & Xie (2019); Geng et al. (2021);

- the parameters (as obtained, e.g., by an Intra-Venous Glucose Tolerance Test) describing the insulin-glucose system of the individuals for which an automated blood glucose regulator has to be designed, Borri et al. (2017); Di Ferdinando et al. (2021);

- the pharmacokinetic parameters of the individuals for which a drug administration protocol has to be designed, Sopasakis et al. (2018);

- the electrocardiogram of patients in cardiac arrest before and after an emergency treatment, Carè et al. (2018);

- stock market values, Pagnoncelli et al. (2012); Calafiore (2013); Ramponi & Campi (2018); Arici et al. (2021);

- inflows and outflows in systems of rivers, Nasir et al. (2018);

- the transfer functions of the different communication channels that can be encountered by a mobile communication device, Carè et al. (2015);

- various demand profiles for selling products, Carè et al. (2014);

- disturbances, Campi et al. (2009b), or actuation errors, Carè et al. (2019), in largely unpredictable environments.

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2.4. Mathematical foundations: a probabilistic framework

Probability is, first and foremost, a measuring tool that allows us to make statements such as "x^* performs well for a *large portion* of situations δ". In fact, if we accept that various instances of δ present themselves according to a probability distribution $\mathbb{P}$, i.e., that

δ is distributed according to $\mathbb{P}$,

then it is *possible* (at least in principle) to *quantify* the portion of the δ's for which x^* does not perform satisfactorily.

In what follows, the probability that the solution x^* does not perform satisfactorily is denoted with $V(x^*)$, which we call the *risk* of the solution x^*.

The risk $V(x^*)$ is a number between 0 and 1 and is an indicator of the extrinsic quality of x^*. A good enough extrinsic quality can then be formalized by means of a condition of the kind $V(x^*) \leq \bar{\epsilon}$, where $\bar{\epsilon}$ is a domain-dependent user-chosen threshold.

It must be remarked that, from the decision-maker point of view, there is a large gap between *accepting that δ is distributed according to some probability $\mathbb{P}$* and *assuming that such $\mathbb{P}$ is known*. In fact, when δ is the outcome of a complex generation mechanism, the availability of a satisfactory description of $\mathbb{P}$ is often precluded. The theory of the scenario approach recognizes this gap and is built upon the premise that

$\mathbb{P}$ exists but is not known to the decision-maker.

$\mathbb{P}$ manifests itself through:

observations $\delta^{(1)}, \ldots, \delta^{(N)}$, which are modeled as independent draws from $\mathbb{P}$ (according to a standard terminology, observations are independent and identically distributed, "i.i.d.").

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While this i.i.d. assumption is limiting[2], it is worth remarking that many applications can be cast within, or drawn back to, this i.i.d. set-up. For example, stock prices at equispaced time intervals are definitely not independent; however, logarithmic return increments are independent according to the Black-Scholes model, Black & Scholes (1973) (the reader is referred to the box "How to get i.i.d. scenarios in practice" for more general strategies to recast a problem into an i.i.d. framework).

Draws $\delta^{(1)}, \ldots, \delta^{(N)}$ are first-hand knowledge on the problem and the scenario approach provides a well-principled framework to estimate $V(x^*)$ from the data. As we shall see in Section 3, the estimate remains correct even when any partial or insecure knowledge on $\mathbb{P}$ that has been used at the time the decision problem was formulated turns out to be incorrect (this is the requirement in point (b) at the end of Section 2.3). Moreover, the theory is grounded on finite-sample results that are rigorously valid for any sample size (the reader is referred to the box "Asymptotic results in the era of data: the tantalizing horizon" for a discussion on the value of finite-sample results).

How to get i.i.d. scenarios in practice.

In many problems, scenarios are naturally i.i.d.; this is the case for example in all problems where data are draws from a population, with myriad applications in machine learning, prediction and classification. Otherwise, various techniques can be used to draw the problem back to the i.i.d. set-up. Here, two families of strategies of wide applicability to recast non-independent data into an independent sample are briefly touched upon.

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[2]Relaxing the assumption of independence is at present an open and thrilling research endeavor. Clearly, without any assumptions, no theory relating observations collected in the past to situations that shall occur in the future is impossible. On the other hand, the assumption of independence can be seen as extreme and can be easily relaxed e.g. into M-dependence, as is done in Campi et al. (2009a). Along more general routes, one can resort to ϕ-mixing or α-mixing assumptions that are a natural generalization of the condition of independence, see e.g. Bosq (1998) or to conditions that limits the mismatch in a conditional sense from independence, as suggested in Dagan et al. (2019) where the Dobrushin's condition has been used.

- **Prediction-based strategies:** Sometimes, at least a rough predictor for a forthcoming observation is available (e.g., the Weather Bureau provides us with weather predictions). Then, any observation at time t can be decomposed into a prediction part, based on the best of our knowledge until time t, and a prediction error. Often, prediction errors at different time instants are only lightly correlated and the sequence of the prediction errors can, at least in first approximation, be treated as an independent sequence.

- **Segmentation:** The states visited by a Markov Chain do not form an independent sequence, but an observed trajectory can be segmented into independent episodes by exploiting the visit of a recurrent state, restart events after entering an absorbing state, etc., see, e.g., Vidyasagar (2014).

Instead, even when the assumption that data are identically distributed is not overall realistic, still it is often an acceptable approximation over relatively short time windows, or after suitable domain-dependent preprocessing such as the removal of seasonal trends. Some additional discussion on this point is provided in Section 4.

Asymptotic results in the era of data: the tantalizing horizon.

Traditionally, statistics has been dominated by asymptotic results. The usage of these results in practice always introduces approximations and is acceptable only when the number of data points is large compared to the dimension of the solution that is being tuned on the data set. The present era where data are ubiquitous and largely available may seem to have lessened the need for finite-sample results and have favored the usage of asymptotics. However, this evaluation turns out to be incorrect. The reason lies in the fact that in our present times the greater availability of data fares hand in hand with the increasing scale of the problems. As

problems become larger-scale, they require solutions of higher dimension to be satisfactorily resolved so that, as new data become available, the decision-maker is tantalized to resort to more articulated and complex solutions sets. This results in a sort of receding horizon that jeopardizes the use of asymptotic statistics.

2.5. The operation of the scenario approach illustrated on a simple example

This section illustrates, and complements, various aspects touched upon in previous sections of this article. It is meant to provide the reader with a more transparent understanding of the operation of the method before delving into the more technically-oriented presentation of Section 3.

Suppose that the severity of a disease can be quantified by a real number y, and yet an accurate assessment of y requires a medical test that is too invasive to be applied on a vast scale. On the other hand, a simpler inspection delivers a number u that carries information on y and

> our aim is to construct an interval predictor that associates to a given value of u a range of values for y.

To do so, we follow a workflow in line with Figure 1.

Leveraging prior knowledge

We start by collecting the opinions of some experts. Most of them express the educated guess that y should increase linearly with u, while a few of them, with a reputation of being contrarians, are more doubtful and say that they could even expect a negative correlation between u and y for values of u that are above average. None of the experts expects that the dispersion of the values of y changes significantly with u. Moving from this latter observation,

> we set out to construct an interval predictor that maps u into intervals $I(u) = [\varphi(u) - \frac{h}{2}, \varphi(u) + \frac{h}{2}]$ of fixed width h, where both

the function $\varphi(\cdot)$ and the width parameter h have to be suitably designed based on further assessments of the problem.

Next we decide to trust the first group of scientists and take $\varphi(u) = a + b \cdot u$, a linear function of u with a and b tunable parameters. Owing to these choices, we are only left to select three parameters, a, b and h, which form the decision variable $x = (a, b, h)$.

The quality criteria

The indicator of the intrinsic quality of an interval predictor $I(u)$ is the width h: the smaller h the more accurate the prediction. While pursuing this intrinsic quality, we also have to keep control on the extrinsic quality, represented by the reliability of the predictor: we must ensure that the portion of patients for which $y \in I(u)$ is large enough. More formally, letting $V(x) = \mathbb{P}\{y \notin I(u)\}$, where $\mathbb{P}$ is the probability according to which patients (corresponding to pairs (u, y)) distribute, we would like that $V(x^*) \leq \bar{\epsilon}$ for a suitably small $\bar{\epsilon}$.

The scenario approach: let the data speak

Suppose that 100 patients are independently drawn from $\mathbb{P}$ and tested (with both the invasive and the simpler test), and the corresponding 100 scenarios $\delta^{(i)} = (u^{(i)}, y^{(i)}), i = 1, \ldots, 100$, are at our disposal.[3] Based on these scenarios, we choose $x^* = (a^*, b^*, h^*)$ by the following rule.

> RULE: the parameters a^*, b^* and h^* are those that yield the interval predictor with minimum width h subject to the condition that $y^{(i)} \in I(u^{(i)}), i = 1, \ldots, 100$.

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[3] In this simulation example the $\delta^{(i)}$ were artifically generated and, for reproducibility, we inform the reader that $u^{(i)}$ is distributed uniformly on $[0, 1]$ and that the corresponding $y^{(i)}$ is obtained from $u^{(i)}$ via relation $y^{(i)} = (0.21 + 1.4u^{(i)} - [u^{(i)}]^2) + 0.02 \cdot \log((1 + w^{(i)})/(1 - w^{(i)}))$ where $w^{(i)}$ is distributed uniformly on $[-1, 1]$ and independent of $u^{(i)}$; however, the scenario approach does not use this information.
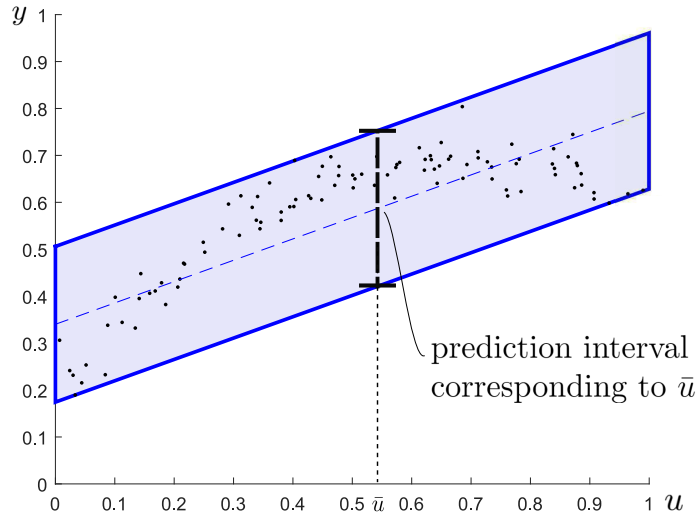
Figure 2: A sample of 100 scenarios: each scenario is a data point in the (u, y) space. The function $\varphi(u) = a^* + b^* \cdot u$ is represented by the dashed line. For each value of u the predicted interval is $I(u) = [\varphi(u) - \frac{h^*}{2}, \varphi(u) + \frac{h^*}{2}]$.

Figure 2 gives the result generated by the RULE using the data that are shown in the same figure.

Assessing the quality of the solution

The intrinsic quality can be assessed from the value of h^* as soon as the predictor is constructed. In the outcome shown in Figure 2, h^* turned out to be only moderately satisfactory (and indeed in this toy example it is even visually clear that the interval predictor includes large portions of empty space).

While h^* is an observable quantity, $V(x^*)$ cannot be directly computed because it depends on probability $\mathbb{P}$, which is not known. An evaluation of $V(x^*)$ can however be performed by using the scenario theory and, to understand the type of results the scenario theory offers, in the next points we first analyze what has been observed in a campaign of simulated examples where $\mathbb{P}$ was artificially

manufactured, and hence it was known.

The distribution of $V(x^*)$

Simulation campaign #1. Our first simulation campaign consists of repeated constructions of the predictor according to the RULE, with different sets of data.

- In $M = 1000$ repetitions, we generated a data set made up of $N = 100$ observations, $(\delta^{(1)}, \ldots, \delta^{(100)})$ according to the same distribution that was used to generate the data points in Figure 2;

- for each one of the 1000 data sets, we constructed a predictor by computing $x^* = (a^*, b^*, h^*)$ according to the RULE;

- for each x^*, we computed $V(x^*)$ (note that we can compute the exact value of $V(x^*)$ because we are running an artificial example where we know the distribution according to which the pairs (u, y) are generated; this would be impossible if data were real data generated from an unknown distribution).

The histogram of the 1000 values of $V(x^*)$ is given in Figure 3. The true distribution of $V(x^*)$ computed analytically is represented by the dashed line in the same figure (the histogram tends to this distribution as $M \to \infty$).

Let us consider the quality threshold $\bar{\epsilon} = 0.11$. In our 1000 simulations, the condition $V(x^*) \leq \bar{\epsilon}$ was always satisfied. By an analytical computation, we found that $\mathbb{P}^{100}\{(\delta^{(1)}, \ldots, \delta^{(100)}) : V(x^*) > \bar{\epsilon}\} = 7.73 \cdot 10^{-4}$ (note that the distribution of $(\delta^{(1)}, \ldots, \delta^{(100)})$ is $\mathbb{P}^{100}$ because scenarios are i.i.d. draws). Since $7.73 \cdot 10^{-4}$ is a small number, it is not surprising that this event did not happen in our 1000 experiments.

The knowledge of the value $7.73 \cdot 10^{-4}$ can be used to make statements like the following one:
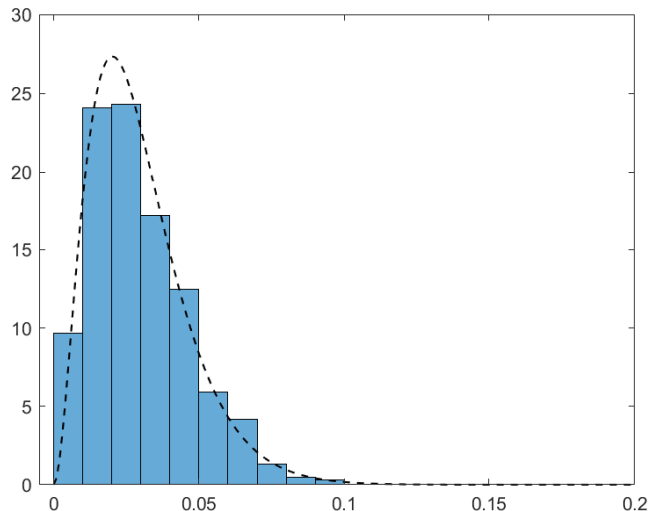
Figure 3: Histogram of $M = 1000$ values of $V(x^*)$. The distribution of $V(x^*)$, to which the histogram tends when $M \to \infty$, is represented by the dashed line.

> if we run an experiment and build a predictor based on 100 scenarios, the resulting predictor will have a risk smaller than 11% with confidence $1 - 7.73 \cdot 10^{-4}$.

A natural question now is how this story changes for a different data generation mechanism, that is, for a different $\mathbb{P}$.

Simulation campaign #2. We made a second simulation campaign as before but, this time, we replaced the probability distribution of (u, y) with a new one: u is uniform over $[0, 1]$ and the distribution of y given u is uniform over $[0.9u, 0.9u + 0.1]$. Unlike the distribution in campaign #1, this distribution fits very well our linear *a priori* belief leading to a small value of h^* (see Figure 4). The histogram of $V(x^*)$ for this second simulation campaign is shown in Figure 5, together with the exact distribution of $V(x^*)$.

Surprisingly, the true distribution is the same as that for campaign #1. This is a symptom of a general fact: any distribution of (u, y) with a density leads exactly to the same distribution of $V(x^*)$. More on the invariant distribution of
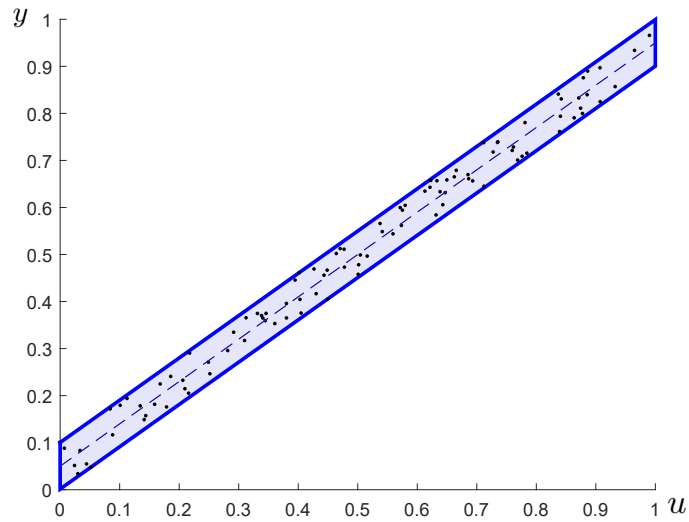
Figure 4: An interval predictor obtained in the second simulation campaign.
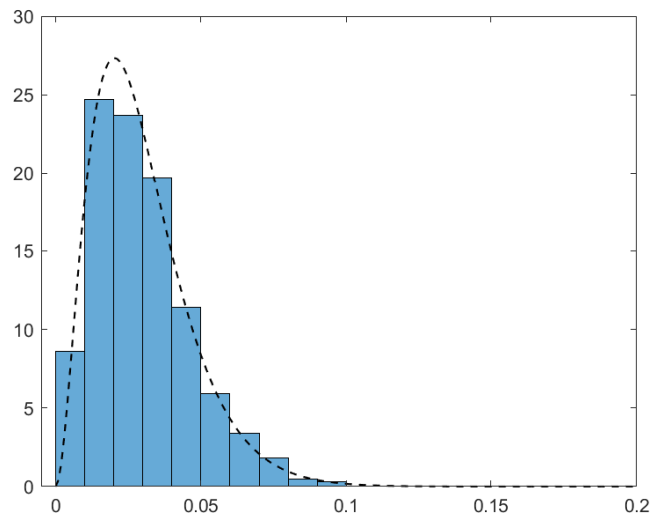

Figure 5: Histogram of $M = 1000$ values of $V(x^*)$ for another distribution of (u, y). The true distribution of $V(x^*)$, to which the histogram tends when $M \rightarrow \infty$, is given by the dashed line and it is exactly the same as that in Figure 3.

$V(x^*)$ will be provided in the next technical Section 3, however, we anticipate that it is a Beta distribution with expected value equal to $\frac{3}{N+1}$. Interestingly, the number 3 at the numerator coincides with the number of optimization variables (a, b, h). This is not a fortuitous circumstance: should we use for example a predictor that has 4 optimization variables (a, b, c, h), $V(x^*)$ would distribute like a Beta with expected value equal to $\frac{4}{N+1}$. In this case, the probability of the event $V(x^*) > 0.11$ would increase to $3.4 \cdot 10^{-3}$.

The separation principle

The above simulation campaigns have revealed a fundamental, and unexpected, property of the RULE: the distribution of $V(x^*)$ remains the same irrespective of the actual mechanism by which data are generated. As a consequence, we can make a judgment on $V(x^*)$ (e.g., $V(x^*) \leq 0.11$ with confidence $1 - 7.73 \cdot 10^{-4}$) and trust it despite the possible incorrectness of the priors which were used in the formulation of our optimization problem. This is a manifestation of the so-called "separation principle".

Going back to our example, in the result in Figure 2 using the prior that u and y are linearly correlated led to a poor result in terms of the width of the prediction interval (which might suggest that our trust in the majority of the experts was misplaced). Nonetheless, the result that $V(x^*) \leq 0.11$ holds with high confidence remains intact. Next, we may want to give a chance to the minority opinion and move to consider a quadratic function $\varphi(u) = a + b \cdot u + c \cdot u^2$ so as to incorporate a possible negative correlation for high values of u. By using again the RULE (extended to the additional parameter c) with the same data set as in Figure 2, we found the much thinner interval predictor in Figure 6 (this is an example of application of point (iv) in Section 2.3). In this latter case, the confidence in the result that $V(x^*) \leq 0.11$ is $1 - 3.4 \cdot 10^{-3}$ (just slightly lower than before as a consequence of having used one more optimization variable).

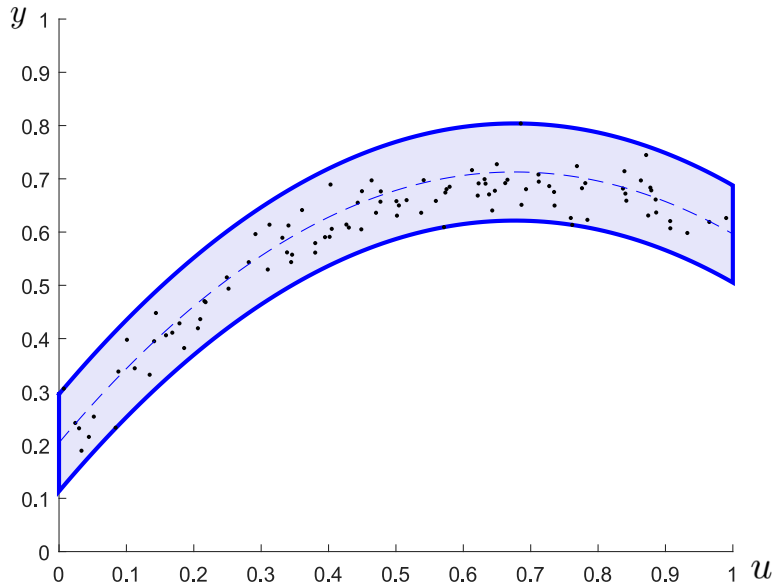*From the simple example to more general problems*

Figure 6: An interval predictor obtained by fitting a function $\varphi(u) = a + b \cdot u + c \cdot u^2$.

Many real life problems differ from the above simple example in some important respects:

- situations are typically described by large dimensional objects and not just (u, y) pairs;

- the impact of prior knowledge cannot always be described as clearly as in this simple example where there is a single decision variable, h, that accounts for the intrinsic quality of the solution, while prior beliefs affect the shape of the predictor through the other decision variables $(a, b, \ldots)$;

- the distribution of $V(x^*)$ is not an invariant as the distribution of data changes.

These differences do not prevent the theory of the scenario approach to be applicable. The last point is of particular interest and we anticipate that the reason why the distribution of $V(x^*)$ was invariant in the example of this section was that the *complexity* (as defined in the next technical section) of the solution was

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the same irrespective of the sample of scenarios at hand. Importantly, when this circumstance does not turn out to be true, universal statements certifying $V(x^*)$ independently of the data generation mechanism are still possible. Moreover, the scenario theory is much more general than what the previous example has shown and includes a rich framework for the evaluation of the risk based on the complexity of the solution when the complexity is not constant, as well as schemes allowing for the removal of some data points with the aim to strike a suitable balance between reliability (extrinsic quality) and accuracy (intrinsic quality).

3. The Theory of the Scenario Approach

To better illustrate the nature of the results, we start from a special but notable set-up (which encompasses as a special case the simple example of Section 2.5) and we will then move gradually towards more general frameworks.

3.1. Convex worst-case optimization

Let $x \in \mathbb{R}^d$ be the decision vector. To make precise the somehow informal description of Section 2.3, let us introduce two real functions $c(x)$ and $f(x, \delta)$ as follows:

- $c(x)$ is a cost function (to be minimized), which is used as a quantifier of the intrinsic quality of x;

- $f(x, \delta)$ models the "regret" for employing x when δ occurs. Condition $f(x, \delta) \leq 0$ indicates a satisfactory performance and, correspondingly, $V(x) = \mathbb{P}\{\delta : f(x, \delta) > 0\}$ is the indicator of the extrinsic quality of x.

In this section, we assume that $\mathcal{X}$ is a convex set; $c(x)$ is convex; and $f(x, \delta)$ is convex in x for all δ (while the dependence of f on δ is arbitrary).[4]

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[4]As we shall see in Section 3.3, recent developments of the scenario approach remove these convexity assumptions.

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The decision process aims at finding the decision with optimal intrinsic quality (i.e., minimum cost $c(x)$) among all the candidate decisions that perform "well" for all the observed scenarios. In formal terms, this leads to the following mathematical program:

$$\min_{x \in \mathcal{X} \subseteq \mathbb{R}^d} c(x)$$
$$\text{subject to: } f(x, \delta^{(i)}) \leq 0, \ i = 1, \ldots, N. \tag{1}$$

As is clear, one primary goal behind using (1) is that of safeguarding against the worst, which motivates enforcing the constraints $f(x, \delta^{(i)}) \leq 0$ for *all* the scenarios. Although suitable to various contexts, this conservative standpoint may result in a poor value of $c(x^*)$ (x^* is the solution to problem (1)), in which case one may want to resort to more flexible alternatives as discussed in Section 3.4.

In the remainder of this Section 3.1 we will assume that, for any N, the feasibility domain of the scenario program (1) has an interior point and that the solution to (1) is unique.[5]

Before proceeding with the description of the theoretical results, it is worth remarking that the example of Section 2.5 fits into this framework of convex worst-case optimization if (a, b, h) is identified with x, and (u, y) with δ. In fact, building a predictor by applying the RULE in Section 2.5 amounts to solving (1) with the cost function $c(x) = h$ and the regret function $f(x, \delta) = |a + b \cdot u - y| - h$; note also that, in that context, the risk $V(x^*)$ is the probability with which a new observation (u, y) falls outside the prediction model of width h^*.[6]

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[5]The uniqueness assumption is introduced to simplify the presentation; if the solution is not unique, a suitable tie-break rule can be introduced, Campi & Garatti (2008); other variations on the theme are possible, see e.g. Calafiore (2010b); Nasir et al. (2016) for contributions about unfeasible problems.

[6]Interval Predictor Models as in Figures 2, 4 and 6 are obtained from convex worst-case optimization problems that have a special structure for which specific theoretical results are available, see in particular Carè et al. (2015) and Carè et al. (2014). For other contributions on Interval Predictor Models, the reader is also referred to Campi et al. (2009a); Calafiore (2010a); Patelli et al. (2013); Crespo et al. (2016); Lacerda et al. (2018); Garatti et al. (2019); Wang et al. (2020).

The scenario theory for convex worst-case optimization

In Campi & Garatti (2008), it is proven that the cumulative distribution function of $V(x^*)$ is (first-order) stochastically dominated by a Beta distribution with parameters $(d, N - d + 1)$. In mathematical terms,

$$\mathbb{P}^N\{V(x^*) \leq v\} \geq 1 - \sum_{i=0}^{d-1} \binom{N}{i} v^i (1-v)^{N-i}. \tag{2}$$

Importantly, this result is *universal*, in the sense that it is valid irrespective of $\mathbb{P}$. An upper-bound $\bar{\epsilon}$ for $V(x^*)$ can then be obtained from (2) as follows. Let β be a small number, say $\beta = 10^{-5}$, and let $\bar{\epsilon}$ be the value that solves the equation

$$\sum_{i=0}^{d-1} \binom{N}{i} v^i (1-v)^{N-i} = \beta, ^7 \tag{3}$$

then $V(x^*) \leq \bar{\epsilon}$ holds with high probability $1 - \beta$. Since the Beta distribution is unimodal and thin tailed, $\bar{\epsilon}$ gets close to its mean (which is $\frac{d}{N+1}$) for a relatively small number of observations. Thus, roughly, $\bar{\epsilon}$ depends on the ratio between d and N.

In applications, $\bar{\epsilon}$ can be used as a relevant information to compare decisions coming from various decision spaces. For instance, in the example of Section 2.5 solving (3) for $\beta = \frac{1}{1000}$ gives $\bar{\epsilon} = 0.108$ when $d = 3$ (linear predictor) and $\bar{\epsilon} = 0.125$ when $d = 4$ (predictor centered around a quadratic function) and these values complement the information coming from the intrinsic quality. The reader is referred to the box "Dealing with multiple comparisons by the union

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[7]This equation can be easily solved by bisection. The MATLAB function betainc-inv(β,d,$N - d + 1$,'upper') solves it and returns directly the value of $\bar{\epsilon}$. An approximate formula that can be useful for a first pencil-and-paper estimation of $\bar{\epsilon}$ is

$$\bar{\epsilon} \leq \frac{1}{N}\left(d - 1 + \ln\frac{1}{\beta} + \sqrt{2(d-1)\ln\frac{1}{\beta}}\right),$$

see Alamo et al. (2015).

bound" for a more detailed discussion on the guarantees that can be attached to a selection made out of competing alternatives.

As already mentioned, in the example of Section 2.5 the distribution of $V(x^*)$ is exactly a Beta distribution with parameters $(d, N - d + 1)$, irrespective of the data generation mechanism, provided that $\mathbb{P}$ has density. Thus, in that case, the inequality "$\leq$" in (2) holds in fact with "$=$", i.e.,

$$\mathbb{P}^N\{V(x^*) \leq v\} = 1 - \sum_{i=0}^{d-1} \binom{N}{i} v^i (1-v)^{N-i} \qquad (4)$$

(the density of this distribution when $N = 100$ and $d = 3$ is the dashed curve in Figures 3 and 5). The existence of problems for which the inequality in (2) is an equality (examples can be found for any d and N) shows that (2) cannot be improved unless it is specialized to subclasses of problems.

The notion of complexity

In the proof of (2) developed in Campi & Garatti (2008), a key role is played by the notion of *complexity*.

> The *complexity* of x^* is an integer $s^* \in \{0, 1, \ldots, N\}$ such that x^* can be obtained by solving a problem similar to (1) that only contains a subsample of scenarios from $\delta^{(1)}, \ldots, \delta^{(N)}$ whose cardinality is s^* while no subsamples of scenarios with cardinality lower than s^* exist that give the same solution x^*.[8]

For an example, we can go back to Figure 2 and note that the same solution (a^*, b^*, h^*) would have been obtained with just 3 points, those that lie on the boundary of the prediction model. Instead, any subsample of scenarios that does not include these 3 points leads to a different solution, so that $s^* = 3$ in

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[8]It is worth noticing that, in the convex set-up, the scenarios that suffice to reconstruct x^* are always a subset of the active constraints, which makes the evaluation of s^* computationally easy.

this case.

The following is a key fact in convex worst-case scenario theory.

Key Fact. *For any problem in the form of* (1), *it holds that* $s^* \leq d$.

Looking at the proof of (2) in Campi & Garatti (2008), one sees that (2) deeply relies on the above Key Fact. Also, the fact that the predictor in Figure 2 turns out to have a complexity s^* that is equal to d $(= 3)$ is not by chance: in Garatti et al. (2019), one can find a proof that this problem belongs to the class of *fully-supported problems*, for which $s^* = d$ happens with probability 1.[9] In Campi & Garatti (2008), it is proven that relation (4) holds for all fully-supported problems and that all other problems are dominated by the fully-supported class in the sense that (2) holds.

Dealing with multiple comparisons by the union bound.

In the prediction problem of Section 2.5, we know that $V(x^*) > \bar{\epsilon}$ happens with probability $\beta = \frac{1}{1000}$, where $\bar{\epsilon} = 0.108$ for $d = 3$ and $\bar{\epsilon} = 0.125$ for $d = 4$. If the selection between $d = 3$ and $d = 4$ is made *a posteriori* (after seeing the prediction interval), we might be advised by an "evil oracle" that indicates a "bad" choice (for which $V(x^*) > \bar{\epsilon}$) whenever one exists. However, no matter how evil the oracle is, the probability that a data set verifies the condition $V(x^*) > \bar{\epsilon}$ for one of the two choices cannot be larger than $\frac{1}{1000} + \frac{1}{1000} = \frac{2}{1000}$. Therefore, the certificate "$V(x^*) \leq \bar{\epsilon}$" is valid with probability at least $1 - 2 \cdot 10^{-3}$. This argument can be extended

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[9]A problem is fully-supported if $s^* = d$ with probability 1 and it is *non-degenerate*; a problem is non-degenerate if, for any N, there is with probability 1 a unique choice of indexes, say $i_1, i_2, \ldots, i_k$ (with $i_1 < i_2 < \ldots < i_k$) from $1, 2, \ldots, N$, such that: (a) problem (1) where only the constraints $f(x, \delta^{(i_j)}) \leq 0$, $j = 1, \ldots, k$, are enforced gives the same solution x^* as with all constraints; (b) if $k > 0$, discarding further indexes from $i_1, i_2, \ldots, i_k$ changes the solution (irreducibility of the set of indexes). (Note that for $k = s^*$ one certainly finds one such set of indexes; the definition of non-degeneracy requires that this set is unique within all subsets of indexes whose cardinality is equal to s^* or larger.) Within convex optimization, degeneracy requires an anomalous accumulation of active constraints to happen and assuming non-degeneracy is therefore reasonable in many applications. *Note for the reader:* the definition of fully-supportedness given in the scenario literature is equivalent to that given here, but sometimes it is formulated differently.

to the case where one chooses a solution from M possibilities leading to the conclusion that $\mathbb{P}^N\{V(x^*) \leq \bar{\epsilon}\} \geq 1 - M\beta$ (where, similarly to the above example with $d = 3$ or $d = 4$, the value of $\bar{\epsilon}$ depends on the choice made). Since small values of β (such as 10^{-7} or 10^{-8}) can be enforced with reasonable sample sizes, the value of $M\beta$ can be easily kept small even when large set of choices are tested out.

3.2. The wait-and-judge perspective

The results in Section 3.1 stand on the observation that the complexity s^* is upper bounded by d (see "Key Fact") and culminate in result (2), which is tight for fully-supported problems, that is, (4) holds. On the other hand, (2) is only an upper bound for non fully-supported problems and the reader is referred to Figure 7 for the distribution of $V(x^*)$ in two non fully-supported examples.
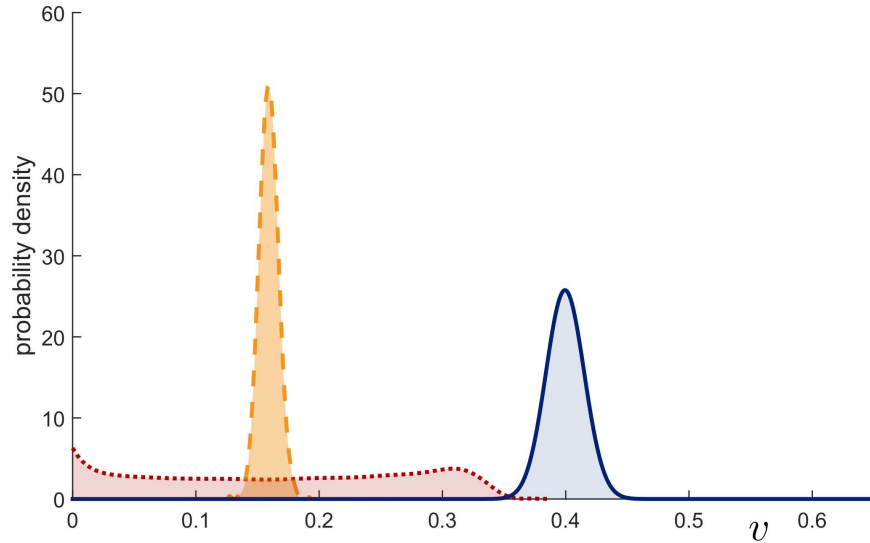

Figure 7: [This figure is taken from Garatti & Campi (2019)] The probability density function of the Beta distribution, as given by the right-hand side of (2) with $N = 1000$ and $d = 400$, is represented in blue continuous line. For a fully-supported problem, $V(x^*)$ distributes as the Beta distribution. The dotted and dashed lines show the distribution of $V(x^*)$ for two examples described in Garatti & Campi (2019): while these densities are dominated by the Beta distribution in the sense of (2), using the Beta distribution to estimate $V(x^*)$ produces conservative evaluations in these two cases.

The conservatism inherent in the bounds that are obtained from the Beta

distribution normally worsens in large scale problems with a high dimensional optimization domain, which is a situation that is encountered with increasing frequency in modern applications.[10]

In what follows, we present results for convex worst-case optimization that establish a connection between $V(x^*)$ and s^* (rather than a connection between $V(x^*)$ and the upper bound d on s^* as in (2)). The use of these results is that one waits until the solution is determined and, at that time, $V(x^*)$ is judged from the complexity s^* that has been measured at the solution (see also the box "Compression and risk: a solid marriage").

Compression and risk: a solid marriage.

The complexity s^* is small when the sample $\delta^{(1)}, \ldots, \delta^{(N)}$ can be slimmed down to a small subsample that is sufficient to reconstruct the solution x^*. Therefore, the complexity is related to the *compressibility* of the sample $\delta^{(1)}, \ldots, \delta^{(N)}$. The idea that the compressibility of the information carried by the data is related to generalization properties is not new in machine learning, see e.g. Rissanen (1978); Littlestone & Warmuth (1986); Rissanen (1986); Ming & Vitányi (1990); Barron et al. (1998); Graepel et al. (2005); Moran & Yehudayoff (2016); Hanneke & Kontorovich (2019). What is new in scenario optimization is the vast applicability of this concept beyond a machine learning context and that compressibility leads to extraordinarily powerful results as in (2), (4). Recent research efforts have been geared towards extending these results to general schemes beyond convex optimization, see e.g. Alamo et al. (2009); Campi (2010); Margellos et al. (2015); Esfahani et al. (2015); Grammatico et al. (2016); Carè et al. (2017); Ramponi & Campi (2018); Carè et al. (2018); Campi et al. (2018); Margellos et al. (2018); Paccagnan & Campi (2019). Part of these results are outlined in

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Sections 3.3, 3.4 and 3.5.

We start by illustrating a simulation example; later, we present the general theory. This section contains results from Campi & Garatti (2018); Garatti & Campi (2019, 2021), to which the reader is referred for the proofs and more details.

Example: orthant that includes random points

A population of points p in a 400-dimensional Euclidean space $\mathbb{R}^{400}$ is distributed according to a probability $\mathbb{P}$. We want to choose an $x \in \mathbb{R}^{400}$ such that $\sum_{j=1}^{400} x_j$ (subscript j denotes component) is minimized (intrinsic criterion) while relation $p_j - x_j \leq 0$, $j = 1, \ldots, 400$, holds with high probability (in other words, the negative orthant with vertex in x contains most of the probabilistic mass of $\mathbb{P}$ – extrinsic criterion).

Worst-case scenario solution

We collected $N = 1000$ points $p^{(1)}, \ldots, p^{(1000)}$ (these are the scenarios), and solved the scenario program:

$$\min_{x \in \mathbb{R}^{400}} \sum_{j=1}^{400} x_j$$
$$\text{subject to: } p_j^{(i)} - x_j \leq 0, \quad j = 1, \ldots, 400 \ \ i = 1, \ldots, 1000, \qquad (5)$$

which is a convex scenario program in the form of (1) with $d = 400$ and $N = 1000$, $c(x) = \sum_{j=1}^{400} x_j$ and $f(x, \delta) = \max_{j=1,\ldots,400}(p_j - x_j)$ (where, clearly, $\delta = p$).

Results of two simulation campaigns

Two simulation campaigns were performed for two different probability distributions $\mathbb{P}_A$ and $\mathbb{P}_B$. For each simulation campaign, we repeated $100\,000$ times

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the sampling of $N = 1000$ scenarios, and computed the corresponding x^* and s^*. Every time, we also computed $V(x^*)$ by exploiting the privilege (due to the fact that we are in a simulated set-up) of knowing the real distribution of the points.
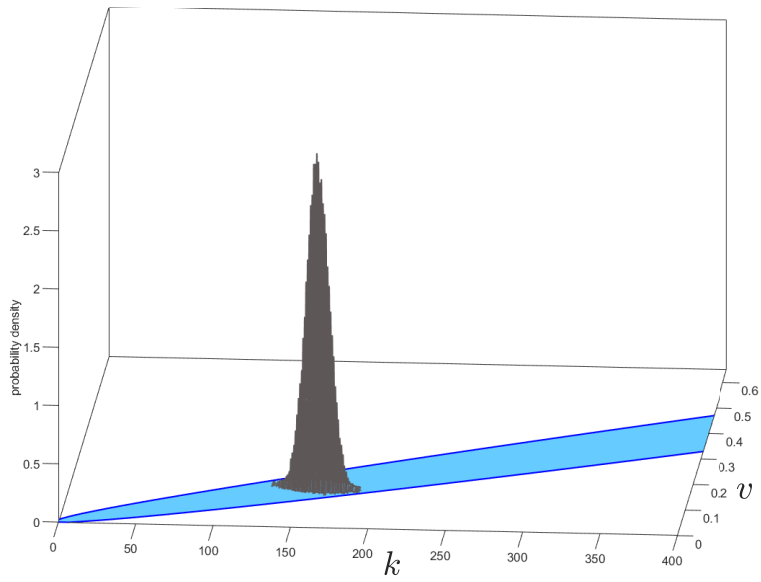

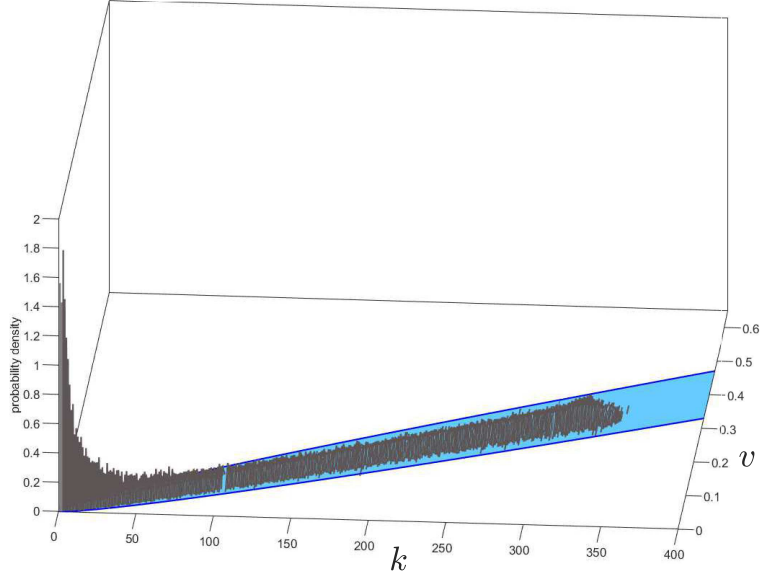Figure 8: Empirical probability distribution of $(s^*, V(x^*))$ when points are generated by $\mathbb{P}_A$.

Figure 9: Empirical probability distribution of $(s^*, V(x^*))$ when points are generated by $\mathbb{P}_B$.

Figure 8 shows the empirical bivariate distribution of the values (k, v) taken by $(s^*, V(x^*))$ over the 100 000 trials in the case of probability $\mathbb{P}_A$. The reader can notice that a blue slanted region is also represented in the (k, v) plane. This region is precisely introduced in the theoretical developments presented below, and, for the time being, we just notice that the distribution of $(s^*, V(x^*))$ appears to be supported on this slanted region. A similar picture for $\mathbb{P}_B$ is given in Figure 9. In this second case, the slanted region is exactly as in the first case and it happens again that the support of the distribution of $(s^*, V(x^*))$ belongs to it.

We next make overt that the marginal distributions of $V(x^*)$ under $\mathbb{P}_A$ and $\mathbb{P}_B$ are those represented in Figure 7 as dashed yellow and dotted red lines respectively. As already discussed in Section 3.2, the Beta distribution of Figure 7 only sets an upper limit to these marginals. The present simulations in Figures 8 and 9 suggest that more information can be gained from the lens of a bivariate point of view where one variable, the risk, is estimated from the other, the com-

plexity (which is a measurable quantity). We anticipate that this result is true in high generality and that the precision in the evaluation of $V(x^*)$ that can be achieved thanks to this new lens is comparable to the precision that comes in fully-supported problems from using relation (4), indeed a remarkable finding.

A general result

As is clear from our simulations, the distribution of the pair $(s^*, V(x^*))$ can take various forms. However, in the two examples that we have just seen this distribution was confined in the slanted region which, for easy reference, is again displayed in Figure 10.
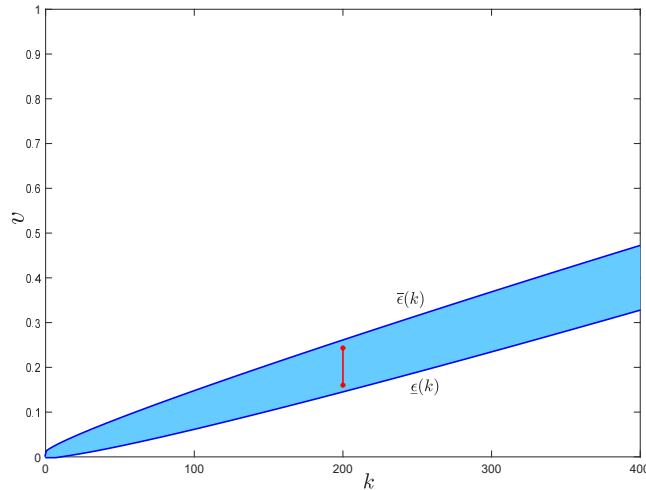

Figure 10: The slanted blue area is the region to which at least $1 - \beta = 99.9\%$ of the probabilistic mass of $(s^*, V(x^*))$ belongs for any non-degenerate convex scenario problem (as defined in Footnote 9) in the form of (1) when $d = 400$ and $N = 1000$. Moreover, for any problem in the form of (1) (degenerate or non-degenerate) at least $1 - \beta = 99.9\%$ of the probabilistic mass of $(s^*, V(x^*))$ lies below the upper boundary of the blue area. For the sake of comparison, in the figure a red interval is represented which is a 99.9% confidence interval of the risk for a fully-supported problem in dimension 200 (the interval has been obtained by using the Beta distribution).

The rule by which the boundaries of the slanted region are constructed is as follows.

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Rule to compute the boundaries of the slanted region

Assume $N > d$. Given a confidence parameter $\beta \in (0, 1)$, for any $k = 0, \ldots, d$ consider the polynomial equation in the t variable

$$\binom{N}{k} t^{N-k} - \frac{\beta}{2N} \sum_{i=k}^{N-1} \binom{i}{k} t^{i-k} - \frac{\beta}{6N} \sum_{i=N+1}^{4N} \binom{i}{k} t^{i-k} = 0. \quad (6)$$

For any $k = 0, 1, \ldots, d$, equation (6) has exactly two solutions in $[0, +\infty)$, which we denote with $\underline{t}(k)$ and $\bar{t}(k)$ ($\underline{t}(k) \le \bar{t}(k)$). Define $\underline{\epsilon}(k) := \max\{0, 1 - \bar{t}(k)\}$ and $\bar{\epsilon}(k) := 1 - \underline{t}(k)$, $k = 0, \ldots, d$. Function $\underline{\epsilon}(k)$ is the lower boundary of the slanted region and $\bar{\epsilon}(k)$ its upper boundary.

A MATLAB code for computing the values of $\underline{\epsilon}(k)$ and $\bar{\epsilon}(k)$ according to this rule is available in the Appendix A of Garatti & Campi (2019).

We now have the following result (it is assumed that the solution x^* exists and is unique, possibly after the use of a suitable tie-break rule).

Wait-and-judge result for convex scenario optimization. *For any non-degenerate problem (as defined in Footnote 9) in the form of* (1), *it holds that*

$$\mathbb{P}^N\{\underline{\epsilon}(s^*) \le V(x^*) \le \bar{\epsilon}(s^*)\} \ge 1 - \beta; \quad (7)$$

moreover, for any problem in the form of (1), *it holds that*

$$\mathbb{P}^N\{V(x^*) \le \bar{\epsilon}(s^*)\} \ge 1 - \beta. \quad (8)$$

Hence, referring to Figure 10, in non-degenerate problems the distribution of $(s^*, V(x^*))$ is confined to the blue slanted region (but a small portion whose probability is no more than β) whereas in degenerate problems the distribution of $(s^*, V(x^*))$ can expand below the lower boundary of the slanted region, while the upper boundary, that sets a limit to $V(x^*)$, is always valid (this latter result is proven in the recent paper Garatti & Campi (2021)).

For the use of this result, the crucial fact to remark is that the quantity on the horizontal axis (complexity) is measurable, while the vertical axis corresponds to the value of the risk $V(x^*)$ that, in real life, is hidden to the decision-maker and can only be estimated. Thus, from Figure 10, one obtains a rule to bound $V(x^*)$ based on the observable s^* and the result in (7) provides sample-dependent bounds of the kind $\underline{\epsilon}(s^*) \leq V(x^*) \leq \bar{\epsilon}(s^*)$ (or just $V(x^*) \leq \bar{\epsilon}(s^*)$, in degenerate cases thanks to (8)) that are valid with high probability $1 - \beta$. The interval $[\underline{\epsilon}(k), \bar{\epsilon}(k)]$ for a given k is comparable to the one that can be generated when working with a fully-supported problem in dimension $d = k$. One of these intervals with $d = 200$, is represented in Figure 10. The quantitative similarity between $[\underline{\epsilon}(k), \bar{\epsilon}(k)]$ and the intervals in the fully-supported case is quite a remarkable fact and reveals the value of the information conveyed by the complexity. From results in Campi & Garatti (2018); Garatti & Campi (2019), one also sees that β impacts on $\underline{\epsilon}(k)$ and $\bar{\epsilon}(k)$ logarithmically so that taking very small values of β enlarges only marginally the interval $[\underline{\epsilon}(k), \bar{\epsilon}(k)]$; moreover, for any k,

$$\bar{\epsilon}(k) - \underline{\epsilon}(k) \to 0$$

as $N \to \infty$ in such a way that the interval shrinks around k/N as N grows, see Campi & Garatti (2020).

3.3. Non-convex worst-case optimization

When the assumption of convexity is dropped, the bound $s^* \leq d$ loses validity. For example, let $x = (x_1, x_2) \in [-1, 1]^2$, $c(x) = x_2$, $\delta \in [-1, 1]$ and $f(x, \delta) = -|x_1 - \delta| - x_2$. An instance of the corresponding scenario program is pictured in Figure 11, and the reader can easily check that any subsample of the 6 observed scenarios yields a solution x^* different from the one obtained with all the scenarios, so that $s^* = 6 > 2 = d$ in this case. Moreover, the same example reveals that a constraint need not be active to be necessary to reconstruct the solution in the non-convex case. This circumstance suggests that degeneracy

becomes quite a common circumstance (the reader may want to think of the case in which one more constraint is added in Figure 11: degeneracy becomes evident whenever this new constraint "shields" the global minimum that opens up after removing one of the 6 constraints already in the figure).
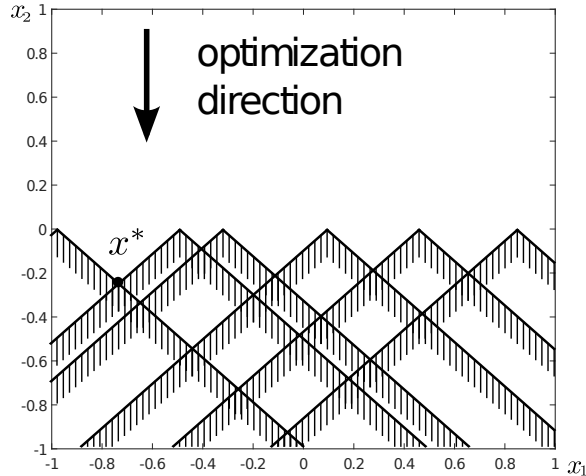

Figure 11: An instance of a non-convex scenario program where $d = 2$ but $s^* = 6$.

Remarkably, the results in the recent paper Garatti & Campi (2021) do not require that $s^* \leq d$ and are applicable even in the degenerate case, leading to the following result (x^* is assumed to exist and to be unique, possibly after the use of a suitable tie-break rule).

Wait-and-judge result for non-convex scenario optimization. *Use equation (6) to compute $\bar{\epsilon}(k)$ for $k = 0, 1, \ldots, N$ (note that the range of k is here extended till N).*[11] *Then, for any problem in the form of (1), where $c(x)$ and $f(x, \delta)$ need not be convex in x and $\mathcal{X}$ is any set, it holds that*

$$\mathbb{P}^N\{V(x^*) \leq \bar{\epsilon}(s^*)\} \geq 1 - \beta. \tag{9}$$

---

[11]The construction for $k = 0, 1, \ldots, N - 1$ is exactly as indicated in the "Rule to compute the boundaries of the slanted region"; for $k = N$, however, equation (6) has only one solution $\bar{t}(N)$, and one defines $\underline{t}(N) = 0$, so that $\bar{\epsilon}(N) = 1$.

The curve $\bar{\epsilon}(k)$ for $N = 1000$ and $\beta = 0.1\%$ is displayed in Figure 12. Note that, for k in the range $[0, 400]$, $\bar{\epsilon}(k)$ is exactly as in Figure 10.
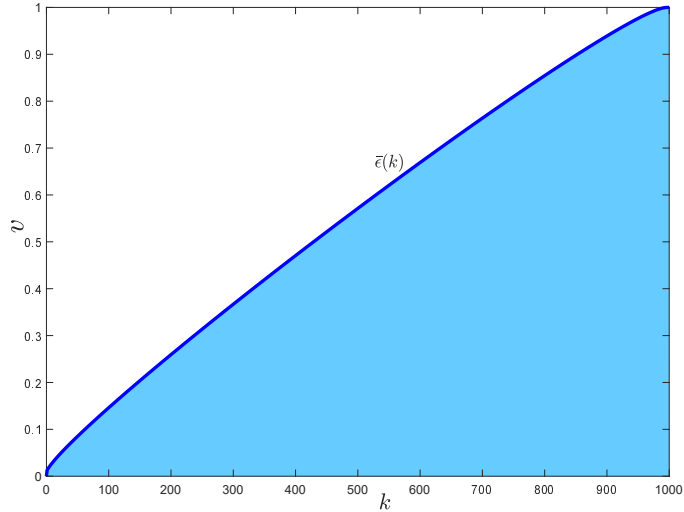

Figure 12: The blue area is the region to which at least $1-\beta = 99.9\%$ of the probabilistic mass of $(s^*, V(x^*))$ belongs for any non-convex scenario problem in the form of (1) when $d = 400$ and $N = 1000$.

3.4. Tuning of the extrinsic quality

In (1), all scenario constraints are rigidly enforced, so expressing an attitude to safeguard against the worst. This, however, can lead to conservative designs: the presence of just one ill scenario can forbid a whole set of candidate solutions and confine the choice to solutions with poor intrinsic quality $c(x)$ (see, e.g., Shapiro et al. (2009); Ramponi (2018); Assif et al. (2020)). On the other hand, in many applications, $c(x)$ and $V(x)$ are seen as conflicting objectives and in this section we focus on a class of decision schemes that allow the decision-maker to compromise between intrinsic and extrinsic quality.

According to the interpretation of $f(x, \delta)$ provided in Section 3.1, a violation of the condition $f(x, \delta^{(i)}) \leq 0$ yields a regret. The first scheme we present relaxes the rigid enforcement of constraints by minimizing a weighted combination of the cost $c(x)$ and the sum of the (positive) regrets $\sum_{i=1}^{N} \left[f(x, \delta^{(i)}) \right]^+$ ($[\cdot]^+$ is positive part, that is, it returns its argument when it is positive and zero otherwise).

This yields the following optimization program where the weight α assigned to the regrets is a tunable trade-off parameter:[12]

$$\min_{\substack{x \in \mathcal{X} \subseteq \mathbb{R}^d, \\ \xi_i \geq 0, i=1,\ldots,N}} c(x) + \alpha \sum_{i=1}^{N} \xi_i$$

$$\text{subject to: } f(x, \delta^{(i)}) \leq \xi_i \ \ i = 1, \ldots, N. \tag{10}$$

Let $(x_\alpha^*, \xi_{i,\alpha}^*)$ be the solution to (10) (assumed to exist and to be unique, possibly after the use of a suitable tie-break rule). In this context, a suitable generalization of the concept of complexity allows one to estimate the extrinsic quality of x_α^* (i.e., quantity $V(x_\alpha^*) = \mathbb{P}\{\delta : f(x_\alpha^*, \delta) > 0\}$).[13] The *generalized complexity* s_α^* is the cardinality of a subsample S of scenarios from $\delta^{(1)}, \ldots, \delta^{(N)}$ defined as follows: S contains all the violated scenarios (i.e., the scenarios for which $f(x_\alpha^*, \delta^{(i)}) > 0$)[14] plus a minimum subsample of the remaining scenarios such that the x-component of the solution to (10) with only S in place remains equal to x_α^*. Then, exactly the same result as in "Wait-and-judge result for non-convex optimization" holds (without requiring any convexity assumption) with the only warning that x^* and s^* must be replaced in the present context by x_α^* and s_α^*.[15] The value of α (which ranges from $\alpha = 0$, corresponding to minimizing $c(x)$ with no concern for the observed scenarios, to $\alpha = \infty$, corresponding to the worst-case approach) influences s_α^* (and, thereby, the bounds on $V(x_\alpha^*)$) and $c(x_\alpha^*)$ and represents a tuning knob in the hands of the decision-maker. By plotting $c(x_{\alpha_1}^*), c(x_{\alpha_2}^*), \ldots$ against the bounds $\bar{\epsilon}(s_{\alpha_1}^*), \bar{\epsilon}(s_{\alpha_2}^*), \ldots$ for $V(x_{\alpha_1}^*), V(x_{\alpha_2}^*), \ldots,$

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[12]The set of all the optimal decisions x_α^* obtained from (10) as a function of α is known as the Pareto frontier of the multi-objective problem of minimizing $c(x)$ and $\sum_{i=1}^{N} \left[f(x, \delta^{(i)})\right]^+$.

[13]The need for a generalization of the concept of complexity can be recognized by noting that problem (10) is not directly in the form of (1), in particular the number of optimization variables x, ξ_i increases with N.

[14]The subsample S is with repetitions, i.e., if two scenarios $\delta^{(i)}$ and $\delta^{(j)}$ for which $f(x_\alpha^*, \delta^{(i)}) > 0$ and $f(x_\alpha^*, \delta^{(j)}) > 0$ turn out to be coincident ($\delta^{(i)} = \delta^{(j)}$), then they both appear in S.

[15]Also in this context a concept of non-degeneracy can be applied leading to the stronger result that $V(x_\alpha^*)$ is both upper and lower bounded, see Garatti & Campi (2019).

one obtains the so-called *cost-risk plot* (see Garatti & Campi (2019) for an example) by which the user can perform a selection of a suitable value for α.

Program (10) is not the only scheme to tune the intrinsic vs. the extrinsic quality. Alternatively, one can discard some of the constraints from the worst-case program (1) and the reader is referred to the papers Campi & Garatti (2011); Garatti & Campi (2013); Picallo & Dörfler (2019); Romao et al. (2020) for studies in this direction. Interestingly, a very precise connection can be traced between the approach of (10) and the discarding strategy. Suppose that $f(x, \delta)$ in (10) can only take the value 0 or 1 (indicator function). Then, as parameter α declines, the solution to (10) explores the optimal solutions that violate an increasing number of scenario constraints, which is the same as computing the optimal solution allowing for the discarding of a certain amount of the constraints.[16] One should however note that solving exactly problem (10) with indicator functions $f(x, \delta)$ is extremely complex even when the cost function $c(x)$ and the sets where $f(x, \delta) = 0$ are convex. The reason is that the indicator functions lift the problem outside the domain of convex optimization. For this reason, the literature on constraint discarding has evolved towards various heuristics covering sub-optimal discarding strategies, e.g., according to greedy procedures. Remarkably, the theoretical results about the risk provided in the references given above remain rigorously valid when these heuristics are used (even though these results do not always provide sharp evaluations). Going back to the approach of (10), one should note that it offers an extra level of flexibility that can accommodate various needs: (i) by the choice of $f(x, \delta)$ one can reflect problem-dependent concepts of regret; (ii) if function $f(x, \delta)$ is selected to be convex in x for any δ, then the problem does not lift outside the domain of convex optimization so that optimal solutions can be easily computed, and the sharp generalization theory touched upon in this section can be applied.

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[16]The optimal solution attains a cost $c(x)$ that has been compared to the optimal cost of a chance-constrained problem in Thorem 6.1 of Campi & Garatti (2011).

3.5. A general theory for decision-making

The scenario theory carries over to an abstract decision-making framework that encompasses all previous set-ups and many others as special cases. Here, we briefly summarize the results in Garatti & Campi (2019, 2021), to which the reader is referred for details.

For any $N = 0, 1, 2, \ldots$ let M_N be a map from any set of scenarios $\delta^{(1)}, \ldots, \delta^{(N)}$ to a decision $z^* \in \mathcal{Z}$, where $\mathcal{Z}$ is a generic decision set. The maps $M_0, M_1, \ldots$ describe the rule according to which decisions are made based on observations in an integrated data-driven set-up as in Figure 1. In order to evaluate whether the extrinsic criterion is met, there is a rule to decide whether a decision z performs "well" when a situation δ occurs. This is expressed in mathematical terms by saying that, to any situation δ, there is associated a set $\mathcal{Z}_\delta \subseteq \mathcal{Z}$ which models the set of the decisions that perform well for δ. The risk of a decision $z \in \mathcal{Z}$ is then defined as $V(z) := \mathbb{P}\{\delta : z \notin \mathcal{Z}_\delta\}$.

No limiting assumptions on the domain of δ and on the map between δ and $\mathcal{Z}_\delta$ are necessary. The freedom of the decision-maker in choosing the maps $M_0, M_1, \ldots$ is also vast as long as the following three properties are satisfied:

- **Permutation invariance**:

  for every N, every $\delta^{(1)}, \ldots, \delta^{(N)}$ and every permutation $(i_1, \ldots, i_N)$ of $(1, \ldots, N)$ it holds that

$$M_N(\delta^{(1)}, \ldots, \delta^{(N)}) = M_N(\delta^{(i_1)}, \ldots, \delta^{(i_N)}).$$

- **Stability in the case of confirmation**:

  for every integers N and n, if $\delta^{(1)}, \ldots, \delta^{(N)}, \delta^{(N+1)}, \ldots, \delta^{(N+n)}$ are such that

$$\forall i \in \{1, \ldots, n\} : \quad M_N(\delta^{(1)}, \ldots, \delta^{(N)}) \in \mathcal{Z}_{\delta^{(N+i)}},$$

  then

$$M_N(\delta^{(1)}, \ldots, \delta^{(N)}) = M_{N+n}(\delta^{(1)}, \ldots, \delta^{(N+n)}).$$

- **Responsiveness to contradiction**:

  for every integers N and n, if $\delta^{(1)}, \ldots, \delta^{(N)}, \delta^{(N+1)}, \ldots, \delta^{(N+n)}$ are such that

  $$\exists i \in \{1, \ldots, n\} : \quad M_N(\delta^{(1)}, \ldots, \delta^{(N)}) \notin \mathcal{Z}_{\delta^{(N+i)}},$$

  then

  $$M_{N+n}(\delta^{(1)}, \ldots, \delta^{(N+n)}) \neq M_N(\delta^{(1)}, \ldots, \delta^{(N)}).$$

Note that the properties of M_N as written above *do not* imply that $z^* \in \mathcal{Z}_{\delta^{(i)}}$, $i = 1, \ldots, N$.

Complexity and degeneracy

Given the scenarios $\delta^{(1)}, \ldots, \delta^{(N)}$, the *complexity* s^* of the decision $z^* = M_N(\delta^{(1)}, \ldots, \delta^{(N)})$ is the cardinality of a minimum subsample S (with repetitions) of scenarios from $\delta^{(1)}, \ldots, \delta^{(N)}$ such that $M_{|S|}(S) = z^*$. The decision-scheme is said to be *non-degenerate* if, for every N, there is with probability 1 a unique choice of indexes, say $i_1, i_2, \ldots, i_k$ (with $i_1 < i_2 < \ldots < i_k$) from $1, 2, \ldots, N$, such that $M_k(\delta^{(i_1)}, \ldots, \delta^{(i_k)}) = M_N(\delta^{(1)}, \ldots, \delta^{(N)})$ while, for $k > 0$, discarding further indexes from $i_1, i_2, \ldots, i_k$ changes the solution.

Guarantees

Similarly to "Wait-and-judge result for non-convex optimization" in Section 3.3, using (6) one computes $\bar{\epsilon}(k)$ for $k = 0, 1, \ldots, N$. Then, equation (9) applies where x^* is replaced in the present context by z^*, while s^* and $V(z^*)$ have to be interpreted according to the definitions of this section. Moreover, under non-degeneracy, one can also compute $\underline{\epsilon}(k)$, $k = 0, 1, \ldots, N$, by extending till $k = N$ the rule for $\underline{\epsilon}(k)$, $k = 0, 1, \ldots, d$, in "Wait-and-judge result for convex optimization" in Section 3.2 and equation (7) holds where again x^* has to be replaced by z^*, while s^* and $V(z^*)$ have to be interpreted according to the def-

initions of this section.

A historical remark

The "responsiveness to contradiction" property was a key property in important generalization results in the history of statistical learning, Littlestone & Warmuth (1986). This property, alone, is sufficient to derive various results within the scenario theory that are applicable to a wide range of decision schemes, Campi et al. (2018). On the other hand, the "stability in the case of confirmation" property plays a special role in obtaining tight bounds (and, under non-degeneracy, small ranges $[\underline{\epsilon}(k), \bar{\epsilon}(k)]$) and, when satisfied, confers considerable added value to the scenario approach. Being naturally satisfied in optimization problems, this property can be recognized as one of the main "secrets" behind the success of the scenario approach in many applications.

4. Conclusions and vision for the future

The increasing availability of data is a resource both for the real-time operation of human-made machines as well as for their off-line design. Although using data can be seen as a heuristic, we are witnessing at various levels attempts to make such heuristics more theoretically-supported, by endowing various techniques with specific guarantees that certify their correct functioning. This endeavor is particularly important in relation to automated systems where machines are required to operate autonomously, or with marginal human intervention, at times in safety-critical conditions.

In this article, we have presented the scenario approach, a general methodology that has come to the fore in recent years for its ability to make designs supported by precise, distribution-free, guarantees. Our main goal with this publication was that of providing an easy access point to various aspects of the scenario methodology, by also positioning its features in relation to current, pressing, challenges posed by the technology of data-driven methods. Numer-

ous are the application contexts where the scenario approach has already proven useful and – while we have decided to leave out of this article detailed presentations – the interested reader has been referred to a list of publications in the box "When should we consider using the scenario approach?". Certainly, this should not give the impression that the scenario approach is a ready methodology for all seasons, and it is common experience that the scenario theory must be adjusted, and at times expanded, to accommodate application needs. We feel appropriate to discuss here a specific field of current, active, research so as to provide a concrete example.

When a human being is found in the condition of cardiac arrest, time is a crucial variable and any hesitation reduces the probability of survival. On the other hand, it is hard for a rescuer with little or no medical training to decide how to act.[17] Hence, the need to provide automatic methods to direct the rescuer and assist her/his decisions. The scenario approach has been applied to this problem with the aim of designing classifiers that act on short ECG (Electro-Cardio-Graphic) traces[18] to determine if an immediate defibrillation shock will be effective to restore spontaneous circulation. In this context, it is important to distinguish between two types of error: erroneously predicting that the defibrillation *will* or *will not* be successful (with very different consequences in practice), and to keep under control their respective probabilities.[19] However, the separation of two types of risks is not germane to the scenario approach, and its use for the defibrillation problem has therefore prompted non-trivial theoretical advances, as discussed in Carè et al. (2018). Still, no claim of final

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[17]Even when a defibrillator is available, there are alternatives to immediately administrating an electrical shock, the most important of which is to give the patient a cardiac massage: a properly given cardiac massage boosts the chances of resuscitation when the defibrillator is subsequently applied. The existence of alternative approaches to shocking has been emphasized by the European Resuscitation Council guidelines: "It is possible to predict, with varying reliability, the success of defibrillation from the fibrillation waveform. [...] It should be possible to prevent the delivery of unsuccessful high energy shocks and minimize myocardial injury", Soar et al. (2021).

[18]All modern defibrillators serve the dual function of shocking and collecting ECGs.

[19]The commonly used terminology *specificity* and *sensitivity* refers to the complement to 1 of these two probabilities of error, i.e., to the corresponding probabilities of success.

success can be pronounced: for this problem precise limits to the specificity and the sensitivity (50% and 95%, respectively, see Neurauter et al. (2007)) need to be guaranteed before a classifier can be deployed on the territory, a performance result that requires more research to be achieved.

The above is just one among many examples where the scenario theory has been adjusted, and expanded in scope, to cover the specific needs of applications. Indeed, accommodating application requirements has been a major engine through the years that has fostered a continuing growth of the scenario theory. Despite all this, the unexplored territory remains vast and we should note that at present the scenario theory is articulated around a paradigm that presents some structural limitations. Two aspects that we deem of primary importance are discussed here, in the hope that this can also stimulate further research.

(a) Problems with partial knowledge.

Throughout this article, and indeed everywhere in the scenario literature, it has been assumed that all uncertain elements are encapsulated in δ, so that, after a scenario has been collected, one acquires a full picture for the particular situation referring to that δ and can ascertain the ensuing performance achieved by any hypothetical decision x. To make a concrete example, in the stock market application of Pagnoncelli et al. (2012); Ramponi & Campi (2018); Arici et al. (2021), δ contains the rate-of-return of all assets in the portfolio, so that, given a scenario referring to the past, the overall return of any possible diversified investment can be computed. While this assumption is completely plausible for this application, still suppose for the sake of the argument that the investor is only provided with the overall return of his own investment (that associated to the diversification that has been actually adopted to buy assets) at the end of each period of investment (from this information, clearly, one cannot figure out the rate-of-return of each single asset) and has to decide how to proceed on investing in the next period. This set-up introduces new, non-trivial, challenges: since no full picture is acquired from the data, one is facing the need to achieve the dual effect of exploring what is unknown while exploiting what

is known. In turn, since now knowledge depends on actions, this rules out the assumption of independence that is inherent in the current formulation of the scenario approach. We believe that studying *active learning schemes* within the scenario approach is of primary importance for various endeavors in control and should rank high in the research agenda.

(b) Problems with varying distribution.

An assumption the scenario approach rests on is that the distribution by which scenarios are drawn remains the same through trials. Often, however, this assumption is only approximately satisfied or holds for a limited amount of time. For example, referring again to an investment problem, the assumption that the rate-of-returns have invariant distribution may be valid only within a restricted time horizon. Moreover, different distributions may be encountered at the time of making a decision and at the later time when the decision is applied, owing to circumstances that are not due to time-variability of the environment but rather to different mechanisms of exposition to observations in the two phases of learning and testing. For example, the emerging field of adversarial machine learning, see, e.g., the survey paper Ortiz-Jiménez et al. (2021), addresses problems of this nature: the input distribution can be varied, even maliciously, by an opponent to mislead the trained classifier. One challenging and important topic within the scenario methodology is to accommodate varying distributions. The research on this topic may proceed at least in two directions: (i) providing guarantees for existing schemes, e.g., worst-case, that remain intact under limited variations of the distribution, see, e.g., Erdogan & Iyengar (2006) for an early contribution in this direction; (ii) designing scenario-based techniques that naturally incorporate robustness features against variability of the distribution while preserving the guarantees on the risk.

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