Computational methods for uncertainty quantification and sensitivity analysis of complex systems
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Computational methods for uncertainty quantification and sensitivity analysis of complex systems

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Labex MS2T - Université Technologique de Compiègne
Outline

1. Introduction to uncertainty quantification

2. Polynomial chaos expansions
   - Polynomial chaos basis
   - Computing the coefficients
   - Post-processing

3. Sensitivity analysis in multiscale problems
   - Single model / independent variables
   - Imbricated models: the issue of correlation
   - Distribution-based sensitivity indices
   - Covariance-based sensitivity indices

4. Application example
   - Mechanical example: composite beam
   - Tolerance analysis
Some common engineering structures

- Cattenom nuclear power plant (France)
- Cormet de Roselend dam (France)
- Military satellite
- Airbus A380
- Bladed disk
Computational models

- Modern engineering has to address problems of increasing complexity in various fields including infrastructures (civil engineering), energy (civil/mechanical engineering), aeronautics, defense, etc.

- Complex systems are designed using computational models that are based on:
  
  - a mathematical description of the physics (e.g. mechanics, acoustics, heat transfer, electromagnetism, etc.)

  - numerical algorithms that solve the resulting set of (e.g. partial differential) equations: finite element-, finite difference-, finite volume- methods, boundary element methods)
Computational models

Simulation models are calibrated and validated through comparison with lab experiments and in situ / full scale measurements. Once they are validated, these models may be run with different sets of input parameters in order to:

- **explore** the design space at low cost
- **optimize** the system w.r.t to cost criteria
- **assess** the robustness of the system w.r.t. uncertainties

**Sources of uncertainty**

- Differences between the **designed** and the **real** system in terms of material/physical properties and dimensions (tolerancing)
- Unforecast **exposures**: exceptional service loads, natural hazards (earthquakes, floods), climate loads (hurricanes, snow storms, etc.)
Global framework for managing uncertainties

**Step A**
Model(s) of the system
Assessment criteria

**Step B**
Quantification of sources of uncertainty

**Step C**
Uncertainty propagation

Random variables

Computational model

Moments
Probability of failure
Response PDF

**Step C’**
Sensitivity analysis

**Introduction**
Polynomial chaos expansions
Sensitivity analysis in multiscale problems
Application example
Global framework for managing uncertainties

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**Moments**
Probability of failure
Response PDF

**Step C’**
Sensitivity analysis
Step A: computational models

- Vector of input parameters: \( \mathbf{x} \in \mathbb{R}^M \)
- Computational model: \( \mathcal{M} \)
- Model response: \( y = \mathcal{M}(\mathbf{x}) \in \mathbb{R}^N \)

- geometry
- material properties
- loading
- analytical formula
- finite element model
- etc.
- displacements
- strains, stresses
- temperature, etc.
Step A: computational models

Vector of input parameters 
\[ \mathbf{x} \in \mathbb{R}^M \]

Computational model \( \mathcal{M} \)

Model response 
\[ \mathbf{y} = \mathcal{M}(\mathbf{x}) \in \mathbb{R}^N \]

- geometry
- material properties
- loading
- analytical formula
- finite element model
- etc.
- displacements
- strains, stresses
- temperature, etc.
Step B: probabilistic models of input parameters

No data exist

- expert judgment for selecting the input PDF’s of \( X \)
- literature, data bases (e.g. on material properties)
- maximum entropy principle

Input data exist

- classical statistical inference
- Bayesian statistics when data is scarce but there is some prior information

Data on output quantities

- inverse probabilistic methods and Bayesian updating techniques
Step B: probabilistic models of input parameters

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Step C: principles of uncertainty propagation

Input parameters: \( \mathbf{x} \in \mathbb{R}^M \)

Computational model: \( \mathcal{M} \)

Model response: \( y = \mathcal{M}(\mathbf{x}) \in \mathbb{R}^N \)

Random variables: \( \mathbf{X} \)

Random response: \( \mathbf{Y} = \mathcal{M}(\mathbf{X}) \)
Step C: principles of uncertainty propagation

Input parameters $\mathbf{x} \in \mathbb{R}^M$

Computational model $\mathcal{M}$

Model response $y = \mathcal{M}(\mathbf{x}) \in \mathbb{R}^N$

Random variables $\mathbf{X}$

Computational model $\mathcal{M}$

Random response $\mathbf{Y} = \mathcal{M}(\mathbf{X})$?
Step C: uncertainty propagation methods

Computational model

Probabilistic model

Mean/std. deviation

Rare event simulation

Response PDF

Probabilistic-computational model

Step A

Step B
Step C: uncertainty propagation methods

Computational model

Step A

Probabilistic model

Step B

Mean/std. deviation

µ

σ

Rare event simulation

P_f

Response PDF
Step C: uncertainty propagation methods

Computational model

Mean/std. deviation

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Response PDF

Step A

Step B
Step C: uncertainty propagation methods
Step C: uncertainty propagation methods

Computational model

Step A

Probabilistic model

Step B

Probabilistic-computational model

Mean/std. deviation

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Probabilistic model

Response PDF

B. Sudret (Chair of Risk & Safety)
Monte Carlo simulation

Paramètres d’entrée

$x_1, x_2, \ldots, x_M$

Modèle de simulation

Quantités d’intérêts

$y = M(x_1, x_2, \ldots, x_M)$
Monte Carlo simulation

Paramètres d’entrée

$x_1, x_2, \ldots, x_M$

Modèle de simulation

$y = \mathcal{M}(x_1, x_2, \ldots, x_M)$

Quantités d’intérêts
Monte Carlo simulation

Principles

- The input random variables are sampled according to their joint PDF $f_X(x)$.
- For each sample $x^{(i)}$, the response $M(x^{(i)})$ is computed (possibly time-consuming).
- The response sample set $M = \{M(x^{(1)}), \ldots, M(x^{(n)})\}^T$ is used to compute statistical moments, probabilities of failure or estimate the response distribution (histogram, kernel density estimation).
Monte Carlo simulation

Advantages

- This is a universal method, i.e. which does not depend on the type of model $M$.
- It is statistically well defined: convergence, confidence intervals, etc.
- It is non intrusive, i.e. it is based on repeated runs of the computational model as a black box.
- It is suited to distributed computing (clusters of PCs).

Drawbacks

- The “scattering” of $Y$ is investigated point-by-point: if two samples $x^{(i)}, x^{(j)}$ are almost equal, two independent runs of the model are carried out.
- The convergence rate is low ($\propto N^{-1/2}$).
Spectral approach

Principle

- The random response \( Y = \mathcal{M}(X) \) is considered as an element of a suitable vector space.
- A basis of this space is built up (with respect to the input joint PDF).
- The response random vector \( Y \) is completely determined by its coordinates in this basis:

\[
Y = \sum_{j=0}^{\infty} y_j \Psi_j(X)
\]

where:

- \( y_j \): coefficients to be computed (coordinates)
- \( \Psi_j(X) \): basis

Assumption: \( Y \) has a finite second moment, \( i.e. \):

\[
\mathbb{E} [Y^2] = \int Y^2(\omega) d\mathbb{P}(\omega) = \int \mathcal{M}^2(x) f_X(x) \, dx < \infty
\]
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4. Application example
Univariate orthogonal polynomials

Definition

- Suppose that the input random vector has independent components.

\[ f_X(x) = \prod_{i=1}^{M} f_{X_i}(x_i) \]

- For each marginal distribution, define:

\[ \langle \phi_1, \phi_2 \rangle_i = \int_{D_i} \phi_1(x) \phi_2(x) f_{X_i}(x_i) \, dx_i \]

- By classical algebra one can build a family of orthogonal polynomials \( \{ P_k^i, \ k \in \mathbb{N} \} \) as follows:

\[ \langle P_j^i, P_k^i \rangle = \int P_j^i(x) P_k^i(x) f_{X_i}(x) \, dx = a_j^i \delta_{jk} \]
Univariate orthogonal polynomials

Classical families

<table>
<thead>
<tr>
<th>Type of variable</th>
<th>Weight function</th>
<th>Orthogonal polynomials</th>
<th>Hilbertian basis $\psi_k(x)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>Uniform</td>
<td>$1] -1,1[ (x)/2$</td>
<td>Legendre $P_k(x)$</td>
<td>$P_k(x) / \sqrt{2k+1}$</td>
</tr>
<tr>
<td>Gaussian</td>
<td>$\frac{1}{\sqrt{2\pi}} e^{-x^2/2}$</td>
<td>Hermite $H^e_k(x)$</td>
<td>$H^e_k(x) / \sqrt{k!}$</td>
</tr>
<tr>
<td>Gamma</td>
<td>$x^a e^{-x} 1_{\mathbb{R}^+}(x)$</td>
<td>Laguerre $L^a_k(x)$</td>
<td>$L^a_k(x) / \sqrt{\Gamma(k+a+1)/k!}$</td>
</tr>
<tr>
<td>Beta</td>
<td>$1] -1,1[ (x) \frac{(1-x)^a(1+x)^b}{B(a)B(b)}$</td>
<td>Jacobi $J^{a,b}_k(x)$</td>
<td>$J^{a,b}_k(x) / \sqrt{\Gamma(k+a+1)}$</td>
</tr>
</tbody>
</table>

Normalization

- The classical orthogonal polynomials are not orthonormal: they may be normalized by dividing by $\sqrt{a_j^i}$:

$$\Psi_j^i = P_j^i / \sqrt{a_j^i} \quad i = 1, \ldots, M, \quad j \in \mathbb{N}$$
Multivariate polynomials

Tensor product of 1D polynomials

- One defines the multi-indices \( \alpha = \{\alpha_1, \ldots, \alpha_M\} \), of degree \( |\alpha| = \sum_{i=1}^{M} \alpha_i \)

- The associated multivariate polynomial reads:

\[
\Psi_\alpha(x) = \prod_{i=1}^{M} \Psi_{\alpha_i}(x_i)
\]

The set of multivariate polynomials \( \{\Psi_\alpha, \ \alpha \in \mathbb{N}^M\} \) forms a basis of the space:

\[
Y = \sum_{\alpha \in \mathbb{N}^M} y_\alpha \Psi_\alpha(X)
\]
Practical implementation

- The input random variables are first transformed into reduced variables (e.g. standard normal variables $\mathcal{N}(0, 1)$, uniform variables on $[-1,1]$, etc.):

$$X = T(\xi) \quad \text{dim} \xi = M \quad \text{(isoprobabilistic transform)}$$

- The model response is cast as a function of the reduced variables and expanded:

$$Y = M(X) = M \circ T(\xi) = \sum_{\alpha \in \mathbb{N}^M} y_\alpha \Psi_\alpha(\xi)$$

- A truncature scheme is selected and the associated finite set of multi-indices is generated, e.g.:

$$\mathcal{A} = \{\alpha \in \mathbb{N}^M : |\alpha| \leq p\} \quad \text{card } \mathcal{A} \equiv P = \binom{M + p}{p}$$
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Regression approach

Principle

The exact (infinite) series expansion is considered as the sum of a truncated series and a residual:

\[ Y = \mathcal{M}(X) = \sum_{j=0}^{P-1} y_j \Psi_j(X) + \varepsilon_P = Y^T \Psi(X) + \varepsilon_P \]

where: \[ Y = \{ y_0, \ldots, y_{P-1} \} \]
\[ \Psi(x) = \{ \Psi_0(x), \ldots, \Psi_{P-1}(x) \} \]

Mean-square minimization

The coefficients are gathered into a vector \( \hat{Y} \), and computed by minimizing the mean square error:

\[ \hat{Y} = \arg \min \mathbb{E} \left[ (Y^T \Psi(X) - \mathcal{M}(X))^2 \right] \]
Regression: discretized solution

The discretized mean-square minimization reads:

$$\hat{Y}^{reg} = \arg \min \frac{1}{n} \sum_{i=1}^{n} (Y^T \Psi(x^{(i)}) - M(x^{(i)}))^2$$

- Select an experimental design
  $$\mathcal{X} = \{x^{(1)}, \ldots, x^{(n)}\}^T$$
  that covers at best the domain of variation of the parameters

- Evaluate the model response for each sample (exactly as in Monte Carlo simulation)
  $$\mathcal{M} = \{M(x^{(1)}), \ldots, M(x^{(n)})\}^T$$

- Compute the experimental matrix
  $$A_{ij} = \Psi_j(x^{(i)}) \quad i = 1, \ldots, n; \quad j = 0, \ldots, P - 1$$

- Solve the least-square minimization problem
  $$\hat{Y} = (A^T A)^{-1} A^T \mathcal{M}$$
Validation of the surrogate model

- The truncated series expansions are convergent in the mean square sense. However one does not know in advance where to truncate (problem-dependent).

- Most people truncate the series according to the total maximal degree of the polynomials, say \( p=2,3,4, \text{ etc.} \). Several values of \( p \) are tested until some kind of convergence is “empirically” observed.

- The recent research deals with the development of error estimates:
  - adaptive integration in the projection approach
  - cross validation in the regression approach

- This has lead to the development of sparse polynomial chaos expansions.
Error estimators
Coefficient of determination

The regression technique is based on the minimization of the mean square error. The generalization error is defined as:

\[ E_{gen} = \mathbb{E} \left[ (\mathcal{M}(X) - \mathcal{M}^{PC}(X))^2 \right] \]

It may be estimated by the empirical error using the already computed response quantities:

\[ E_{emp} = \frac{1}{n} \sum_{i=1}^{n} \left( \mathcal{M}(x^{(i)}) - \mathcal{M}^{PC}(x^{(i)}) \right)^2 \]

The coefficient of determination \( R^2 \) is often used as an error estimator:

\[ R^2 = 1 - \frac{\hat{E}_{emp}}{\hat{\mathbb{V}}[\mathcal{Y}]} \quad \hat{\mathbb{V}}[\mathcal{Y}] = \frac{1}{n} (\mathcal{M}(x^{(i)}) - \bar{Y})^2 \]

This error estimator may lead to overfitting.
Error estimators
Leave-one-out cross validation

Principle

- In statistical learning theory, cross validation consists in splitting the experimental design $\mathcal{Y}$ in two parts, namely a training set (which is used to build the model) and a validation set.

- The leave-one-out technique consists in using each point of the experimental design as a single validation point for the meta-model built from the remaining $n - 1$ points.

- $n$ different meta-models are built and the error made on the remaining point is computed, then mean-square averaged.
For each \( x^{(i)} \), a polynomial chaos expansion is built using the following experimental design: \( \mathcal{X} \setminus x^{(i)} = \{ x^{(j)}, j = 1, \ldots, n, j \neq i \} \), denoted by \( \mathcal{M}^{PC \setminus i}(.) \).

The predicted residual is computed in point \( x^{(i)} \):

\[
\Delta_i = \mathcal{M}^{PC \setminus i}(x^{(i)}) - \mathcal{M}(x^{(i)})
\]

The PRESS coefficient (\textit{predicted residual sum of squares}) is evaluated:

\[
PRESS = \sum_{i=1}^{n} \Delta_i^2
\]

The leave-one-out error and related \( Q^2 \) error estimator are computed:

\[
E_{LOO} = \frac{1}{n} \sum_{i=1}^{n} \Delta_i^2 \quad Q^2 = 1 - \frac{E_{LOO}}{\hat{V}[Y]}
\]
Post-processing of polynomial chaos expansions

Reminder

Polynomial chaos

\[ Y = M(X) = M \circ T(\xi) \]

Truncated series

\[ Y_{PC} = \sum_{\alpha \in A} y_{\alpha} \Psi_{\alpha}(\xi) \]

- The computed coefficients ("coordinates" of the random variable in the Hilbertian basis) are not the quantities of interest.
- Depending on the situation, the PDF, the statistical moments or quantiles of \( Y \) are of interest (e.g. low quantiles in structural reliability analysis).

The PC expansion must be post-processed in order to get relevant information on the model response.
Statistical moments
Mean value and variance

From the orthonormality of the polynomial chaos basis one gets:

\[\mathbb{E}[\Psi_\alpha] = 0 \quad \mathbb{E}[\Psi_\alpha \Psi_\beta] = 0\]

Mean value

\[\hat{\mu}_Y = y_0\]

The estimated mean value is the first term of the series.

Variance

\[\hat{\sigma}_Y^2 = \sum_{\alpha \in \mathcal{A} \setminus 0} y_\alpha^2\]

The estimated variance is computed as the sum of the squares of the remaining coefficients.
Principle

- The polynomial series expansion may be considered as a stochastic response surface, i.e. an analytical function of the input variables $\xi$ (after some isoprobabilistic transform), that may be sampled easily using Monte Carlo simulation.

- A large sample set of reduced variables is drawn $\xi$, say of size $n_{sim} = 10^5 - 10^6$:

$$\mathcal{X}_{sim} = \{\xi^{(j)}, \ j = 1, \ldots, n_{sim}\}$$

- The truncated series is evaluated onto this sample:

$$\mathcal{Y}_{sim} = \left\{ \sum_{\alpha \in A} y_{\alpha} \Psi_{\alpha}(\xi^j), \ j = 1, \ldots, n_{sim} \right\}$$

- The obtained sample set is plotted as an histogram or by kernel density smoothing.
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Sensitivity analysis

Sobol’ decomposition

- Sensitivity analysis aims at quantifying what are the input parameters (or combinations thereof) that influence the most the response variability.
- Global sensitivity analysis relies on so-called variance decomposition techniques.

Consider a model $M : x \in [0, 1]^M \rightarrow M(x) \in \mathbb{R}$. The Sobol’ decomposition reads:

$$M(x) = M_0 + \sum_{i=1}^{M} M_i(x_i) + \sum_{1 \leq i < j \leq M} M_{ij}(x_i, x_j) + \cdots + M_{12...M}(x)$$

where:

- $M_0$ is the mean value of the function
- $M_i(x_i)$ are univariate functions
- $M_{ij}(x_i, x_j)$ are bivariate functions
- etc.
Sensitivity analysis

**Sobol’ decomposition**

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where:
- $\mathcal{M}_0$ is the mean value of the function
- $\mathcal{M}_i(x_i)$ are univariate functions
- $\mathcal{M}_{ij}(x_i, x_j)$ are bivariate functions
- etc.
Sobol’ indices
Independent input variables

Variance decomposition

- Assume $X_i \sim U(0, 1), i = 1, \ldots, M$ (possibly after some isoprobabilistic transform)

- Due to the orthogonality of the decomposition:

$$D \equiv \text{Var} \left[ \mathcal{M}(X) \right] = \mathbb{E} \left[ (\mathcal{M}(X) - \mathcal{M}_0)^2 \right]$$

$$= \mathbb{E} \left[ \left( \sum_{\{i_1, \ldots, i_s\} \subset \{1, \ldots, M\}} \mathcal{M}_{i_1 \ldots i_s}(X_{i_1}, \ldots, X_{i_s}) \right)^2 \right]$$

$$= \sum_{\{i_1, \ldots, i_s\} \subset \{1, \ldots, M\}} \mathbb{E} \left[ \mathcal{M}_{i_1 \ldots i_s}^2(X_{i_1}, \ldots, X_{i_s}) \right]$$
Sobol’ indices

Partial variance

- Consider:

\[ D_{i_1...i_s} = \int_{[0,1]^s} \mathcal{M}_{i_1...i_s}^2(x_{i_1}, \ldots, x_{i_s}) \, dx_{i_1} \ldots dx_{i_s} \]

- Then:

\[ \text{Var} [\mathcal{M}(X)] = \sum_{i=1}^{M} D_i + \sum_{1 \leq i < j \leq M} D_{ij} + \ldots + D_{12...M} \]

- The Sobol’ indices are obtained by normalization:

\[ S_{i_1...i_s} = \frac{D_{i_1...i_s}}{D} \]

They represent the fraction of the total variance \( \text{Var} [Y] \) that can be attributed to each input variable \( i \) \( (S_i) \) or combinations of variables \( \{i_1 \ldots i_s\} \).
Sobol’ decomposition vs. polynomial chaos expansions

The truncated polynomial series may be sorted so as to emphasize the uni-, bi-, etc.- variate functions of the Sobol’ decomposition, namely:

\[ M_0 = y_0 \]

\[ M_i(x_i) = \sum_{\alpha \in A_i} y_{\alpha} \Psi_{\alpha}(x) \quad A_i = \{ \alpha : \alpha_i > 0, \alpha_j \neq i = 0 \} \]

\[ \ldots \]

\[ M_{i_1 \ldots i_s}(x_{i_1}, \ldots, x_{i_s}) = \sum_{\alpha \in A_{i_1, \ldots, i_s}} y_{\alpha} \Psi_{\alpha}(x) \]

\[ A_{i_1 \ldots i_s} = \{ \alpha : \alpha_k > 0 \iff k \in (i_1, \ldots, i_s) \} \]

The PC expansion readily provides a functional decomposition
Link with PC expansions

Computation of the Sobol’ indices

- The partial variances $D_{i_1...i_s}$ are obtained by summing up the square of selected PC coefficients.

\[
D_i = \sum_{\alpha \in A_i} y_{\alpha}^2 \quad A_i = \{ \alpha : \alpha_i > 0, \alpha_j \neq i = 0 \}
\]

\[
D_{i_1...i_s} = \sum_{\alpha \in A_{i_1,...,i_s}} y_{\alpha}^2 \quad A_{i_1...i_s} = \{ \alpha : \alpha_k > 0 \Leftrightarrow k \in (i_1, \ldots, i_s) \}
\]

- The Sobol’ indices come after normalization:

\[
S_{i_1...i_s} = \frac{D_{i_1...i_s}}{D}
\]

Once the PC expansion is available, the full set of Sobol’ indices are obtained for free!
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The computational workflow consists in an imbrication of models (multiscale/multiphysics analysis).

Models of different levels have common parameters (e.g. geometrical dimensions for thermal, mechanical, acoustics problems).

The output of micro-scale models is the input of the macro-scale models.
Impact on sensitivity analysis

Specificity of multi-level imbricated models

● The PDF of the parameters of the intermediate levels (margins + dependence structure (correlations)) is implicitly defined by uncertainty propagation.

● The appropriate statistical tools are non-parametric representation of the marginals and copula theory.

Wand & Jones (1995); Nelsen (1999)

● Sensitivity indices have to be defined in the case of dependent input variables.

● Two approaches have emerged recently:
  ● Distribution-based sensitivity indices
  ● Covariance decomposition

Caniou (2012)
Borgonovo (2007, 2011)
Li & Rabitz (2010)
**Distribution-based sensitivity indices**

**Heuristics**

If the PDF of the model response $Y = \mathcal{M}(X)$ is influenced by a given input parameter $X_i$ then the **conditional** distribution $f_{Y|X_i}(y)$ shall significantly differ from $f_Y(y)$.

Shift between $f_Y$ and $f_{Y|X_i}$:

$$s(X_i) = \int_{D_Y} |f_Y(y) - f_{Y|X_i}(y)| \, dy$$

Borgonovo’s $\delta$ importance measure: expected shift

$$\delta_i \equiv \frac{1}{2} \mathbb{E} [s(X_i)] = \frac{1}{2} \int_{D_{X_i}} \left[ \int_{D_Y} |f_Y(y) - f_{Y|X_i}(y)| \, dy \right] f_{X_i}(x_i) \, dx_i$$

Schemes are taken from Caniou (2012)
Estimating the shift from PDFs

(Conditional) distributions are estimated by kernel smoothing techniques based on massive Monte Carlo simulation:

\[
\hat{f}_Y(y) = \frac{1}{N_{MC}h} \sum_{l=1}^{N_{MC}} K\left(\frac{y - \mathcal{M}(y^{(l)})}{h}\right)
\]

e.g. \(K(t) = e^{-t^2/2}/\sqrt{2\pi}\)

\(h\) is the bandwidth parameter, e.g. \(h \propto N_{MC}^{-1/5}\).

Shift estimation (inner integral)

\[s(x_i) = \int_{D_Y} |f_Y(y) - f_{Y|x_i}(y)| \, dy\]

Expected shift (outer integral)

\[E[s(X_i)] = \int_{D_{X_i}} s(X_i) f_{X_i}(x_i) \, dx_i\]

Computed by Gaussian quadrature

\[\approx \sum_{q=1}^{N_Q} \omega_q s(x_i^{(q)})\]
(Conditional) distributions are estimated by kernel smoothing techniques based on massive Monte Carlo simulation:

\[ \hat{f}_Y(y) = \frac{1}{N_{MC}} \sum_{l=1}^{N_{MC}} K\left( \frac{y - M(y^{(l)})}{h} \right) \]

\[ e.g. \quad K(t) = e^{-t^2/2}/\sqrt{2\pi} \]

\( h \) is the bandwidth parameter, e.g. \( h \propto N_{MC}^{-1/5} \).

**Shift estimation (inner integral)**

\[ s(x_i) = \int_{D_Y} |f_Y(y) - f_Y|_{X_i}(y)| \, dy \]

**Expected shift (outer integral)**

\[ E[s(X_i)] = \int_{D_{X_i}} s(X_i) \, f_{X_i}(x_i) \, dx_i \]

Computed by Gaussian quadrature

\[ \approx \sum_{q=1}^{N_Q} \omega_q s(x_i^{(q)}) \]
Estimating the shift from CDFs

Remarks

- The two nested integrals of functions derived from kernel PDF estimates are extremely expensive to compute and prone to numerical noise.
- The inner integral may be computed analytically though.

Alternative scheme

\[ s(X_i) = 2 \mathbb{P} (f_Y(y) > f_{Y|X_i}(y)) - 2 \mathbb{P} (f_Y(y) < f_{Y|X_i}(y)) \]

\[ \delta_i = \mathbb{E}_{X_i} [F_Y(y_2) - F_Y(y_1) + F_{Y|X_i}(y_1) - F_{Y|X_i}(y_2)] \]

where \((y_1, y_2)\) are the intersection points of the PDF and the conditional PDF, i.e. are the roots of:

\[ f_Y(y) = f_{Y|X_i}(y) \]
Covariance-based sensitivity indices

Suppose a functional decomposition of the model exists, e.g. :

\[ M(x) = \sum_{u \subset \{1, \ldots, M\}} M_u(x_u) \]

NB: the summands \( M_u(x_u) \) may not be orthogonal with respect to the probability measure associated with \( f_X \).

Covariance decomposition

\[
\text{Var} [M(X)] = \text{Var} \left[ \sum_{u \subset \{1, \ldots, M\}} M_u(X_u) \right] \\
= \sum_{u \subset \{1, \ldots, M\}} \text{Var} [M_u(X_u)] + \sum_{u \subset \{1, \ldots, M\}} \sum_{v \subset u} \text{Cov} [M_u(X_u); M_v(X_v)]
\]

NB: in case of independence and Sobol’ decomposition, the second term is zero.
Total sensitivity index

\[ S^{(T)}_u = \text{Cov} \left[ M_u(X_u), M(X) \right] / \text{Var} \left[ Y \right] \]

Structural sensitivity index

\[ S^{(S)}_u = \text{Var} \left[ M_u(X_u) \right] / \text{Var} \left[ Y \right] \]

Correlative sensitivity index

\[ S^{(C)}_u = S^{(T)}_u - S^{(S)}_u \]

- \( S^{(C)}_u = 0 \) on the case of independent input parameters and classical Sobol’ functional decomposition.

- These indices are denoted respectively by \( S_{p_j} \), \( S_{p_j}^a \) and \( S_{p_j}^b \) in Li, Rabitz et al. (2010).
Covariance-based indices using PC expansions

Principle

Functional decomposition

A truncated polynomial expansion is computed assuming that the parameters are independent:

$$M(x) \approx \sum_{\alpha \in A} y_{\alpha} \Psi_{\alpha}(x)$$

• The terms are grouped according to their input parameters so as to build the functional decomposition:

$$M_u(x_u) = \sum_{\alpha \in A_u} y_{\alpha} \Psi_{\alpha}(x) \quad A_u = \{ \alpha : k \in u \Leftrightarrow \alpha_k > 0 \}$$
Covariance-based indices using PC expansions

Monte Carlo estimators

- A sample set \( \mathcal{X} = \{ x^{(i)} , i = 1, \ldots , N \} \) of size \( N \) is drawn according to the joint PDF \( f_x(x) \).

- Classical estimators are used:

\[
\hat{\text{Var}} [ Y ] = \frac{1}{N - 1} \sum_{i=1}^{N} \left( M(x^{(i)}) - \mu_M \right)^2 \\
\hat{S}_u^{(T)} = \frac{1}{N - 1} \sum_{i=1}^{N} \left( M_u(x^{(i)}) - \mu_{M_u} \right) \left( M(x^{(i)}) - \mu_M \right) / \hat{\text{Var}} [ Y ] \\
\hat{S}_u^{(S)} = \frac{1}{N - 1} \sum_{i=1}^{N} \left( M_u(x^{(i)}) - \mu_{M_u} \right)^2 / \hat{\text{Var}} [ Y ] \\
\hat{S}_u^{(C)} = \hat{S}_u^{(T)} - \hat{S}_u^{(S)}
\]

NB: all the functions are multivariate polynomials. \( N = 10^5 \) evaluations may be carried out in a matter of seconds.
Outline

1. Introduction to uncertainty quantification
2. Polynomial chaos expansions
3. Sensitivity analysis in multiscale problems
4. Application example
   - Mechanical example: composite beam
   - Tolerance analysis
A composite beam of length $L$ and section $b \times h$ is made of a fraction $f$ of carbon fibers $(E_f, \rho_f)$ and a fraction $(1 - f)$ of epoxy matrix $(E_f, \rho_f)$.

The beam is loaded by its dead weight $q = \rho_{\text{hom}}gbh$:

Of interest is the sensitivity of the maximum midspan deflection $v$ to its input parameters:

$$v = \frac{5}{384} \frac{qL^4}{E_{\text{hom}}I}$$
Nested model and input variables

Models

\[ E_{\text{hom}} = f E_f + (1 - f) E_m \]
\[ I = \frac{bh^3}{12} \]
\[ q = [f \rho_f + (1 - f) \rho_m] g h \]
\[ v = \frac{5}{384} \frac{qL^4}{E_{\text{hom}} I} \]

Basic random variables

<table>
<thead>
<tr>
<th>Parameter</th>
<th>Distribution</th>
<th>Mean</th>
<th>Coefficient of variation</th>
</tr>
</thead>
<tbody>
<tr>
<td>( L )</td>
<td>Lognormal</td>
<td>2 m</td>
<td>1%</td>
</tr>
<tr>
<td>( b )</td>
<td>Lognormal</td>
<td>10 cm</td>
<td>3%</td>
</tr>
<tr>
<td>( h )</td>
<td>Lognormal</td>
<td>1 cm</td>
<td>3%</td>
</tr>
<tr>
<td>( E_f )</td>
<td>Lognormal</td>
<td>300 GPa</td>
<td>15%</td>
</tr>
<tr>
<td>( E_m )</td>
<td>Lognormal</td>
<td>10 GPa</td>
<td>15%</td>
</tr>
<tr>
<td>( \rho_f )</td>
<td>Lognormal</td>
<td>1800 kg·m(^{-3})</td>
<td>3%</td>
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<tr>
<td>( \rho_m )</td>
<td>Lognormal</td>
<td>1200 kg·m(^{-3})</td>
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<tr>
<td>( f )</td>
<td>Lognormal</td>
<td>0.5</td>
<td>10%</td>
</tr>
</tbody>
</table>

B. Sudret (Chair of Risk & Safety)
Labex MS2T - Seminar
February 14th, 2013 48 / 58
Output of the intermediate variables

- PC expansions of the intermediate variables $E_{\text{hom}}, q, I$
  
  \[ E_{\text{hom}} = \sum_{\alpha \in A_E} e_{\alpha} \Psi_{\alpha}(X) \]
  
  \[ q = \sum_{\alpha \in A_q} q_{\alpha} \Psi_{\alpha}(X) \]
  
  \[ I = \sum_{\alpha \in A_I} i_{\alpha} \Psi_{\alpha}(X) \]

- Non parametric estimation of the margins (kernel smoothing with $10^4$ points):

\[ \hat{f}_I(i) \]
\[ \hat{f}_q(q) \]
\[ \hat{f}_{E_{\text{hom}}}(e) \]
Output of the intermediate variables
Correlation structure

- Estimation of the dependence structure: Gaussian copula based on the Spearman’s rank correlation coefficients.

\[ \rho_S(q, E) = 0.20 \]
\[ \rho_S(q, I) = 0.71 \]

Last level uncertainty propagation
- Isoprobabilistic transform of \( L, I, q, E_{\text{hom}} \) into standard normal variates \( U \)
- PC expansion of \( v \) and computation of distribution-based (resp. covariance-based) sensitivity indices
Results - Distribution-based indices
CDF and conditional CDF of the maximal deflection

\[ F_Y(y), F_{Y|X_1}(y) \]
\[ \delta_1 = 0.015 \]
\[ F_Y(y), F_{Y|X_2}(y) \]
\[ \delta_2 = 0.003 \]
\[ F_Y(y), F_{Y|X_3}(y) \]
\[ \delta_3 = 0.546 \]
\[ F_Y(y), F_{Y|X_4}(y) \]
\[ \delta_4 = 0.129 \]
## Results

<table>
<thead>
<tr>
<th>Parameter</th>
<th>$\delta^{CDF}$</th>
<th>$S$</th>
<th>$S^{U}$</th>
<th>$S^{C}$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$q$</td>
<td>0.015</td>
<td>-0.08</td>
<td>0.09</td>
<td>-0.17</td>
</tr>
<tr>
<td>$L$</td>
<td>0.003</td>
<td>0.01</td>
<td>0.01</td>
<td>0.00</td>
</tr>
<tr>
<td>$E_{\text{hom}}$</td>
<td>0.546</td>
<td>0.89</td>
<td>0.94</td>
<td>-0.05</td>
</tr>
<tr>
<td>$I$</td>
<td>0.129</td>
<td>0.18</td>
<td>0.30</td>
<td>-0.12</td>
</tr>
<tr>
<td>$\Sigma$</td>
<td>0.693</td>
<td>1.00</td>
<td>1.34</td>
<td>-0.34</td>
</tr>
</tbody>
</table>

$v = \frac{5}{384} \frac{qL^4}{E_{\text{hom}}I}$

- The homogenized Young’s modulus and the moment of inertia are the most important parameters.
- The distribution- and covariance-based indices give consistent results.
- The cov.-based index of $q$ is negative due to the strong correlation between $q$ and $I$ which have opposite influences in the maximal displacement.
Outline

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Tolerance analysis

- In industrial mass production, assembled products are made of individual parts that are prone to uncertainty in their dimensions due to the manufacturing processes.

- **Non-assembly problems** may occur when some parts’ real dimensions differ (even slightly) from their theoretical values.

- Predicting the probability of non-assembly **prior** to the mass production is of crucial importance. Probabilistic methods derived from structural reliability analysis have been recently proposed.
Connector contact clearance

The axial deviation of an electrical connector pin is considered.

The deviation is computed as a non linear function of 14 geometrical dimensions describing the geometry of the two parts and their respective position.

Each dimension is characterized by a Gaussian variable.

Since different surfaces are machined during the same operation, their dimensions are highly correlated, e.g. \( \rho_S = 0.8 \) between the three different diameters of the pin.

After Gayton et al., Méca. Indus. (2011)
## Results

<table>
<thead>
<tr>
<th>Parameter</th>
<th>$\delta^{CDF}$</th>
<th>$S$</th>
<th>$S^U$</th>
<th>$S^C$</th>
</tr>
</thead>
<tbody>
<tr>
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<td>0.00</td>
<td>0.00</td>
<td>0.00</td>
<td>0.00</td>
</tr>
<tr>
<td>$cm_3$</td>
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<td>0.00</td>
<td>0.00</td>
<td>0.00</td>
</tr>
<tr>
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<td>0.01</td>
</tr>
<tr>
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<td>0.15</td>
<td>0.15</td>
<td>0.00</td>
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<tr>
<td>$cm_6$</td>
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<td>0.02</td>
<td>0.00</td>
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<tr>
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<tr>
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<tr>
<td>$ima_2$</td>
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<td>0.00</td>
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<tr>
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<tr>
<td>$\Sigma$</td>
<td>0.68</td>
<td>1.00</td>
<td>0.81</td>
<td>0.19</td>
</tr>
</tbody>
</table>

Sorted by $S^U$.

A set of **5 dimensions** explain 94% of the variance of the pin deviation.

Three of them are correlated input variables for which the correlated contribution of the cov.-based indices is larger than the uncorrelated part.

The $\delta$-indices are consistent.

The results may be used for tolerance allocation, *i.e.* allocate smaller tolerance intervals to important dimensions.

---

*After Caniou (2012)*
Complex systems are nowadays modelled by computational workflows that involve various submodels / physics / scales of description.

The existing uncertainty propagation methods have to be adapted to the specific characteristics of these models:

- large CPU demand: need for surrogate models, e.g. polynomial chaos expansions
- statistical dependence between the meaningful intermediate parameters of the workflow

Sensitivity indices have been developed for the case of models with dependent input parameters and can be used in the context of complex workflows:

- distribution-based $\delta$-indices
- covariance-based indices
Conclusions

- Polynomial chaos expansions whose efficiency in moment-, reliability- and sensitivity analysis is well established show a new feature in this context, namely the \textit{functional decomposition} of the computational model.

- Further work is required to make these methods available to the industry in a convenient format.

\textbf{Thank you very much for your attention!}
Conclusions

- Polynomial chaos expansions whose efficiency in moment-, reliability- and sensitivity analysis is well established show a new feature in this context, namely the *functional decomposition* of the computational model.

- Further work is required to make these methods available to the industry in a convenient format.

Thank you very much for your attention!
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