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Assessing Fire Safety using Complex Numerical Models with a Bayesian Multi-fidelity Approach

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ABSTRACT

Nowadays, fire safety engineers are increasingly relying on sophisticated numerical simulators, typically based on Computational Fluid Dynamics (CFD) solvers, to conduct their analyses. However, the complexity of these numerical models often limits drastically the number of simulations that can be afforded, making traditional methods of safety analysis difficult or impossible to apply. This paper proposes a statistical method to evaluate a quantity of interest with an expensive simulator while saving computation time. The method is based on Bayesian statistics and multi-fidelity. We use Gaussian process regression to construct a Bayesian model of the complex simulator. This model is based on a multi-fidelity approach, which consists in simulating at different levels of accuracy, for instance by varying the spatial discretization in a CFD solver. We illustrate the method on an example of fire safety analysis, where the quantity of interest is the probability of exceeding a tenability threshold in a building on fire.

KEYWORDS: *risk assessment, statistics, numerical experiments, meta-model, multi-fidelity*

NOMENCLATURE LISTING

General notations

f^{env}	distribution of environmental inputs
$f_{x,t}^{\text{sim}}$	distribution of output at (x,t)
\mathcal{N}	normal distribution
t	fidelity input
t_{HF}	highest level of fidelity
x	$= (x_e, x_s)$, physical inputs
x_e	environmental inputs
x_s	scenario inputs
Z	output of the simulator

Bayesian modeling

k	covariance function
m	mean function
n	number of observations
(x_i)	n points of observations
(x_i, z_i)	n observations
(z_i)	n values of observations
n	computed with n observations
λ	variance of the stochastic effect
μ	value of a constant mean function
ξ	mean response of the simulator

Estimation of conformity

n_p	number of sample paths
n_x	number of values of environmental inputs
$P_f(x,t)$	probability of failure at the input (x,t)
$P_f(x_s)$	probability of failure at the scenario x_s
z^{crit}	critical safety threshold

Application

A_f	fire area (m^2)
H_f	smoke radiative heat flux ($\text{kW} \cdot \text{m}^{-2}$)
P_{atm}	atmospheric pressure (bar)
\dot{Q}''	maximal heat release rate per unit area ($\text{kW} \cdot \text{m}^{-2}$)
q_{fd}	total released energy per unit area ($\text{MJ} \cdot \text{m}^{-2}$)
T_{amb}	ambient temperature ($^{\circ}\text{C}$)
T_c	temperature at 1.8 m ($^{\circ}\text{C}$)
T_{ext}	external temperature ($^{\circ}\text{C}$)
t_{xyz}	mesh size (cm)
V	visibility (m)
Y_{soot}	soot yield ($\text{kg} \cdot \text{kg}^{-1}$)
α	fire growth acceleration ($\text{kW} \cdot \text{s}^{-2}$)

ACRONYMS

CFD	Computational Fluid Dynamics	GP	Gaussian Process
DOE	Design Of Experiments	LHS	Latin Hypercube Sampling
FDS	Fire Dynamics Simulator	OpenFOAM	Open Field Operation And Manipulation

INTRODUCTION

The use of sophisticated numerical simulators, often based on Computational Fluid Dynamics (CFD) solvers such as Fire Dynamics Simulator (FDS) or Open Field Operation And Manipulation (OpenFOAM), is becoming more and more common in fire safety analyses. These simulators can be used to test various fire scenarios, which would be too difficult or impossible to test using real experiments [1]. However, they are usually very time-consuming, which limits the number of numerical experiments that can be afforded. Consequently, traditional mathematical methods for fire safety analysis based on Monte-Carlo methods [2; 3] become impractical, and fire safety engineers have to assess the safety of a building from a small number of simulation results. Moreover, operator dependence may cause different conclusions on a same problem, and complicate the comparison of experiments.

This paper proposes a new approach to help carry out fire safety analyses using complex, expensive-to-run numerical simulators, with a limited budget of simulation time. This approach comes from the statistical literature on the design and analysis of computer experiments (see, e.g., [4] and references therein), and is based on two key ingredients: *Bayesian statistics* and *multi-fidelity*. Bayesian statistics, on the one hand, is a mathematical framework for data analysis, where results are expressed in term of probability statements [5]. Each unknown quantity of interest is equipped with a probability distribution, which describes current knowledge of the quantity, and is updated whenever additional data becomes available (see [6] for an overview). The Bayesian method presented in this text is based on Gaussian process regression, which is a well-established approach in the statistical literature for modeling complex computer experiments. Multi-fidelity, on the other hand, is a technique that can be used when fast approximations of the simulator are available [7–11]. Typically, this is the case when the accuracy of the simulation can be tuned by changing the mesh size of the spatial discretization and/or its time step. A simulation with a coarser mesh describes the same physical phenomenon with a lower fidelity, but runs faster than a high-fidelity simulation. The multi-fidelity approach combines low- and high-fidelity simulations to obtain information about the quantities of interest faster than with high-fidelity simulations only.

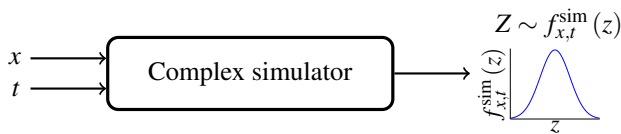


Fig. 1. Diagram of a numerical experiment. A simulator is described as a black box. The implementation is ignored, only the inputs and output are considered.

Let us consider the simulator as a *black box*, taking physical inputs x and a fidelity input t , and returning an output Z (see Figure 1). The fidelity input t controls level of the fidelity of simulations [9]. The physical input x represent physical quantities (temperatures, fire properties, boundary conditions. . .). In our approach, several forms of uncertainty affecting the output must be distinguished [12]. First, some of the physical inputs can be imperfectly known in the scenario under study. This lack of knowledge is modeled by a probability distribution, which conveys a first form of uncertainty, called *input uncertainty*. Second, the output Z (which may be the heat flux, the visibility. . .) is assumed stochastic. A typical example is the case of a deterministic CFD solver with random initial or boundary conditions, or a random forcing term. Thus, the output at a given input point (x, t) follows a certain distribution $f_{x,t}^{\text{sim}}(z)$. This constitutes a second form of uncertainty, called *output uncertainty*. Finally, we consider a third form of uncertainty, the *Bayesian modeling uncertainty*, which represents the lack of knowledge about the simulator itself, i.e., about the output distributions $f_{x,t}^{\text{sim}}(z)$. This form of uncertainty is progressively reduced when the number of simulations grows, and with it the available knowledge about the simulator. The method proposed in this paper will provide, along with estimations of the quantities of interest, a quantification of the uncertainty about these estimations resulting from these three sources of uncertainty.

The paper is organized as follows. The second section provides a short introduction to Gaussian process regression, and discusses the assumption of stochasticity about the output. The third section presents the main

contribution of this paper, which is a methodology based on Bayesian multi-fidelity modeling to estimate virtually any quantity of interest using a limited budget of simulation time, and quantify the associated uncertainty. The fourth section illustrates the methodology with an example of conformity assessment, where the quantity of interest is the probability of exceeding a tenability threshold in a given period of time after the start of a fire (this probability, also called probability of failure, can be for instance the probability that the mean heat flux exceeds $2.5 \text{ kW} \cdot \text{m}^{-2}$ or that the mean temperature at 1.8 m exceeds 60°C). The final section concludes the paper with a summary of our contributions and discussion of future work.

BAYESIAN APPROACH TO MODELING THE OUTPUT OF THE SIMULATOR

Gaussian process regression

The objective of this section is to provide a short introduction to Gaussian process regression. Here, for the sake of simplicity, the input t corresponding to the mesh size is not considered. There is no loss of generality in doing so at this point.

Let $(x_i, z_i)_{1 \leq i \leq n}$ be a data set of n observations of the simulator, where the x_i 's are inputs, and the z_i 's are the associated outputs. Each x_i is a vector whose components correspond to the physical inputs of the simulator (temperatures, fire area...). Each z_i corresponds to a scalar output of the simulation (e.g., the visibility) associated to x_i . To model the fact that the output of the simulator is stochastic, we assume that the observed response is a random variable Z with a Gaussian probability density function f_x^{sim} such that

$$f_x^{\text{sim}}(z) = \frac{1}{\sqrt{2\pi\lambda(x)}} e^{-\frac{1}{2} \frac{(z-\xi(x))^2}{\lambda(x)}}, \quad z \in \mathbb{R}, \quad (1)$$

where $\xi(x)$ and $\lambda(x)$ stand respectively for the mean and variance of the stochastic output of the simulator at x . In this section, we assume that λ is known, and ξ is unknown.

To deal with the fact that ξ is unknown, we assume that ξ is a Gaussian Process (GP), which will be denoted by

$$\xi \sim \mathcal{GP}(\mathbf{m}, \mathbf{k}), \quad (2)$$

where \mathbf{m} and \mathbf{k} are respectively the mean and covariance functions of the Gaussian process. The mean function is assumed to be constant and unknown. The covariance function is chosen by the user in such a way to describe how he/she expects two values $\xi(x)$ and $\xi(x')$ to be similar if x and x' are close to each other. In particular, the covariance function specifies the smoothness of the simulator. From a Bayesian point of view, the Gaussian process represents our prior knowledge about the unknown mean response of the simulator.

Once our Bayesian model is set up, the next step is to condition the model on the observations $(x_i, z_i)_{1 \leq i \leq n}$. Bayes' theorem yields the posterior distribution of ξ . Under the assumptions of a constant mean function, $\mathbf{m}(x) = \mu$, with a uniform prior on μ , and known covariance and output variance functions \mathbf{k} and λ , the posterior distribution of $\xi(x)$ is Gaussian, with mean

$$\mathbf{m}_n(x) = \widehat{\mu} + \mathbf{k}^T(x) [\mathbf{K} + \mathbf{\Lambda}]^{-1} (\mathbf{z} - \widehat{\mu} \mathbf{1}) \quad (3)$$

and variance

$$\sigma_n^2(x) = \mathbf{k}(x, x) - \mathbf{k}^T(x) [\mathbf{K} + \mathbf{\Lambda}]^{-1} \mathbf{k}(x) + \frac{\left(1 - \mathbf{1}^T [\mathbf{K} + \mathbf{\Lambda}]^{-1} \mathbf{k}(x)\right)^2}{\mathbf{1}^T [\mathbf{K} + \mathbf{\Lambda}]^{-1} \mathbf{1}}, \quad (4)$$

where $\mathbf{z} = (z_i)$ are the outputs, $\mathbf{K} = (\mathbf{k}(x_i, x_j))$ is the covariance matrix of the observations, $\mathbf{\Lambda} = \text{Diag}(\lambda(x_i))$ is the diagonal matrix of variances of stochastic effects at observation points, $\mathbf{k}(x) = (\mathbf{k}(x_i, x))$ is the covariance vector between observations and the point x , $\mathbf{1}$ a $n \times 1$ is vector of ones, and

$$\widehat{\mu} = \left(\mathbf{1}^T [\mathbf{K} + \mathbf{\Lambda}]^{-1} \mathbf{1}\right)^{-1} \mathbf{1}^T [\mathbf{K} + \mathbf{\Lambda}]^{-1} \mathbf{z}$$

is the generalized least squares estimate of μ . We refer the reader to Stein [13] or Forrester et al. [14] for more details about how these expressions can be established. Note that the function \mathbf{m}_n can be seen as an estimator

of ξ , and the function σ_n can be seen as a measure of our lack of knowledge about ξ . The function m_n is also called a meta-model, or an emulator, of the mean response of the simulator. The computation of the posterior distribution of ξ is usually carried out using dedicated softwares. The Small (Matlab/Octave) Toolbox for Kriging (STK) [15] has been used in this paper. Once the posterior distribution is obtained, it can be used to estimate quantities of interest, and in particular, a probability of exceeding a safety threshold.

Example

This section illustrates the approach on a simple example. We consider a numerical model with a scalar input x and a stochastic scalar output Z . The quantity of interest is the mean response of the numerical model.

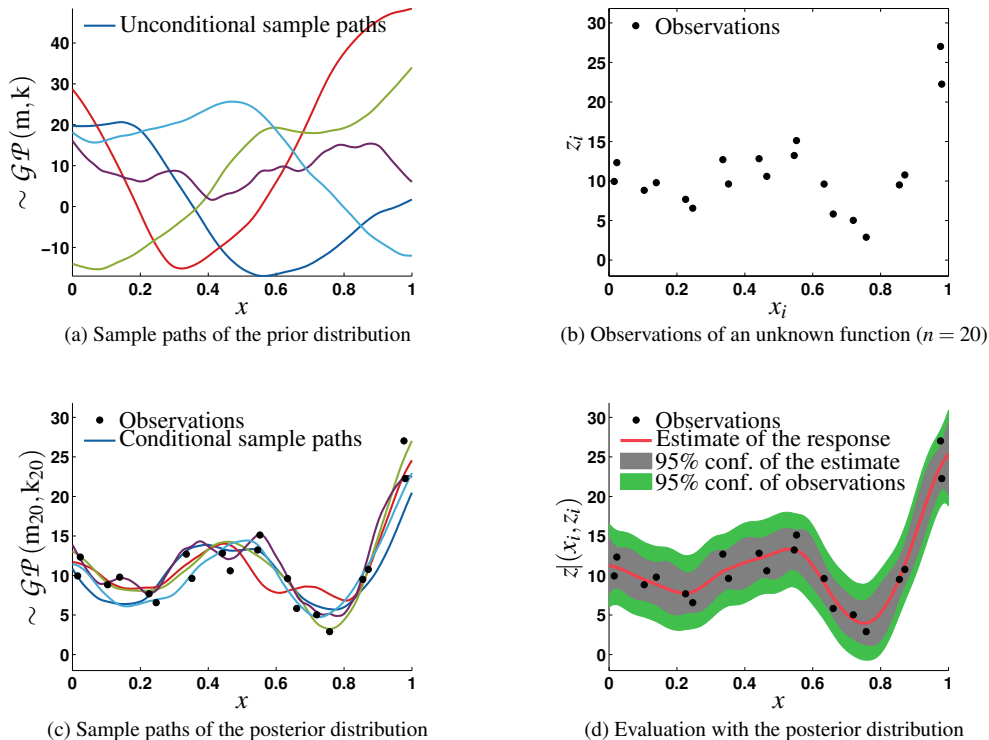


Fig. 2. Illustration of Gaussian process regression. (a) The unconditioned Gaussian process. (b) Observations of the numerical model. (c) The conditioned Gaussian process. (d) An estimate of the mean response.

The first step of the Bayesian approach is to define a prior distribution. Although the numerical model is seen as a black box, some assumptions can be made about it, for instance about its smoothness. These assumptions are translated into a Gaussian process prior about ξ . Figure 2a shows five sample paths of ξ , which represent what the mean response could be prior to any observation.

The second step is to observe the numerical model. Thus, n inputs x_1, \dots, x_n are selected. Then, for each input x_i , the numerical model is run, and returns a result z_i . Figure 2b shows $n = 20$ observations (x_i, z_i) corresponding to the stochastic outcome of the numerical model.

Conditioning the prior distribution on the observations produces the posterior distribution of the output of the numerical model. Figure 2c shows five sample paths of the posterior distribution. The sample paths of the posterior distribution have smoothness properties similar to those of the prior distribution, but they now “capture” the observations.

Finally, Figure 2d represents an estimate of the unknown mean response. The solid line is the function m_n (Eq. (3)), which is an estimator of the unknown mean response. The inner area around the line is a 95% confidence interval of the estimate, which is proportional to σ_n (Eq. (4)). The outer area is a 95% confidence

interval of future observations, taking into account the stochastic effect. In other words, a new observation of the unknown numerical model has a 95% chance to be in this area according to the posterior distribution. Confidence intervals are sometimes called credible intervals by Bayesian statisticians.

Remarks on the stochastic nature of the output

We have assumed until now that the output Z of the simulator at a given input point is stochastic. As explained in the introduction, a typical example is the case of a deterministic CFD solver with random initial or boundary conditions. Stochastic simulators have the property that running several simulations at the same point yields different output values. This calls for two remarks.

First, it is important to note that smooth deterministic simulators can be seen as a special case of the above model (when $\lambda(x)$ becomes equal to zero everywhere). The main difference, in this case, is that conditional sample paths (see Figure 2c) interpolate between the observed values. Thus, the general methodology proposed in this paper can actually be used with such simulators as well, with some minor modifications of the underlying statistical model.

Second, we emphasize that the same statistical model can also be used in the case of a deterministic simulator, when the output can be decomposed as the sum of the smooth term and a very irregular (visually, “noisy”) term. In such a situation, running two simulations with almost—but not strictly—identical inputs yields different outputs, exactly as in the case of a truly stochastic simulator. Using an additional Gaussian “noise term” to represent this local variability is a well-established practice in the statistical literature (see, e.g., Gramacy and Lee [16]). This term, sometimes called *nugget effect* in reference to the geostatistical literature, corresponds to the variance $\lambda(x)$ in our model. From a mathematical point of view, as long as no replicated simulations (i.e., simulation at the same input x) are performed, this is perfectly equivalent to the model described earlier in this section except for the prediction of a future simulation output at an already observed location.

PROPOSED METHODOLOGY

Multi-fidelity

Because of the complexity of CFD computer models, the number of observations that can be made is limited. In this section, we will detail how, by making specific assumptions about the Bayesian model presented above, we can derive a multi-fidelity model.

Multi-fidelity is a solution to reduce the uncertainty in the posterior distribution at a small budget of computation time. In a simulator using a CFD solver based on a field model, such as FDS or OpenFOAM, the accuracy of a simulation is controlled by the spatial discretization. A high-density mesh improves space accuracy (hence should also increase fidelity to physical reality) but also increases computation time. The multi-fidelity approach consists in simulating at different levels of fidelity, instead of only the highest one. Thus, the idea is to combine fast low-fidelity simulations (by changing the mesh size) with slow high-fidelity simulations to improve the quality of the posterior model.

In the literature, multi-fidelity often deals with only two levels of fidelity [7; 8]. In this article, we work with a continuum of levels of fidelity indexed by a positive scalar t , following Picheny and Ginsbourger [9] and Tuo et al. [10]. The level t can be seen as an input of the simulator, as the input vector x (see Figure 1). In this work, since the level of fidelity t is related to the mesh size, fidelity increases when t tends to zero, even if this level of spatial discretization is impossible to simulate. We denote by t_{HF} the highest available level of fidelity (lowest value of t , corresponding to the finest meshes). This ideal level of fidelity t_{HF} should be selected in order to adequately resolve the fluid flow and fire dynamics. In the case of a CFD code, where t corresponds to the size of mesh cells, Lin et al. [17] recommend to use $t_{HF} \leq D^*/10$, with D^* the characteristic fire diameter.

The first subsection proposes a prior Gaussian-process-based multi-fidelity model of the simulator to describe the relationship between the levels of spatial discretization. Then, the second subsection describes how the numerical experiments are selected. As the observations improve the knowledge about the simulator, the distribution modeling the simulator is updated by conditioning. This posterior distribution can be used to estimate any quantity of interest. The third subsection shows an example of such an estimation, with the

estimation of a probability of failure. Note that, even if we focus only on the highest level of fidelity t_{HF} , this posterior model could be used at any level of fidelity, in other words, for each spatial resolution.

Multi-fidelity prior model

The simulators that we aim to describe have a fidelity input t and a stochastic output Z . Following the model of stochastic simulator described Eq. (1), we suppose that the output Z at (x, t) follows a normal probability distribution, whose parameters depend on (x, t) :

$$f_{x,t}^{\text{sim}}(z) = \frac{1}{\sqrt{2\pi\lambda(x,t)}} \cdot e^{-\frac{1}{2} \cdot \frac{(z-\xi(x,t))^2}{\lambda(x,t)}}, \quad (5)$$

where $\xi(x, t)$ and $\lambda(x, t)$ are respectively the mean and the variance of the output at (x, t) , which are both assumed unknown now. The next paragraphs present briefly our prior models about ξ and λ .

The prior distribution that we propose for the mean response ξ was simultaneously introduced, with minor differences, by Picheny and Ginsbourger [9] and Tuo et al. [10] for deterministic simulators. To build this prior model, the mean response ξ is assumed to be the sum of an ideal response ξ_0 and a numerical error ε :

$$\xi(x, t) = \xi_0(x) + \varepsilon(x, t), \quad (6)$$

where the ideal response ξ_0 corresponds to the case where t tends to zero, whereas the numerical error ε stems from the numerical approximations of the equations of physics. Thus, it is assumed that $\varepsilon(x, t = 0) = 0$ for all x . Moreover, the two processes ξ_0 and ε are supposed independent, which means that knowing the physical phenomena represented by the simulator (ξ_0) does not help to know how the simulator approximate these phenomena (ε). Following the suggestion of Tuo et al. [10], we propose to use the following covariance for the process ε :

$$k_\varepsilon((x, t), (x', t')) = \min\{t, t'\}^L \tilde{k}_\varepsilon(x, x'), \quad (7)$$

where k_ε is the covariance function of the numerical error ε , and \tilde{k}_ε is a covariance function defined on the space of the physical inputs x . The term $\min\{t, t'\}^L$ is such that the numerical error decreases when t decreases, i.e., when fidelity increases. Notice that the numerical error and the stochastic effect are two distinct parts of the simulator. The simulator would still be stochastic, even if the numerical error was zero.

The prior distribution that we propose for the noise variance λ is described in detail by Stroh et al. [18]. In order to simplify the model, we suppose that the noise variance depends only on the levels of spatial discretization, so that λ depends only on the fidelity levels t . Then, the noise variance is assumed to be different at each level, but have the same order of magnitude. These assumptions are summed up in a log-normal prior distribution.

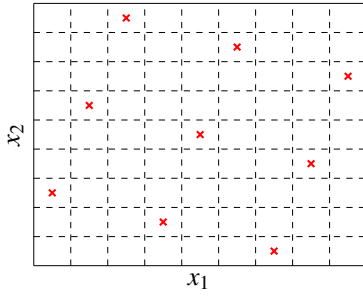
Design of Experiments

This section presents how numerical experiments are selected once the Bayesian model has been set.

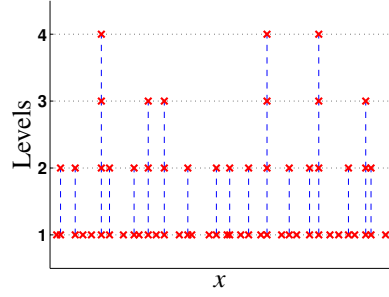
In our multi-fidelity framework, we choose the set of observations, also called Design Of Experiments (DOE), in such a way that the computation time remains low. It is well-known that to construct a good Bayesian model, the DOE should be space-filling (see Pronzato and Müller [19] for a review). One of the most popular method to obtain a space-filling DOE is to use maximin Latin Hypercube Sampling (LHS). A design is LHS, when the inputs are regularly spaced along all one-dimensional axis (see Figure 3a for an example of space-filling LHS). More details can be found in Santner et al. [4].

In our multi-fidelity approach, since low-fidelity is cheaper than high-fidelity, low-fidelity should be more often observed than high-fidelity. To this purpose, we use nested LHS. A design is said to be nested, when each point observed at an high-fidelity level is also observed at lower-fidelity levels (see Figure 3b).

In order to construct a nested LHS design, we use the algorithm developed by Qian [20], which produces nested LHS design, with an additional maximin optimization at each level, to obtain a good space-filling.



(a) A space-filling LHS, $n = 9$, $d = 2$



(b) A nested design, $n = 36 + 18 + 6 + 3$, $d = 1$, $S = 4$

Fig. 3. Two examples of design. (a) A Latin Hypercube Sampling. (b) A nested design (the lowest level of fidelity is below). n is the number of points, d the dimension, S the number of levels of fidelity.

This method imposes a constraint on the size of the design: the number of design points at a given level must be a multiple of the number of points at the next (finer) level.

Estimation of the conformity of a building using the multi-fidelity model

Given the results of the simulations, we can compute the posterior distribution which can be used in turn to estimate quantities of interest, and therefore assess the conformity of a building during a fire. In this paper, conformity is assessed by comparing the output Z to a critical threshold z^{crit} . More precisely, we estimate the probability that Z exceeds the threshold z^{crit} in various fire scenarios, in order to find which scenarios are the most dangerous.

The procedure to estimate the probability of failure at a given fire scenario is based on two steps. First, we consider a pointwise probability of failure defined as the probability that the output Z at (x, t) exceeds its safety threshold z^{crit} (for instance, the probability that the heat flux predicted by the simulator at a given fire scenario, weather conditions and spatial discretization exceeds $2.5 \text{ kW} \cdot \text{m}^{-2}$),

$$P_f(x, t) = \mathbb{P}(Z > z^{\text{crit}}) = \int_{z^{\text{crit}}}^{+\infty} f_{x,t}^{\text{sim}}(z) dz, \quad (8)$$

where $f_{x,t}^{\text{sim}}(z)$ denotes the output density at (x, t) (see Introduction). Recall from (5) that Z is assumed to follow a normal distribution. Given ξ and λ , the pointwise probability of failure can be computed as

$$P_f(x, t) = \Phi\left(\frac{\xi(x, t) - z^{\text{crit}}}{\sqrt{\lambda(x, t)}}\right), \quad (9)$$

where Φ is the normal cumulative distribution function.

Until here, all components of the input x have been treated without distinction. Now, two kinds of inputs are distinguished: the environmental inputs x_e on the one hand, and the fire scenario inputs x_s on the other hand (in such a way that $x = (x_e, x_s)$). The environmental inputs can vary in time, which is modeled by a distribution $f^{\text{env}}(x_e)$ representing the variations of the environmental inputs. We would like to express the probability of failure only as a function of the scenario inputs x_s , to assess their impact on the conformity of the building during a fire. Of course, our objective is to estimate the conformity at high-fidelity level, so that the fidelity parameter is fixed to t_{HF} . Then, the probability of failure can be written as

$$P_f(x_s) = \int P_f(x_s, x_e, t_{HF}) f^{\text{env}}(x_e) dx_e. \quad (10)$$

This second type of probability of failure depends on the output and input distributions. Since, $P_f(x_s)$ depends on ξ and λ , $P_f(x_s)$ has a posterior distribution which depends on those of ξ and λ .

The posterior distribution of $P_f(x_s)$ has no closed-form expression. Consequently, numerical approximations must be used. We resort to a Monte-Carlo method. To estimate the probability of failure given a scenario x_s , we begin by drawing randomly n_x values of environmental inputs $(x_{e,i})$ from the distribution f^{env} . Then, we draw randomly n_p sample paths of ξ and λ from the posterior distribution (see Figure 2c). For each path, the probability of failure can be estimated by computing the pointwise probability of failure, and averaging on the environmental input samples:

$$p_j(x_s) = \frac{1}{n_x} \sum_{i=1}^{n_x} P_f^{(j)}(x_{e,i}, x_s, t_{HF}), \quad 1 \leq j \leq n_p. \quad (11)$$

The p_j 's are approximate samples of the posterior distribution of $P_f(x_s)$.

From these samples, a point estimate and an uncertainty can be computed. We use respectively median and 95% confidence interval. The uncertainty of the estimate takes account on the uncertainty of the environmental inputs (when n_x values of environmental inputs are generated), of the output (when pointwise probability of failure are computed), and of the posterior Bayesian model (when n_p sample paths are drawn from posterior distribution). This algorithm has to be applied for each scenario of interest.

Summary of the methodology

Consider a fire safety study which has to be effected on a numerical field model. The numerical model is a stochastic complex simulator, taking physical inputs x and a fidelity input t , and returning an output Z . The inputs x must be the most significant on the output. Other parameters must be set, and are supposed known and constant. The simulations should ideally be run at the level of fidelity t_{HF} . The inputs x are divided in two categories: environmental inputs x_e and scenario inputs x_s , $x = (x_e, x_s)$. A distribution $f^{\text{env}}(x_e)$ on environmental inputs is specified. In order to estimate any quantity of interest, the procedure consists of:

1. Define different levels of fidelity, which must be lower fidelity than t_{HF} . Generate a nested Latin Hypercube Sampling design $(x_i, t_i)_{1 \leq i \leq n}$ according to the computation budget. For each point (x_i, t_i) of the design, run the corresponding simulation, and get the output z_i .
2. Condition the Gaussian process which models the simulator to the observations $(x_i, t_i, z_i)_{1 \leq i \leq n}$.
3. Compute the posterior distribution of the quantity of interest. Deduce an estimate and an uncertainty. For instance, the method described above can be used to estimate a probability of failure.

The methodology has two main advantages. First, as the simulator is described as a black box, the method is not intrusive, and therefore, easily applicable on many types of simulators. Second, the distinction between scenario and environment inputs does not impact the conditioning of Gaussian process. Consequently, once the posterior distribution is computed, the scenario x_s and the distribution $f^{\text{env}}(x_e)$ can be changed as wished. If scenario and/or environment distributions are modified, only the last step of the methodology must be relaunched. This fact allows to draw curves of probabilities of failure according to various configurations.

APPLICATION

Description of the case

To illustrate the methodology, we present an example of fire safety analysis. We consider a real building of the Laboratoire National de métrologie et d'Essais (LNE), which is a 20 m by 12 m by 16 m rectangular parallelepiped. The thickness and the thermal properties of the walls are defined in accordance with the present structure. It has two open doors at the floor level, 2 m by 1 m, and two natural smoke removal systems in the ceiling, 2 m by 2 m (Figures 4a and 4b). The fire source is located in the center of the test hall, and is described by a Heat Release Rate (HRR) evolution in three steps which are growing, steady state and decreasing (Figure 4c). Concerning the boundary conditions, the gas in the computational domain is set still with ambient temperature, and, at the free sides, a static pressure boundary condition is employed.

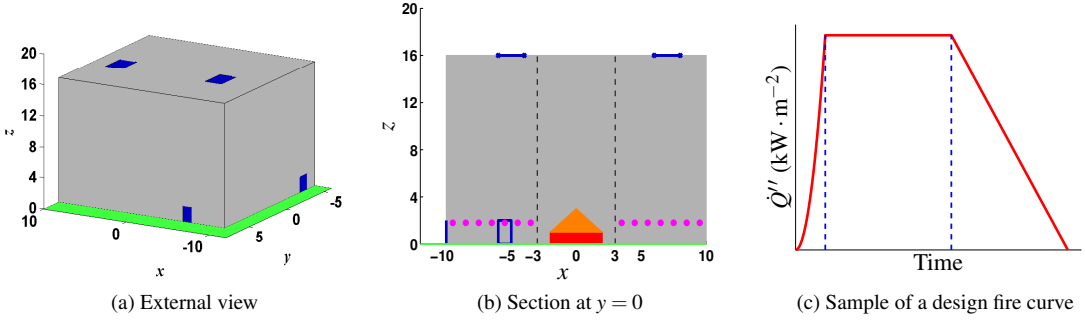


Fig. 4. (a) External view of the considered building. (b) Section of the building. The triangle in the middle is the fire, the bullet points are the sensors. (c) Time evolution of the Heat Release Rate.

The conformity of the smoke control system is estimated by comparing tenability conditions to critical thresholds (see Table 1), determined according to the ISO standard 13571:2012 [21]. The three studied outputs are the radiative heat flux H_f , the temperature at 1.8 m T_c , and the visibility V . A simulation of FDS returns these three outputs. During a simulation, sensors are put in the whole building at 1.8 m above the ground, except above the plume of the fire (see Figure 4b). The distance between the sensors is 1 m. Each sensor returns a time evolution of the three outputs. A post-treatment transforms these time evolutions into one scalar value per output: the values on all sensors are averaged at each time step, and then the extremal value of this time-dependent average is computed (maximum for the heat flux and the temperature, minimum for the visibility). The outputs are analyzed separately.

Outputs	Signification	Failure
H_F ($\text{kW} \cdot \text{m}^{-2}$)	Radiative Heat Flux	$> 2.5 \text{ kW} \cdot \text{m}^{-2}$
T_c ($^{\circ}\text{C}$)	Temperature at 1.8 m	$> 60^{\circ}\text{C}$
V (m)	Visibility	$< 5 \text{ m}$

Table 1. List of outputs studied during this experiment.

The study is carried out using FDS version 5.5.3. It is a CFD model of fire-driven fluid flow, developed by the National Institute of Standards and Technology (NIST), which solves an approximation of the Navier-Stokes equations appropriate for low-Mach number using finite difference methods [22].

Strictly speaking, the FDS-based simulator used in this application is a stochastic simulator (random fields are initialized, inside FDS, with a pseudo-random number generator) that has been turned into a deterministic simulator by fixing the seed of the generator. It can be observed, however, that the outputs of this simulator are still very irregular functions of its physical inputs. This justifies, as explained previously, the use of a stochastic simulator model (more precisely, a model with a nugget effect).

We focus on eight particular input variables, which are the most influential ones. Three of them are considered as environmental inputs (external temperature, atmospheric pressure, and ambient temperature), while the five other are categorized as scenario inputs (fire growth acceleration, fire area, maximal surface power, total surface energy, and soot yield). Moreover, the size t_{xyz} of the mesh cells is used as the fidelity input in our multi-fidelity approach. The time step is determined automatically from the mesh size t_{xyz} by FDS, in order to satisfy the Courant-Friedrichs-Lewy (CFL) stability conditions [22, Section 6.4.7]. See Table 2 for more information.

In this particular study, the number of mesh cells per meter is constrained by the geometry of the building to be an integer (this constrained could be relaxed by using a less regular, adapted mesh). Consequently, the mesh size t_{xyz} must be equal to $100 \text{ cm}/n$, where n is a positive integer. Moreover, based on computation-time

Random inputs		Signification	Distributions
T_{ext}	(°C)	External temperature	$T_{ext} (°C) \sim \mathcal{N}(10, 6.66^2)$
P_{atm}	(bar)	Mean atmospheric pressure	$P_{atm} (\text{bar}) \sim \mathcal{N}(1, 0.00666^2)$
T_{amb}	(°C)	Initial ambient temperature	$T_{amb} (°C) T_{ext} (°C) \sim \mathcal{N}\left(22.5 + 2 \frac{(T_{ext} - 10)}{6.66}, 1.5^2\right)$
Scenario inputs		Signification	Ranges
α	(kW · s ⁻²)	Growth acceleration	[0.011338; 0.2] kW · s ⁻²
A_f	(m ²)	Fire area	[1; 20] m ²
\dot{Q}''	(kW · m ⁻²)	Maximal Heat Release Rate	[300; 600] kW · m ⁻²
q_{fd}	(MJ · m ⁻²)	Total surface released energy	[300; 600] MJ · m ⁻²
Y_{soot}	(kg · kg ⁻¹)	Soot yield	[0.01227; 0.1227] kg · kg ⁻¹
Fidelity input		Signification	Values
t_{xyz}	(cm)	Mesh size	{100; 50; 33; 25; 20} cm

Table 2. List of inputs studied during this experiment. \mathcal{N} is the normal distribution.

Levels	100 cm	50 cm	33 cm	25 cm	20 cm	Total
Number of simulations in the design	270	90	30	10	0	400
Approximate computation time	4.7 min	65 min	6.2 h	21 h	2.3 days	21.4 days
Comparison with level 20 cm	1/700	1/50	1/8.8	1/2.6	1	9.5

Table 3. Description of the levels. Remarks: the computation time depends on the hardware, and the simulations can be parallelized.

considerations, $t_{HF} = 20$ cm is chosen as the highest fidelity level. Five levels of fidelity are thus considered: 100 cm, 50 cm, 100/3 cm (hereafter referred to as 33 cm), 25 cm and 20 cm.

Numerical experiments are performed according to a nested LHS with four levels of observations: 270 simulations at level 100 cm, 90 at 50 cm, 30 at 33 cm, 10 at 25 cm, which correspond to a total computation time of approximately 21.4 days (see Table 3). No observations are made at the highest level $t_{HF} = 20$ cm, which is only used for validation purposes.

Results

This section presents results obtained with our methodology. First, as an example of the kind of results that can be obtained from the posterior distribution. Figure 5 presents the estimates of the three outputs along the fire area A_f axis, at high-fidelity level $t_{xyz} = 20$ cm (other inputs being fixed at $T_{ext} = 10$ °C, $P_{atm} = 1$ bar, $T_{amb} = 22.5$ °C, $\alpha = 0.1057$ kW · s⁻², $\dot{Q}'' = 460$ kW · m⁻², $q_{fd} = 450$ MJ · m⁻² and $Y_{soot} = 0.027$ kg · kg⁻¹). The solid line is the mean function of the posterior distribution, which is our estimator of ξ (defined in Eq. (1)). The inner area around the line is the 95% confidence interval of the estimate. The outer area is the 95% confidence interval of future simulations, which also takes into account the output variability measured by the variance λ (defined in Eq. (1)). We can observe that, as expected, the building becomes more dangerous when the fire area grows.

Let us now consider the evaluation of conformity for some fire scenarios. As written in previous sections, the building conformity can be assessed by the probability of exceeding the safety threshold (Table 1). We let the fire area A_f vary between 1 m² and 20 m², with one scenario every 0.25 m². Other scenario inputs are fixed to $\alpha = 0.1057$ kW · s⁻²; $\dot{Q}'' = 460$ kW · m⁻²; $q_{fd} = 450$ MJ · m⁻²; $Y_{soot} = 0.027$ kg · kg⁻¹. Thus, there are 77 fire scenarios for which the probability of failure is estimated. We use the method described in the previous section to estimate probability of exceeding the threshold, with $n_p = 10^6$ sample paths and $n_x = 700$ values of environmental inputs. Estimates are presented in Figure 6. The solid line is the estimate of the probability of exceeding the threshold at the given scenario. The area is the 95% confidence interval. Once again, we

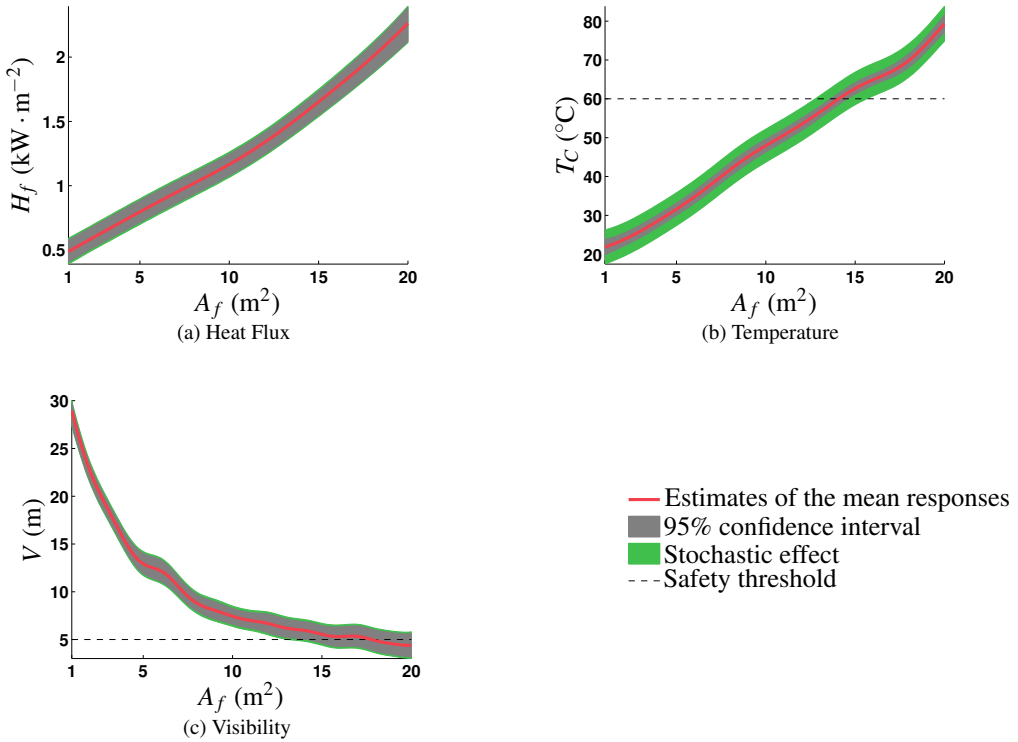


Fig. 5. Estimates of the three outputs as function of the fire area $A_f \in [1; 20]$ m². (a) Heat Flux H_f . (b) Temperature T_c . (c) Visibility V . The safety threshold of H_f is never reached.

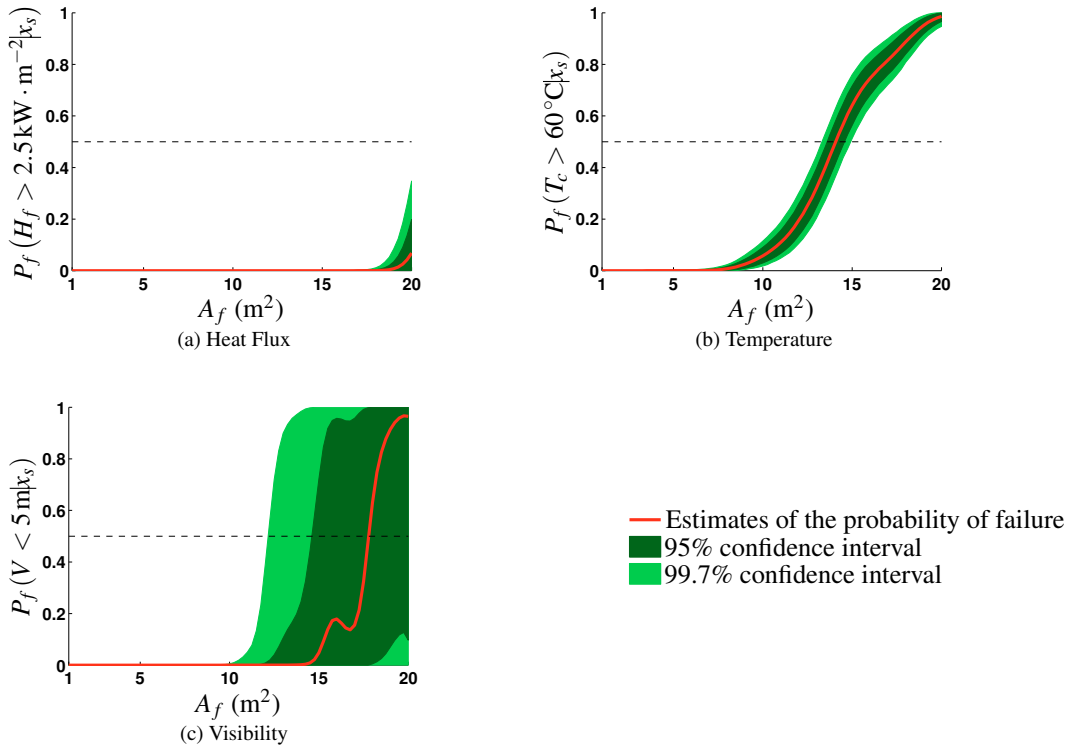


Fig. 6. Estimates of probabilities of failure as function of the scenario variable $A_f \in [1; 20]$ m², the fire area. Each figure corresponds to one output: (a) Heat Flux H_f , (b) Temperature T_c and (c) Visibility V .

find that the place becomes more dangerous when fire becomes bigger. These graphs show also the critical scenario of the building. Thus, other scenario inputs being fixed, the building is safe until the fire area A_f is greater than 12 m^2 . We can also observe on the visibility, Figure 6c, that estimation of probability of failure becomes very inaccurate around 17 m^2 . In fact, the width of the confidence interval becomes equal to 1, implying that nothing can be deduced from the data about the value of the probability P_f at this point. To fix this, additional simulations must be added to reduce the uncertainty. Of course, the inputs of these simulations should be properly selected to provide as much information as possible. The methods for selecting the inputs to observe are called sequential design of experiments.

In this subsection, the presented results depend on the fire area A_f . However, the results (distribution of outputs, probability of failure...) can be analyzed according to any inputs (\dot{Q}' , α ..., or many of them). The Bayesian approach provides an estimate and a corresponding measure of uncertainty for any quantity of interest.

The relevance of our method can be seen by comparing computation time (see Table 3). In fact, we can observe that our multi-level design takes $270 \times \frac{4.7}{60 \times 24} + 90 \times \frac{65}{60 \times 24} + 30 \times \frac{6.2}{24} + 10 \times \frac{21}{24} = 21.4$ days to be evaluated. In such a delay, about $\frac{21.4}{2.3} = 9.5$ high-fidelity simulations could be run, which is too few to draw any conclusion about output values at unobserved locations or probability of failure. The other parts of our procedure (selecting a DOE, building the posterior model, deducing results) last together about one hour.

Validation on additional data

The previous part presented the estimates obtained with a Bayesian multi-fidelity approach. This part compares estimates with outputs of high-fidelity simulations, to assess their accuracy. In fact, there is no simulation on the highest level of fidelity (see Table 3). However, we use our method to study a non-observed level. Consequently, we must ensure that the method can estimate a non-observed level.

To do this, we launch 100 new simulations at 20 cm, the highest fidelity level. These new data are designed according to an optimized maximin LHS, but are not involved in the construction of the conditional Gaussian processes. The previously built posterior distribution estimates the outputs at the inputs of these high-fidelity simulations. Thus, estimates (with uncertainty) can be compared with high-fidelity simulations.

Figure 7 presents the comparison between simulations and estimates at the highest level of fidelity, for the three outputs. On these figures, estimates and simulations are distributed around the $y = x$ line. The estimates seems good on the whole input space. Observe that a few uncertainty bars do not cross the $y = x$ line, which is expected since they are 95% confidence intervals.

The method can also be checked by comparing the estimate of the probability of failure with the estimate obtained with another method. To do it, the following fire scenario is chosen: $\alpha = 0.1057 \text{ kW} \cdot \text{s}^{-2}$, $A_f = 14 \text{ m}^2$, $\dot{Q}' = 460 \text{ kW} \cdot \text{m}^{-2}$, $q_{fd} = 450 \text{ MJ} \cdot \text{m}^{-2}$, $Y_{soot} = 0.027 \text{ kg} \cdot \text{kg}^{-1}$. This fire scenario is present on the curves presented Figure 6. The results obtained by our method for this specific scenario are compared with estimates of the probabilities of failure by Monte-Carlo, using 150 high-fidelity simulations.

Figure 8 presents the different results returned by both methods. For each figure, the left result is obtained with our methodology and the uncertainty is measured as a 95% confidence interval, and the right one is computed with Monte-Carlo and the uncertainty is measured as a 95% confidence interval. For the smoke radiative Heat Flux, both methods estimate a probability of failure at 0. For the Temperature, both results are similar. For the Visibility, our methodology returns a larger 95% interval than the Monte-Carlo method. As emphasized earlier, some additional simulations are needed to learn on the Visibility. In the three cases, the results of both methods are compatible.

Recall that the results presented for our methodology are based on a data set which took 21.4 days to be computed. In contrast, the design used by Monte-Carlo method takes $150 \times 2.3 = 345$ days (see Table 3). So, our method is $\frac{345}{21.4} = 16$ time faster than a direct Monte-Carlo method. Moreover, the Monte-Carlo method gives a probability of failure for one scenario, while our methodology can provide estimates of the probability of failure for any scenario.

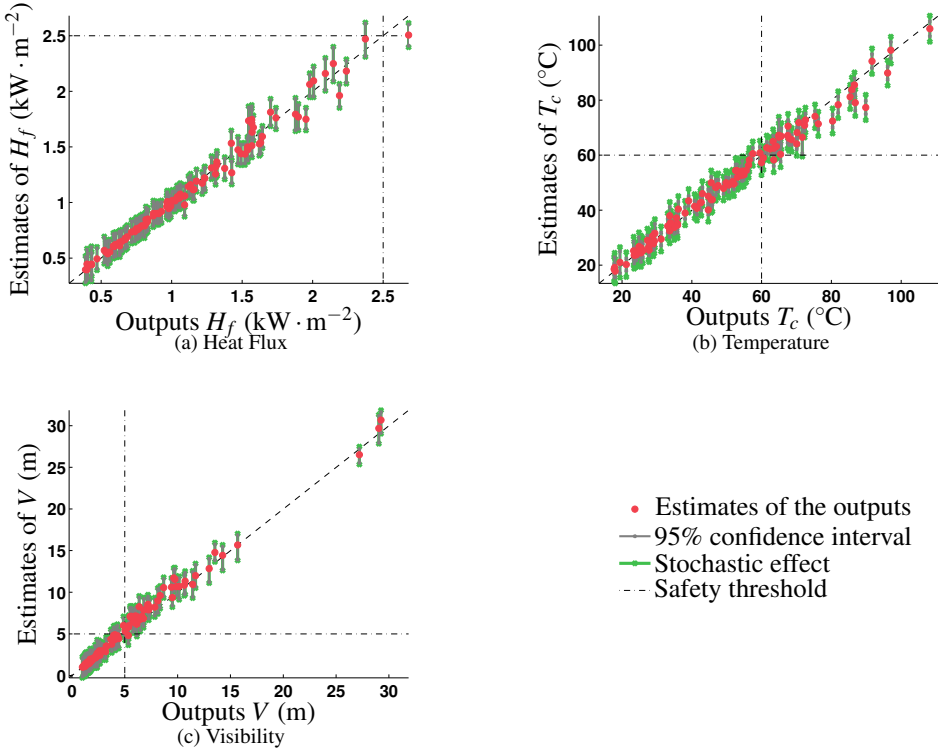


Fig. 7. Comparison between estimates (ordinate) and high-fidelity simulations (abscissa), for the three outputs: (a) Heat Flux H_f , (b) Temperature T_c and (c) Visibility V . The dashed diagonals are the lines $y = x$.

CONCLUSION

This article provides a method to assess the fire safety of a building using a complex, expensive-to-run numerical simulator. The proposed method is non-intrusive, and treats the simulator as a black box, using a Bayesian multi-fidelity approach. More exactly, we use a Gaussian process model to efficiently learn about the simulator with a limited budget of simulation time. The prior Gaussian model, which represents our assumption about the simulator, and observations of the simulator are combined to construct a posterior distribution, which is used to estimate quantities of interest, such as, in this paper, the probability of exceeding a safety threshold.

Because of the complexity of the CFD model, the number of observation is limited. The multi-fidelity approach is a way to reduce the uncertainties of estimates for a given budget computation time. The idea consists in combining fast low-fidelity and slow high-fidelity simulations of the simulator based on a CFD solver to construct the posterior distribution.

The methodology was tested on a fire safety study. By comparing the computation time, the Bayesian multi-fidelity approach was found to be, on this example, one order of magnitude faster than a standard one-level Monte Carlo approach with comparable accuracy. The developed Bayesian models can estimate the expected outputs of the simulator at the highest level of fidelity, and probabilities of failure at given fire scenarios. Moreover, the posterior distributions provide a quantification of uncertainty for the quantities of interest. Our methodology assesses safety quicker and with more confidence than with traditional analyses.

The proposed methodology, being non-intrusive by nature, can in principle be applied to virtually any simulator, and therefore to much more complicated examples than the one used in this paper as an illustration. Some care is required, however, in order to ensure that the underlying prior model faithfully represents available knowledge about the simulator at hand. It is thus important to complement the methodology with tools for model validation (e.g., cross-validation or posterior predictive checks) and assessment of robustness with respect to modeling assumptions. These directions will be investigated in future work.

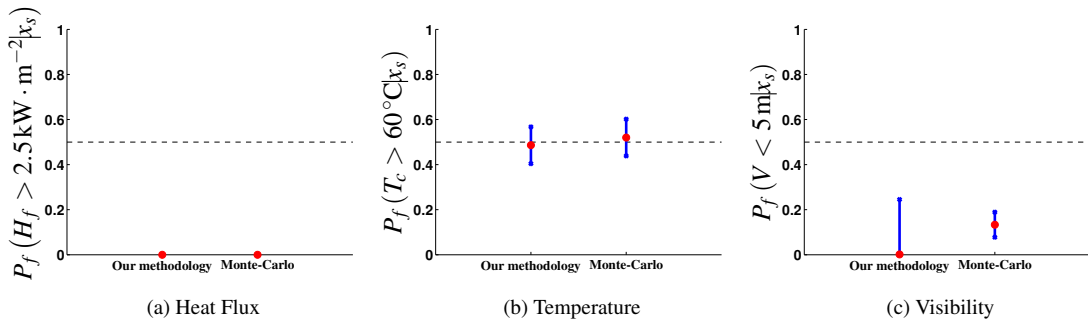


Fig. 8. Comparison between the estimates of probabilities of failure returned by the method (left on figures) and by Monte-Carlo (right on figures). The bullet points are estimates, and the bars are 95% uncertainties.

The methodology can also be improved regarding how the simulation points are selected. Indeed, the design of experiment that we proposed follows general recommendations of the statistical literature, but additional observations of the numerical model could be selected for the specific purpose of improving the estimation of a particular quantity of interest. Future work will thus also focus on the sequential design of experiments in such a multi-fidelity setting.

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