Title:	Introduction: Sensitivity Analysis
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Introduction: Sensitivity Analysis

Abstract

Sensitivity analysis provides users of mathematical and simulation models with tools to appreciate the dependency of the model output from model input, and to investigate how important is each model input in determining its output. All application areas are concerned, from theoretical physics to engineering and socio-economics. This introductory paper provides the sensitivity analysis aims and objectives in order to explain the composition of the overall "Sensitivity Analysis" chapter of the Springer Handbook. It also describes the basic principles of sensitivity analysis, some classification grids to understand the application ranges of each method, a useful software package and the notations used in the chapter papers. This section also offers a succinct description of sensitivity auditing, a new discipline that tests the entire inferential chain including model development, implicit assumptions and normative issues, and which is recommended when the inference provided by the model needs to feed into a regulatory or policy process. For the "Sensitivity Analysis" chapter, in addition to this introduction, eight papers have been written by around twenty practitioners from different fields of application. They cover the most widely used methods for this subject: the deterministic methods as the local sensitivity analysis, the experimental design strategies, the sampling-based and variance-based methods developed from the 1980s and the new importance measures and metamodel-based techniques established and studied since the 2000s. In each paper, toy examples or industrial applications illustrate their relevance and usefulness.

Keywords: Computer Experiments, Uncertainty Analysis, Sensitivity Analysis, Sensitivity Auditing, Risk Assessment, Impact Assessment

Introduction

In many fields such as environmental risk assessment, behavior of agronomic systems, structural reliability or operational safety, mathematical models are used for simulation, when experiments are too expensive or impracticable, and for prediction. Models are also used for uncertainty quantification and sensitivity analysis studies. Complex computer models calculate several output values (scalars or functions) that can depend on a high number of input parameters and physical variables. Some of these input parameters and variables may be unknown, unspecified, or defined with a large imprecision range. Inputs include engineering or operating variables, variables that describe field conditions, and variables that include unknown or partially known model parameters. In this context, the investigation of computer code experiments remains an important challenge.

This computer code exploration process is the main purpose of the Sensitivity Analysis (SA) process. SA allows the study of how uncertainty in the output of a model can be apportioned to different sources of uncertainty in the model input [51]. It may be used to determine the input variables that contribute the most to an output behavior, and the non-influential inputs, or to ascertain some interaction effects within the model. The SA process entails the computation and analysis of the so-called sensitivity or importance indices of the input variables with respect to a given quantity of interest in the model output. Importance measures of each uncertain input variable on the response variability provide a deeper understanding of the modeling in order to reduce the response uncertainties in the most effective way [57], [30], [23]. For instance, putting more efforts on knowledge of influential inputs will reduce their uncertainties. The underlying goals for SA are model calibration, model validation and assisting with the decision making process. This chapter is for engineers, researchers and students who wish to apply SA techniques in any scientific field (physics, engineering, socioeconomics, environmental studies, astronomy, etc.).

Several textbooks and specialist works [56], [3], [10], [12], [9], [59], [21], [11], [2] have covered most of the classic SA methods and objectives. In parallel, a scientific conference called SAMO ("Sensitivity Analysis on Model Output") has been organized every three years since 1995 and extensively covers SA related subjects. Works presented at the different SAMO conferences can be found in their proceedings and several special issues published in international journals (mainly in "Reliability Engineering and System Safety").

The main goal of this chapter is to provide an overview of classic and advanced SA methods, as none of the referenced works have reported all the concepts and methods in one single document. Researchers and engineers will find this document to be an up-to-date report on SA as it currently stands, although this scientific field remains very active in terms of new developments. The present chapter is only a snapshot in time and only covers well-established methods.

The next section of this paper provides the SA basic principles, including elementary graphic methods. In the third section, the SA methods contained in the chapter are described using a classication grid, together with the main mathematical notations of the chapter papers. Then, the SA-specialized packages developed in the R software environment are discussed. To finish this introductory paper, a process for the sensitivity auditing of models in a policy context is discussed, by providing seven rules that extend the use of SA. As discussed in Saltelli et al [61], SA, mandated by existing guidelines as a good practice to use in conjunction with mathematical modeling, is insufficient to ensure quality in the treatment of scientific uncertainty for policy purposes. Finally, the concluding section lists some important and recent research works that could not be covered in the present chapter.

Basic principles of sensitivity analysis

The first historical approach to SA is known as the local approach. The impact of small input perturbations on the model output is studied. These small perturbations occur around nominal values (the mean of a random variable, for instance). This deterministic approach consists of calculating or estimating the partial derivatives of the model at a specific point of the input variable space [68]. The use of adjoint-based methods allows models with a large number of input variables to be processed. Such approaches are particularly well-suited to tackling uncertainty analysis, SA and data assimilation problems in environmental systems such as those in climatology, oceanography, hydrogeology, etc. [3], [48], [4]. To overcome the limitations of local methods (linearity and normality assumptions, local variations), another class of methods has been developed in a statistical framework. In contrast to local SA, which studies how small variations in inputs around a given value change the value of the output, global sensitivity analysis ("global" in opposition to the local analysis) does not distinguish any initial set of model input values, but considers the numerical model in the entire domain of possible input parameter variations [57]. Thus, the global SA is an instrument used to study a mathematical model as a whole rather than one of its solution around parameters specific values.

Numerical model users and modelers have shown high interest in these global tools that take full advantage of the development of computing equipment and numerical methods (see Helton [20], de Rocquigny et al [9] and [11] for industrial and environmental applications). Saltelli et al [58] emphasized the need to specify clearly the objectives of a study before performing an SA. These objectives may include:

- The factors prioritization setting, which aims at identifying the most important factors. The most important factor is the one that, if fixed, would lead to the greatest reduction in the uncertainty of the output;
- The factors fixing setting, which aims at reducing the number of uncertain inputs by fixing unimportant factors. Unimportant factors are the ones that, if fixed to any value, would not lead to a significant reduction of the output uncertainty. This is often a preliminary step before the calibration of model inputs using some available information (real output observations, constraints, etc.);
- The variance cutting setting, which can be a part of a risk assessment study. Its aim is to reduce the output uncertainty from its initial value to a lower pre-established threshold value;

• The factors mapping setting, which aims at identifying the important inputs in a specific domain of the output values, for example which combination of factors produce output values above or below a given threshold.

In a deterministic framework, the model is analyzed at specific values for inputs, and the space of uncertain inputs may be explored in statistical approaches. In a probabilistic framework instead, the inputs are considered as random variables $\mathbf{X} = (X_1, \ldots, X_d) \in \mathbb{R}^d$. The random vector \mathbf{X} has a known joint distribution, which reflects the uncertainty of the inputs. The computer code (also called "model") is denoted $G(\cdot)$, and for a scalar output $Y \in \mathbb{R}$, the model formula writes

$$Y = G(\mathbf{X}) \ . \tag{1}$$

Model function G can represent a system of differential equations, a program code, or any other correspondence between \mathbf{X} and Y values that can be calculated for a finite period of time. Therefore, the model output Y is also a random variable whose distribution is unknown (increasing the knowledge of it is the goal of the uncertainty propagation process). SA statistical methods consist of techniques stemming from the design of experiments theory (in the case of a large number of inputs), the Monte Carlo techniques (to obtain precise sensitivity results) and modern statistical learning methods (for complex CPU time-consuming numerical models).

For example, to begin with the most basic (and essential) methods, simple graphical tools can be applied on an initial sample of inputs/output $(x_1^{(i)}, \ldots, x_d^{(i)}, y^{(i)})_{i=1..n}$ (sampling strategies are numerous and are described in many other chapters of this handbook). To begin with, simple scatterplots between each input variable and the model output can allow the detection of linear or non-linear input/output relation. Figure 1 gives an example of scatterplots on a simple function with three input variables and one output variable. As the two-dimensional scatterplots do not capture the possible interaction effects between the inputs, the cobweb plots [32] can be used. Also known as parallel coordinate plots, cobweb plots allow to visualize the simulations as a set of trajectories by joining the value (or the corresponding quantile) of each variables' combination of the simulation sample by a line (see Figure 2): Each vertical line represents one variable, the last vertical line representing the output variable. In Figure 2, the simulations leading to the smallest values of the model output have been highlighted in red. This allows to immediately understand that these simulations correspond to combinations of small and large values of the first and second inputs respectively.

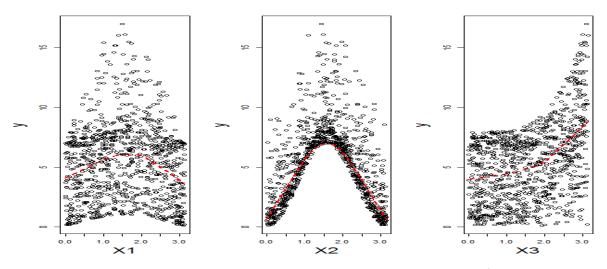


Fig. 1. Scatterplots of 200 simulations on a numerical model with three inputs (in abscissa of each plot) and one output (in ordinate). Dotted curves are local-polynomial based smoothers.

Moreover, from the same sample $\left(x_1^{(i)}, \ldots, x_d^{(i)}, y^{(i)}\right)_{i=1..n}$, quantitative global sensitivity measures can be easily estimated, as the linear (or Pearson) correlation coefficient and the rank (or Spearman) correlation coefficient [56]. It is also possible to fit a linear model explaining the behaviour of Y given the values of **X**, provided that the sample size n is sufficiently large (at least n > d). The main indices are then the Standard Regression Coefficients $\mathrm{SRC}_j = \beta_j \sqrt{\frac{\mathrm{Var}(x_j)}{\mathrm{Var}(Y)}}$, where β_j is the linear regression coefficient associated to X_j . SRC_j^2 represents a share of variance if the linearity

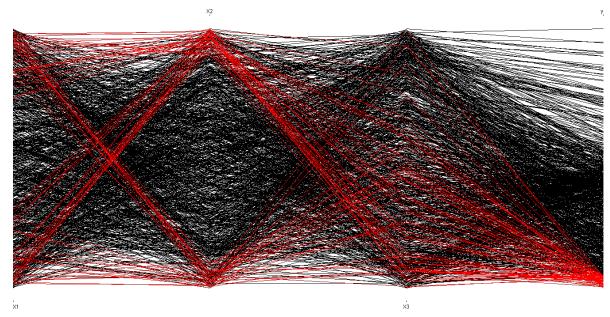


Fig. 2. Cobweb plot of 200 simulations of a numerical model with three inputs (first three columns) and one output (last column).

hypothesis is confirmed. Amongst many simple sensitivity indices, all these indices are included in the so-called sampling-based sensitivity analysis methods (see the description of the content of this chapter in the next section and Helton et al [23]).

Methods contained in the chapter

Three families of methods are described in this chapter, based on which is the objective of the analysis:

 First, screening techniques aim to a qualitative ranking of input factors at minimal cost in the number of model evaluations. Paper 2 (see <u>Variational Methods</u>, written by Maëlle Nodet and Arthur Vidard) introduces the local SA based on variational methods, while Paper 3 (see <u>Design of Experiments for Screening</u>, written by Sue Lewis and David Woods) makes an extensive review on design of experiments techniques, including some screening designs and numerical exploration designs specifically developed in the context of computer experiments;

- 2. Second, sampling based methods are described. Paper 4 (see <u>Weights and Importance in Composite Indicators: Mind the Gap</u>, written by William Becker, Paolo Paruolo, Michaela Saisana and Andrea Saltelli) shows how, from an initial sample of input and output values, quantitative sensitivity indices can be obtained by various methods (correlation, multiple linear regression, non-parametric regression) and applied in analyzing composite indicators. In Paper 5 (see <u>Variance-based Sensitivity Analysis: Theory and Estimation Algorithms</u>, written by Clémentine Prieur and Stefano Tarantola), the definitions of the variance-based importance measures (the so-called Sobol' indices) and the algorithms to calculate them will be detailed. In Paper 6 (see <u>Derivative based Global Sensitivity Measures</u>, written by Sergeï Kucherenko and Bertrand Iooss), the global SA based on derivatives sample (the DGSM indices) are explained, while in Paper 7 (see <u>Moment Independence Importance Measures and a Common Rationale</u>, written by Emanuele Borgonovo and Bertrand Iooss), the moment independent and reliability importance measures are described.
- 3. Third, in depth exploration of model behavior with respect to inputs variation can be carried out. Paper 8 (see <u>Metamodel-based Sensitivity Analysis: Polynomial Chaos and Gaussian Process</u>, written by Loïc Le Gratiet, Stefano Marelli and Bruno Sudret) includes recent advances made in the modeling of computer experiments. A metamodel is used as a surrogate model of the computer model with any SA techniques, when the computer model is too CPU time-consuming to allow a sufficient number of model calculations. Special attention is paid to two of the most popular metamodels: the polynomial chaos expansion and the Gaussian process model, for which the Sobol' indices can be efficiently obtained. Finally, Paper 9 (see <u>Sensitivity Analysis of Spatial and/or Temporal Phenomena</u>, written by Amandine Marrel, Nathalie Saint Geours and Matthias De Lozzo) extends the SA tools in the context of temporal and/or spatial phenomena.

All the SA techniques explained in this chapter present a trade-off between the number of model computations required and the assumed model complexity. Figure 3 proposes a coarse classification of the global method families described before. This figure shows how to place a method depending on its required number of computations and its underlying hypothesis on the complexity of the G model. For example "Non monotonic" means that the method can be applied to non monotonic models, as of course to monotonic and linear ones. A distinction is made between screening techniques (identification of non-influential variables among a large number of variables) and more precise variance-based quantitative methods. As most of the methods have a dimension-dependent cost in terms of required model evaluations, another distinction is made with the few methods whose costs are dimension-independent.

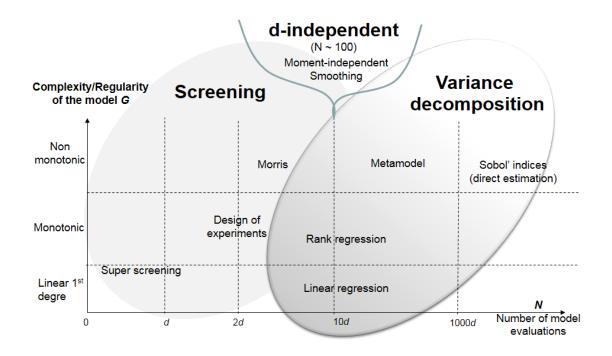


Fig. 3. Coarse classification of main global SA methods in terms of required number of model evaluations and model complexity.

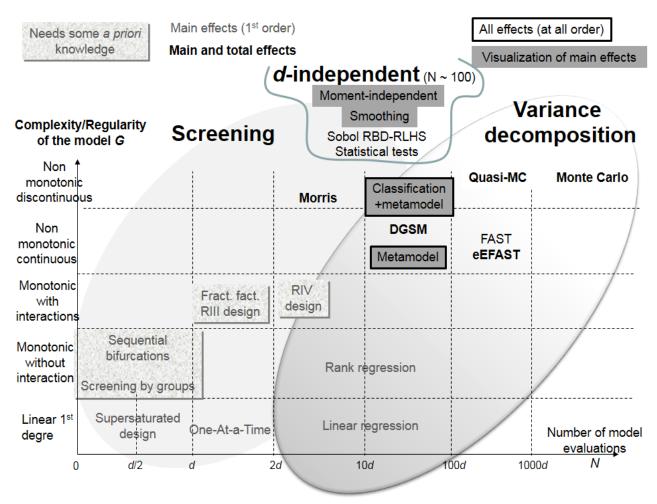


Fig. 4. Classification of global SA methods in terms of required number of model evaluations and model complexity.

With the same axes than the previous figure, Figure 4 proposes a more accurate classification of the classic global SA methods described in the present SA chapter. Note that this classification is not exhaustive and does not take full account of ongoing attempts to improve the existing methods. Overall, this classification tool has several levels of reading:

• positioning methods based on their cost in terms of the number of model calls. These methods linearly depend on the dimension (number of inputs) for most of the methods, except for the moment-independent measures (estimated with given-data approaches), smoothing methods, Sobol-RBD (Random Balance Design), Sobol-RLHS (Replicated Latin Hypercube Sampling) and statistical tests;

- positioning methods based on assumptions about model complexity and regularity;
- distinguishing the type of information provided by each method;
- identifying methods which require some prior knowledge about the model behavior.

Each of these techniques corresponds to different categories of problems met in practice. One should use the simplest method that is adapted to the study's objectives, the number of numerical model evaluations that can be performed, and the prior knowledge on the model's regularity. Each sensitivity analysis should include a validation step, which helps to understand if another method should be applied, if the number if model evaluations should be increased, and so on. Based on the characteristics of the different methods, some authors [9], [26] have proposed decision trees to help the practitioner to choose the most appropriate method for their problem and model.

Finally, the different papers of this chapter use the same mathematical notations, which are summarized below:

$G(\cdot)$	Numerical model
N, n	Sample sizes
d	Dimension of an input vector
p	Dimension of an output vector
$\mathbf{x} = (x_1, \dots, x_d)$	Deterministic input vector
$\mathbf{X} = (X_1, \dots, X_d)$) Random input vector
x_j, X_j	Deterministic and random input variable
\mathbf{x}^{T}	Transpose of \mathbf{x}
$\mathbf{x}_{\sim i} = \mathbf{x}_{-i} = \mathbf{x}_{\overline{i}}$	$= (X_1, \ldots, X_{i-1}, X_{i+1}, \ldots, X_d)$
$\mathbf{x}^{(i)}$	Sample vector of \mathbf{x}
$y, Y \in \mathbb{R}$	Deterministic and random output variable when $p = 1$
$y^{(i)}$	Sample of y
$f_X(\cdot)$	Density function of a real random variable X
$f_X(\cdot)$ $F_X(\cdot)$	Density function of a real random variable X Distribution function of X
$F_X(\cdot)$	Distribution function of X
$F_X(\cdot)$ μ_X, σ_X	Distribution function of X Mean and standard deviation of X
$F_X(\cdot)$ μ_X, σ_X $x_\alpha = q_\alpha(X)$	Distribution function of X Mean and standard deviation of X α -quantile of X
$F_X(\cdot)$ μ_X, σ_X $x_\alpha = q_\alpha(X)$ $V = \operatorname{Var}(Y)$	Distribution function of X Mean and standard deviation of X α -quantile of X Total variance of the model output Y

Specialized R software packages

From a practical point of view, the application of SA methods by researchers, engineers, end-users and students is conditioned by the availability of an easy-to-use software. Several software include some SA methods (see the Software chapter of the Springer Handbook) but only a few are specialized on the SA issues. In this section, the **sensitivity** package of the R environment is presented [49]. Its development has started since 2006 and the several contributions that this package has received have made it particularly complete. It includes most of the methods presented in the papers of this chapter.

The R software is a powerful tool for knowledge diffusion in the statistical community. The open source and availability have made R the software of choice for many statisticians in education and industry. The characteristics of R are the following:

- R is easy to run and install on main operating systems. It can be efficiently used with an old computer, with a single workstation and with one of the most recent supercomputer;
- R is a programming language which is interpreted and object-oriented, and which contains vector operators and matrix computation;
- The main drawbacks of R are its virtual memory limits and non-optimized computation times. To overcome these problems, compiled languages as fortran or C are much more efficient, and can be introduced as compiled codes inside R algorithms requiring huge computation time.
- R contains a lot of built-in statistical functions;
- R is extremely well-documented with built-in help system;
- R encourages the collaboration, the discussion forums and the creation of new packages by researchers and students. Thousands of packages are made available on the CRAN website (http://cran.r-project.org/).

All these benefits highlight the interest to develop specific softwares in R and many packages have been developed on SA. For instance, the FME package contains basic SA and local SA methods (see <u>Variational Methods</u>), while the **spartan** package contains basic methods for exploring stochastic numerical models. For global SA, The **sensitivity** package includes a collection of functions for factor screening, sensitivity indices estimation and reliability sensitivity analysis of model outputs. It implements:

- A few screening techniques as the sequential bifurcations and the Morris method (see <u>Design of Experiments for Screening</u>). Note that the R package planor allows to build fractional factorial design (see <u>Design of Experiments for Screening</u>);
- The main sampling-based procedures as linear regression coefficients, partial correlations, rank transformation (see <u>Weights and Importance in Composite</u> <u>Indicators: Mind the Gap</u>). Note that the ipcp function of the R package iplots provides an interactive cobweb graphical tool (see Figure 2), while the R package **CompModSA** implements various nonparametric regression procedures for SA (see Weights and Importance in Composite Indicators: Mind the Gap);
- The variance-based sensitivity indices (Sobol' indices), by various schemes of the so-called pick-freeze method, the Extended-FAST method and the replicated orthogonal array-based Latin hypercube sample (see <u>Variance-based Sensitivity</u> <u>Analysis: Theory and Estimation Algorithms</u>). The R package fast is fully devoted to the FAST method;
- The Poincaré constants for the Derivative-based Global Sensitivity Measures (DGSM) (see Derivative based Global Sensitivity Measures);
- The sensitivity indices based on Csiszar f-divergence and Hilbert-Schmidt Independence Criterion of Da Veiga [6] (see <u>Moment Independence Importance</u> Measures and a Common Rationale);
- The reliability sensitivity analysis by the Perturbation Law-based Indices (PLI) (see Moment Independence Importance Measures and a Common Rationale);
- The estimation of the Sobol' indices with a Gaussian process metamodel (see Metamodel-based Sensitivity Analysis: Polynomial Chaos and Gaussian

<u>Process</u>) with a Gaussian process metamodel coming from the R package DiceKriging. Note that the R package tgp performs the same job using treed Gaussian process and that the R package GPC allows to estimate the Sobol' indices by building a polynomial chaos metamodel (see <u>Metamodel-based Sensitivity</u> Analysis: Polynomial Chaos and Gaussian Process);

- Sobol' indices for multidimensional outputs: Aggregated Sobol' indices and functional (1D) Sobol' indices (see <u>Sensitivity Analysis of Spatial and/or Temporal Phenomena</u>). Note that the R package multisensi is fully devoted to this subject while the R package safi implements new SA methods of models with functional inputs;
- The Distributed Evaluation of Local Sensitivity Analysis (DELSA) described in Rakovec et al [50].

The sensitivity package has been designed to work either models written in R than external models such as heavy computational codes. This is achieved with the input argument model present in all functions of this package. The argument model is expected to be either a function or a predictor (i.e. an object with a predict function such as Im). The model is invoked once for the whole design of experiment. The argument model can be left to NULL. This is referred to as the decoupled approach and used with external computational codes that rarely run on the statistician's computer. Examples of use of all the sensitivity functions can be found using the R built-in help system.

As a global and generic platform allowing to include all the methods of these different R packages, the mtk package [70] has recently been proposed. It is an object-oriented framework which aims at dealing with external simulation platforms and managing all the different tasks of uncertainty and sensitivity analyses. Finally, the ATmet

and pse packages interface several sensitivity package functions for, respectively, metrology applications and parameter space exploration.

Sensitivity auditing

It may happen that a sensitivity analysis of a model-based study is meant to underpin an inference, and to certify its robustness, in a context where the inference feeds into a policy or decision making process. In these cases the framing of the analysis itself, its institutional context, and the motivations of its author may become a matter of great importance, and a pure SA - with its emphasis on parametric uncertainty - may be seen as insufficient. The emphasis on the framing may derive inter-alia from the relevance of the policy study to different constituencies that are characterized by different norms and values, and hence by a different story about 'what the problem is' and foremost about 'who is telling the story'. Most often the framing includes more or less implicit assumptions, which could be political (e.g. which group needs to be protected) all the way to technical (e.g. which variable can be treated as a constant).

These concerns about how the story is told and who tells it are all the more urgent in a climate as today's where science's own quality assurance criteria are under scrutiny due to a systemic crisis in reproducibility [25] and the explosion of the blogosphere invites more open debates on the scientific basis of policy decisions [42]. The Economist, a weekly magazine, has entered the fray by commenting on the poor state of current scientific practices and devoting its cover to "How Science goes wrong". It adds that, '*The false trails laid down by shoddy research are an unforgivable barrier to understanding*' (The Economist [66], p. 11). Among the possible causes of such a predicament is a process of hybridization [33] of fact and values, and of public and private institutions and actors. Thus the classical division of roles among science, providing tested fact, and policy, providing legitimized norms, becomes arduous to maintain. As an additional difficulty, according to Grundmann [19], 'One might suspect that the more knowledge is produced in hybrid arrangements, the more the protagonists will insist on the integrity, even veracity of their findings'.

In order to take these concerns into due consideration the instruments of SA have been extended to provide an assessment of the entire knowledge and model generating process. This approach has been called sensitivity auditing. It takes inspiration from NUSAP, a method used to qualify the worth of quantitative information with the generation of 'Pedigrees' of numbers [17], [69]. Likewise, sensitivity auditing has been developed to provide pedigrees of models and model-based inferences [61], [53], [54].

Sensitivity auditing has been especially designed for an adversarial context, where not only the nature of the evidence, but also the degree of certainty and uncertainty associated to the evidence, will be the subject of partian interests. Sensitivity auditing is structured along a set of seven rules/imperatives:

- 1. Check against the rhetorical use of mathematical modeling. Question addressed: is the model being used to elucidate or to obfuscate?;
- Adopt an 'assumption hunting' attitude. Question addressed: what was 'assumed out'? What are the tacit, pre-analytical, possibly normative assumptions underlying the analysis?;
- 3. Detect Garbage In Garbage Out (GIGO). Issue addressed: artificial deflation of uncertainty operated in order to achieve a desired inference at a desired level of confidence. It also works on the reverse practice, the artificial inflation of uncertainties, e.g. to deter regulation;
- 4. Find sensitive assumptions before they find you. Issue addressed: anticipate criticism by doing careful homework via sensitivity and uncertainty analyses before publishing results.

- 5. Aim for transparency. Issue addressed: stakeholders should be able to make sense of, and possibly replicate, the results of the analysis;
- 6. Do the right sums, which is more important than 'Do the sums right'. Issue addressed: is the viewpoint of a relevant stakeholder being neglected? Who decided that there was a problem and what the problem was?
- 7. Focus the analysis on the key question answered by the model, exploring the entire space of the assumptions holistically. Issue addressed: don't perform perfunctory analyses that just 'scratch the surface' of the system's potential uncertainties.

The first rule looks at the instrumental use of mathematical modeling to advance one's agenda. This use is called rhetorical, or strategic, like the use of Latin by the elites and clergy before the Reformation. At times the use of models is driven a simple pursuit of profit; according to Stiglitz [64], this was the case for the modelers 'pricing' the derivatives at the root of the sub-prime mortgages crisis:

[...] Part of the agenda of computer models was to maximize the fraction of, say, a lousy sub-prime mortgage that could get an AAA rating, then an AA rating, and so forth, [...] This was called rating at the margin, and the solution was still more complexity, p. 161.

At times this use of models can be called 'ritual', in the sense that it offers a false sense of reassurance. An example is Fisher [13] (quoting Szenberg [65]): Kenneth Arrow, one of the most notable Nobel Laureates in economics, has his own perspective on forecasting. During World War II, he served as a weather officer in the U.S. Army Air Corps and worked with a team charged with the particularly difficult task of producing month-ahead weather forecasts. As Arrow and his team reviewed these predictions, they confirmed statistically what you and I might just as easily have guessed: The Corps' weather forecasts were no more useful than random rolls of a die. Understandably, the forecasters asked to be relieved of this seemingly futile duty. Arrow's recollection of his superiors' response was priceless: "The commanding general is well aware that the forecasts are no good. However, he needs them for planning purposes" Szenberg [65].

The second rule about 'assumption hunting' is a reminder to look for what was assumed when the model was originally framed. Modes are full of *caeteris paribus* assumptions, meaning that e.g. in economics the model can predict the result of a shock to a given set of equations assuming that all the rest - all other input variables and inputs - remains equal, but in real life *caeteris* are never *paribus*, meaning by this that variables tend to be linked with one another, so that they can hardly change in isolation.

Furthermore at times the assumption made by modelers do not to withstand scrutiny. A good example of assumption hunting is from John Kay [28], where the author takes issue with modeling used in transport policy. This author discovered that among the input values assumed (and hence fixed) in the model was 'average car occupancy rates, differentiated by time of day, in 2035'. The point is that such assumptions are very difficult to justify. This comment was published in the Financial Times (where John Kay is a columnist) showing that at present times controversies that could be called epistemological evade the confines of academia and populate the media. Rule three is about artificially exaggerating or playing down uncertainties wherever convenient. The tobacco lobbies exaggerated the uncertainties about the health effects of smoking according to Oreskes and Conway [44], while advocates of the death penalty played down the uncertainties in the negative relations between capital punishment and crime rate [34]. Clearly the latter wanted the policy, in this case the death penalty, and were interested in showing that the supporting evidence was robust. In the former case the lobbies did not want regulation (e.g. bans on tobacco smoking in public places) and were hence interested in amplifying the uncertainty in the smoking-health effect causality relationship.

Rule four is about 'confessing' uncertainties before going public with the analysis. This rule is also one of the commandments of applied econometrics according to Kennedy [29]: 'Thou shall confess in the presence of sensitivity. Corollary: Thou shall anticipate criticism'. According to this rule a sensitivity analysis should be performed before the result of a modeling study are published. There are many good reasons for doing this, one being that a carefully performed sensitivity analysis often uncovers plain coding mistakes or model inadequacies. The other is that most often than not the analysis reveal uncertainties that are larger than those anticipated by the model developers. Econometrician Edward Learner in discussing this [34] argues that 'One reason these methods are rarely used is their honesty seems destructive'. In Saltelli and d'Hombres [52], the negative consequences of doing a sensitivity analysis *a-posteriori* are discussed. The case is the first review of the cost of offsetting climate change done by Nicholas Stern of the London School of Economics (the so called Stern Review) which was criticized by William Nordhaus, of the University of Yale, on the basis of large sensitivity of the estimates upon the discount factors employed by Stern. Sterns own sensitivity analysis, published as an annex to the review, revealed according to the authors in Saltelli and d'Hombres [52] that while the discount factors were not

the only important factors determining the cost estimate, the estimates were indeed very uncertain. On the large uncertainties of integrated assessment models of climate's impact, see also Saltelli et al [62].

Rule five is about presenting the results of the modeling study in a transparent fashion. Both rules originate from the practice of impact assessment, where a modeling study presented without a proper SA, or as originating from a model which is in fact a black box, may end up being rejected by stakeholders [52]. Both rules four and five suggest that reproducibility may be a condition for transparency and that this latter may be a condition for legitimacy. This debate on Sciences transparency is very much alive in the US, in the dialectic relationship between the US Environmental Protection Agency (EPA) and the US Congress (especially the Republican Party) which objects to EPAs regulations on the basis that these are based on 'secret science'.

Rule six, about doing the right sum, is not far from the 'assumptions hunting' rule; it is just more general. It deals with the fact that often an analyst is set to work on an analysis arbitrarily framed to the advantage of a party. Sometime this comes via the choice of the discipline selected to do the analysis. Thus an environmental impact problem may be framed through the lenses of economics, and presented as a cost benefit or risk analysis, while the issue has little to do with costs or benefits or risks and a lot to do with profits, controls, and norms. An example is in Marris et al [41] on the issue of GMOs, mostly presented in the public discourse as a food safety issue while the spectrum of concerns of GMO opponents - including lay citizens - appears broader. According to Winner [71] (p. 138-163), ecologists should not be led into the trap of arguing about the 'safety' of a technology after the technology has been introduced. They should instead question the broader power, policy and profit implications of that introduction and its desirability. Rule seven is about avoiding perfunctory sensitivity analyses. As discussed in Saltelli et al [60], an SA where each uncertain input is moved at a time while leaving all other inputs fixed is perfunctory. A true SA should make an honest effort at activating all uncertainties simultaneously, leaving the model free to display its full nonlinear and possibly non-additive behaviour. A similar point is made in Sam L. Savage's book 'The flaw of averages' [63].

In conclusion, these rules are meant to help an analyst to anticipate criticism. In drafting these rules the authors in [61], [53], [54] have tried to put themselves in the shoes of a modeler also based on their own experience and tried to imagine a modelbased inference feeding into an impact assessment. What questions and objections may be received by the modeler? Here is a possible list:

- "You treated X as a constant when we know it is uncertain by at least 30%"
- "It would be sufficient for a 5% error in X to make your statement about Z fragile"
- "Your model is but one of the plausible models you neglected model uncertainty"
- "You have instrumentally maximized your level of confidence in the results"
- "Your model is a black box why should I trust your results?"
- "You have artificially inflated the uncertainty"
- "Your framing is not socially robust"
- "You are answering the wrong question"

The reader may easily check that a careful go through the sensitivity auditing checklist should provide ammunition to anticipate objections of this nature.

Sensitivity auditing can then be seen as a user guide to criticize the model-based studies and SA is a part of this guide. In the following section, we go back to sensitivity analysis in order to conclude and give some perspectives.

Conclusion

This introductory paper has presented the SA aims and objectives, the SA and sensitivity auditing basic principles, the different methods explained in the chapter papers by positioning them in a classification grid, useful R software packages and the notations used in the chapter papers. The chapter's Editor, Bertrand Iooss, would like to sincerely thank all authors and co-authors of the "Sensitivity Analysis" chapter for their efforts and the quality of their contributions. Dr. Jean-Philippe Argaud (EDF R&D), Dr. Géraud Blatman (EDF R&D), Dr. Nicolas Bousquet (EDF R&D), Dr. Sébastien da Veiga (Safran), Dr. Hervé Monod (INRA), Dr. Matieyendou Lamboni (INRA) and Dr. Loïc Le Gratiet (EDF R&D) are also greatly thanked for their advices on the different chapter papers.

The papers in this chapter the most widely used methods for sensitivity analysis: the deterministic methods as the local sensitivity analysis, the experimental design strategies, the sampling-based and variance-based methods developed from the 1980s and the new importance measures and metamodel-based techniques established and studied since the 2000s. However, with such a rich subject, choices had to be made for the different chapter papers and some important omissions are present. For instance, the robust Bayesian analysis [1], [24] is not discussed, while it is nevertheless a great ingredient in the study of the sensitivity of Bayesian answers to assumptions and uncertain calculation inputs. Moreover, in the context of non-probabilistic representation of uncertainty (as in the interval analysis, the evidence theory and the possibility theory), a small amount of SA methods has been developed [22]. This subject is deferred to a future review work.

Due to their very new or incomplete nature, several other SA issues are not discussed in this chapter. For instance, estimating total Sobol' indices at low cost remains a problem of primary importance in many applications (see Saltelli et al [60] for a recent review on the subject). Second-order Sobol indices estimations have recently been considered by Fruth et al [16] (by way of total interaction indices) and Tissot and Prieur [67] (by way of replicated orthogonal-array LHS). The latter work offers a powerful estimation method because the number of model calls is independent of the number of inputs (as in the spirit of permutation-based technique [38], [37]). However, for high-dimensional models (several hundreds inputs), estimation biases and computational costs remain considerable; De Castro and Janon [8] have proposed to introduce modern statistical techniques based on variable selection in regression models. Owen [46] has introduced generalized Sobol' indices allowing to compare and search efficient estimators (as the new one found in Owen [45]).

Another mathematical difficulty is the consideration of the dependence between the inputs X_i (i = 1 ... d). Non-independence between inputs in SA has been discussed by many authors such as Saltelli and Tarantola [55], Jacques et al [27], Xu and Gertner [72], Da Veiga et al [7], Li et al [36], Kucherenko et al [31], Mara and Tarantola [39] and Chastaing et al [5]. Despite all these works, much confusion still exists in practical applications. In practice, it would be useful to be able to measure the influence of the potential dependence between some inputs on the output quantity of interest.

Note that most of the works in this chapter focus on SA relative to the overall variability of model output (second order statistics). In practice, one can be interested in other quantities of interest, such as the output entropy, the probability that the output exceeds a threshold and a quantile estimation [56], [15], [35]. This is an active area of research as shown in this chapter (see <u>Moment Independence Importance Measures and a Common Rationale</u>), but innovative and powerful ideas have recently been developed by Owen et al [47] and Geraci et al [18] using higher order statistics, Fort et al [14] using contrast functions and Da Veiga [6] using a kernel point of view.

Finally, in some situations, the computer code is not a deterministic simulator but a stochastic one. This means that two model calls with the same set of input variables lead to different output values. Typical stochastic computer codes are queuing models, agent-based models, and models involving partial differential equations applied to heterogeneous or Monte Carlo-based numerical models. For this type of code, Marrel et al [40] have proposed a first solution for dealing with Sobol' indices. Moutoussamy et al [43] have also tackled the issue in the context of the metamodel building of the output probability density function. Developing relevant SA methods in this context will certainly be subject of future works.

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