A multivariate interval approach for inverse uncertainty quantification with limited experimental data

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Abstract

This paper introduces an improved version of a novel inverse approach for the quantification of multivariate interval uncertainty for high dimensional models under scarce data availability. Furthermore, a conceptual and practical comparison of the method with the well-established probabilistic framework of Bayesian model updating via Transitional Markov Chain Monte Carlo is presented in the context of the DLR-AIRMOD test structure. First, it is shown that the proposed improvements of the inverse method alleviate the curse of dimensionality of the method with a factor up to 10^5 . Furthermore, the comparison with the Bayesian results revealed that the selection of the most appropriate method depends largely on the desired information and availability of data. In case large amounts of data are available, and/or the analyst desires full (joint)-probabilistic descriptors of the model parameter uncertainty, the Bayesian method is shown to be the most performing. On the other hand however, when such descriptors are not needed (e.g., for worst-case analysis), and only scarce data are available,

Preprint submitted to Mechanical Systems and Signal Processing

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the interval method is shown to deliver more objective and robust bounds on the uncertain parameters. Finally, also suggestions to aid the analyst in selecting the most appropriate method for inverse uncertainty quantification are given. *Keywords:* Multivariate interval uncertainty, Uncertainty Quantification, DLR-AIRMOD, Bayesian model updating, Limited data

1. Introduction

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In general engineering practice, the knowledge on a structure is usually incomplete, be it due to inherent variable model parameters or a lack of knowledge on the true parameter values [1, 2]). Hence, representing these model parameters as deterministic quantities might prove to be inadequate when a reliable and economic design is pursued, as a large degree of conservatism is needed to prevent premature failure and corresponding maintenance or insurance costs. This over-conservatism not only impairs the economic cost of producing the component; it also leads to unnecessary weight increase, which is impermissible in high-performance sectors such as machinery design, aerospace or automotive. In the last few decades, highly advanced techniques including probabilistic [3],

possibilistic [4] or imprecise probabilistic methods [5] have been introduced to include non-determinism efficiently in these design models.

In order for these tools to deliver a realistic quantification of the nondeterminism in the responses of the design model, the description of the nondeterministic parameters of the model should be made objectively and accurately. Since not all parameters (such as e.g. connection stiffness values or heterogeneous material properties) are trivial to measure directly, inverse uncertainty quantification (UQ) techniques have been introduced. Following inverse

²⁰ UQ, the responses of the structure are measured and used to infer knowledge on the non-determinism in the model parameters. As concerns inverse UQ in a probabilistic sense, the class of Bayesian methods is considered the standard approach [6], even for random fields [7]. However, in the context of limited, insufficient, vague or ambiguous data, the prior estimation of the joint probability density function of the non-deterministic parameter values is subjective. Moreover this estimate influences the quantified result to a large extent when insufficient independent measurement data are available.

Inverse UQ methods for the identification and quantification of multivariate interval uncertainty usually minimise a squared \mathcal{L}_2 -norm over the difference

³⁰ between the interval boundaries of respectively a measurement data set and the prediction of the FE model [8, 9]. Application of most of these techniques is prone to ill-conditioning and non-uniqueness when no special care is taken in the definition of the identification problem [10]. Also alternative approaches using Kriging predictor models were introduced recently [11] and compared to stochastic model updating [12].

Recently, a novel methodology for the identification of multivariate interval uncertainty was introduced by some of the authors in [13], with an extension to interval fields in [14] and [15]. This method is based on the convex hull concept for the representation of dependent uncertain output quantities of an

- ⁴⁰ interval FE model, and iteratively minimises the discrepancy between the convex hull of these uncertain output quantities with the convex hull over a set of replicated measurement data. However, since the computation of a convex hull follows an exponential time complexity with its dimension, the dimension of these convex hulls should be reduced as to allow for applying this method to
- ⁴⁵ large-scale problems. Dimension reduction is a topic that is quickly emerging in the fields of big data and machine learning, where datasets are often too highdimensional to be handled directly. In this context, a broad range of techniques based on for instance covariance matrix decompositions [16], manifold learning approaches [17], or active subspace methods [18] have been introduced in recent
- ⁵⁰ years. In the context of the inverse quantification of multi-dimensional interval uncertainty, the application of such dimension reduction methods is an underexplored domain.

Finally, whereas the literature on comparing forward UQ in a probabilistic and non-probabilistic context is abundant (see e.g., [19] or [20] for a more recent treatment), such a practical comparison for inverse approaches is severely lacking in literature. The objective of this paper is therefore twofold. First, an improved version of a recently proposed interval method [13, 14] is presented in the sense that by reducing the dimension of the corresponding convex hulls, more challenging problems can be tackled. In addition, an objective comparison

of Bayesian uncertainty quantification methods [21], which are most commonly applied in a probabilistic context is provided and suggestions for choosing the most appropriate technique based on the data are made.

The paper is structured as follows. Section 2 introduces the extensions to the novel method for the identification and quantification of multivariate interval uncertainty. It is illustrated that the exponential time complexity of computing the objective function is relaxed by projecting the convex hull onto lower-dimensional subspaces of an orthogonal basis with a dimension equal to the effective dimension of the convex hulls. Section 3 presents the reader with the concept of Bayesian uncertainty quantification. Both techniques are critically

⁷⁰ compared and a conceptual comparison is given in section 4. Section 5 presents a case study comparing the applicability of both methods to the well-known DLR-AIRMOD [22] case. Specifically, it is studied how both methods perform in terms of obtained information, computational cost and accuracy, depending on the size of the dataset. Finally, section 6 lists the conclusions of this work.

75 2. Multivariate interval quantification

This section introduces the interval finite element method and the method used for the identification and quantification of multivariate interval uncertainty based on indirect measurement data. In the following, a model parameter θ having interval uncertainty is denoted θ^{I} . Vectors are expressed as lower-case boldface characters θ . Interval parameters are either represented using the

boldface characters θ . Interval parameters are either represented using the bounds of the interval $\theta^I = [\underline{\theta}; \overline{\theta}]$ or the centre point $\hat{\theta} = \frac{\underline{\theta} + \overline{\theta}}{2}$ and the interval radius $r_x = \frac{\overline{\theta} - \underline{\theta}}{2}$.

2.1. The interval finite element method

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Let \mathcal{M} be a deterministic Finite Element model that is used to solve a (set of) differential equations for $\boldsymbol{z}^{\boldsymbol{m}} \in \mathbb{R}^d$ through the vector valued function operator g:

$$\mathcal{M}(\boldsymbol{\theta}): \boldsymbol{z}^{\boldsymbol{m}} = g(\boldsymbol{\theta}), \qquad g: \mathbb{R}^k \mapsto \mathbb{R}^d \tag{1}$$

with $\boldsymbol{\theta} \in \mathcal{F} \subset \mathbb{R}^k$ the vector of model parameters and \mathcal{F} the sub-domain of feasible parameters (e.g., non-negative contact stiffness).

The uncertainty that is attributed to $\boldsymbol{\theta}$ is modeled as an interval vector $\boldsymbol{\theta}^{I} \in \mathcal{F}^{I} \subset \mathbb{IR}^{k}$, with \mathbb{IR}^{k} the space of k-dimensional interval vectors. Note that due to the orthogonality of all θ_{i} , i = 1, ..., k, all model parameters are considered independent by definition. Therefore, $\boldsymbol{\theta}^{I}$ can also be represented as a k-dimensional hypercube.

The interval FE method generally aims at finding a solution set \tilde{z} containing the extreme realizations of z^m given the hyper-cubic parameter uncertainty. In general, \tilde{z} is not hyper-cubic, but spans a non-convex region in \mathbb{R}^d , since \mathcal{M} provides coupling between z_i^m , i = 1, ..., d. Therefore, it is commonly approximated by an uncertain realization set \tilde{z}^m , which is obtained by propagating qwell selected deterministic realizations z_j^m of the hyper-cubic uncertain input parameters θ^I :

$$\tilde{\boldsymbol{z}}^{m} = \left\{ \boldsymbol{z}_{j}^{m} \mid \boldsymbol{z}_{j}^{m} = \mathcal{M}(\boldsymbol{\theta}_{i}); \boldsymbol{\theta}_{i} \in \boldsymbol{\theta}^{I}; i = 1, \dots, q \right\}$$
(2)

In general, $\tilde{z}^m \cong \tilde{z}$ when q is taken sufficiently large, or sampled *intelligently* (e.g., following the Cauchy deviates method [23]). When \mathcal{M} is a strictly monotonic FE model, the Transformation Method [24] provides an exact mapping from θ^I to \tilde{z} , albeit needing 2^k deterministic model evaluations for the solution of a single interval problem.

A mathematical handle to the boundaries of \tilde{z}^m is provided by the convex hull \mathcal{C}^m of the realization set \tilde{z}^m . \mathcal{C}^m can be considered as a set of vertices bounding the uncertain realization set:

$$\mathcal{C}^{m} = \left\{ \sum_{j=1}^{q} \beta_{j} \boldsymbol{z}_{j}^{m} \mid (\forall j : \beta_{j} \ge 0) \land \sum_{j=1}^{q} \beta_{j} = 1 \; ; \; \boldsymbol{z}_{j}^{m} \in \tilde{\boldsymbol{z}}^{m} \right\}$$
(3)

with β a vector of weighting factors, such that all elements β_j are non-negative and sum to one. Such convex hull has also a multi-dimensional volume \mathcal{V}_m . Alternatively, \mathcal{C}^m can also be represented as a set of h_m *d*-dimensional linear inequalities that describe the boundaries of \mathcal{C}^m :

$$\mathcal{C}^m \equiv [f_1, ..., f_{h_m}]^T \equiv \mathbf{A}_m (\boldsymbol{z}^m)^T - \mathbf{b}_m \ge 0$$
(4)

with $\mathbf{A}_m \in \mathbb{R}^{h_m \times d}$, $\mathbf{b}_m \in \mathbb{R}^{h_m}$ and h_m the number of half-spaces f_i , $i = 1, ..., h_m$. An illustration of such convex hull over 12 arbitrary realisations \mathbf{z}_j^m of two responses z_1 and z_2 is given in figure 1.



Figure 1: Convex hull and bounding halfspaces over 12 realisations z_j^m of two arbitrary responses z_1 and z_2 using five half-spaces f_i .

2.2. Multivariate interval quantification

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This section presents an improved version of the method presented in [13, 14]. Specifically, the reduction of the computational expense is extended as to accommodate high-dimensional FE models. First the basis method is introduced for the readers convenience.

As a basis for the inverse quantification of multivariate interval uncertainty, experimental data \mathcal{D} of the model responses \boldsymbol{z} are obtained by experimentally testing t times a physical replica of the numerical model. These measurements are then used to construct a measurement set \tilde{z}^e , and the non-determinism that is present in these replicated measurements is bounded by its convex hull C^e and corresponding d-dimensional volume \mathcal{V}_e in analogy to eq. (3). These computations are performed using the QuickHull algorithm [25].

A specific issue arises when d becomes large (say d > 10). In that case, computation of a single convex hull might take prohibitively long since the worst-case time complexity QuickHull is:

$$\mathcal{O}\left(\lfloor v_c^{\frac{d}{2}} \rfloor / \lfloor \frac{d}{2} \rfloor!\right) \tag{5}$$

with v_c the number of vertices of \mathcal{C}^m [25].

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To accommodate this, the dimension of the vector space \mathbb{R}^d in which these convex hulls are computed, is reduced using the method presented in [14]. This method is briefly recalled in the following paragraphs. Specifically, an orthogonal basis \mathcal{B} is constructed in \mathbb{R}^{d_r} , with $d_r \ll d$. This basis is defined as:

$$\mathcal{B} = \operatorname{span}\{\phi_{e,d-d_r}, \phi_{e,d-d_r+1}, \dots \phi_{e,d}\}$$
(6)

with d_r chosen such that all non-zero dimensions are included in \mathcal{B} , and ϕ_e the eigenvectors corresponding to the d_r largest eigenvalues of the covariance matrix Ξ_e of the measurement data set \tilde{z}^e . These eigenvectors are obtained by following singular value decomposition:

$$\boldsymbol{\Xi}_{\boldsymbol{e}} = \boldsymbol{\Phi}_{\boldsymbol{e}} \boldsymbol{\Lambda}_{\boldsymbol{e}} \boldsymbol{\Phi}_{\boldsymbol{e}}^{T} \tag{7}$$

with $\Lambda_e \in \mathbb{R}^{d \times d}$ the diagonal matrix of the ordered eigenvalues $\lambda_1 \leq \lambda_2 \leq \cdots \leq \lambda_d$ of Ξ_e , and $\Phi_e \in \mathbb{R}^{d \times d}$ a matrix containing the orthogonal eigenvectors $\phi_{e,j} \in \mathbb{R}^d$, $j = 1, \ldots, d$. Standard Matlab routines are used to perform the decomposition, which make use of Blas and LAPACK libraries. Note that since Ξ_e is symmetric and positive definite by definition, the result of eq. (7) is exactly the same as when performing an eigenvalue decomposition. The covariance

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matrix Ξ_e is specifically constructed as:

$$\mathbf{\Xi}_{e} = \begin{bmatrix} var(\mathbf{z}_{e1}^{t}) & cov(\mathbf{z}_{e1}^{t}, \mathbf{z}_{e2}^{t}) & \dots & cov(\mathbf{z}_{e1}^{t}, \mathbf{z}_{ed}^{t}) \\ cov(\mathbf{z}_{e2}^{t}, \mathbf{z}_{e1}^{t}) & var(\mathbf{z}_{e2}^{t}) & \dots & cov(\mathbf{z}_{e2}^{t}, \mathbf{z}_{ed}^{t}) \\ \vdots & & \ddots & \vdots \\ cov(\mathbf{z}_{ed}^{t}, \mathbf{z}_{e1}^{t}) & cov(\mathbf{z}_{e1}^{t}, \mathbf{z}_{e2}^{t}) & \dots & var(\mathbf{z}_{ed}^{t}) \end{bmatrix}$$
(8)

with $\mathbf{z}_{ej}^t \in \mathbb{R}^t$, j = 1, ..., d a vector containing all measured replica for a single response, and t the number of replicated measurements. Finally, the number of needed eigenvectors d_r is then selected as:

$$\sum_{i=1}^{d_r} \frac{\lambda_{e,i}}{tr(\mathbf{\Xi}_e)} \ge 1 - \epsilon \tag{9}$$

- with ϵ the reduction error, which is selected based on its convergence (see [14]). The quantity d_r is defined as the effective dimension of the convex hull. The underlying idea of this concept is that the combination of d system responses not necessarily represents a d-dimensional manifold in \mathbb{R}^d , e.g. due to a high degree of dependence. In that case, d_r is lower than d and a reduction of the dimension is obtained. Note that it is not the statistical correlation between the measured
- replica that is decomposed, but rather the dependence between the model responses. Obviously, when all model responses are completely independent, this reduction does not decrease the dimension of the problem.

As such, instead of computing the convex hulls C^e and C^m directly in \mathbb{R}^d , the constituting sets \tilde{z}^e and \tilde{z}^m are first projected onto \mathcal{B} , and the reduced convex hulls $C^{e,r}$ and $C^{m,r}$ are computed in that basis. Since the quantification is based on minimising the discrepancy between these convex hulls, all computations for the quantification are performed in \mathcal{B} .

However, when realistic numerical models containing thousands of degree'sof-freedom and/or measurement sets consisting of numerous responses are considered, the effective dimension d_r still might be prohibitively large (i.e., $d_r >$ 10). Therefore, as an extension to this previously presented dimension reduction method, it is proposed to further reduce the dimension of the vector space \mathbb{R}^{d_r} in which both convex hulls are defined. Specifically, both $\tilde{\boldsymbol{z}}_m^r$ and $\tilde{\boldsymbol{z}}_e^r$ are further projected onto d_r^+ -dimensional subspaces, defined by a lower-dimensional orthogonal basis $\mathcal{B}_i^+ \subset \mathcal{B}, \ i = 1, \ldots, \binom{d_r}{d_r^+}$, constructed as a subset of \mathcal{B} , with $d_r^+ \ll d_r$ and $\binom{d_r}{d_r^+}$ the binomial coefficient. Specifically, the i^{th} orthogonal subspace basis \mathcal{B}_i^+ is defined as:

$$\mathcal{B}_i^+ = \operatorname{span}\{\phi_{m,\mathcal{I}_i(1)}, \phi_{m\mathcal{I}_i(2)}, \dots \phi_{m,\mathcal{I}_i(d_r^+)}\}$$
(10)

with \mathcal{I}_i an index set containing the d_r^+ indices for the i^{th} , $i = 1, ..., \binom{d_r}{d_r^+}$ subspace of \mathcal{B} . Then, the reduced sets $\tilde{\boldsymbol{z}}_e^r$ and $\tilde{\boldsymbol{z}}_m^r$ are projected on these subspaces, i.e. $\tilde{\boldsymbol{z}}_{\mathcal{B}_i^+}^e$ and $\tilde{\boldsymbol{z}}_{\mathcal{B}_i^+}^m$, and their respective convex hulls $\mathcal{C}_{\mathcal{B}_i^+}^e$ and $\mathcal{C}_{\mathcal{B}_i^+}^m$ with corresponding d_r^+ -dimensional volumes \mathcal{V}_e and \mathcal{V}_m are computed using the functionality of the QHULL library. By applying this reduction scheme, the time complexity of the computation of the convex hull, which is originally defined as in eq. (5), becomes:

$$\mathcal{O}\left(\lfloor v_c^{\frac{d_r^+}{2}} \rfloor / \lfloor \frac{d_r^+}{2} \rfloor! \times \begin{pmatrix} d_r \\ d_r^+ \end{pmatrix}\right) \tag{11}$$

Therefore, as long as d_r^+ is a very small number, the computational cost of 130 calculating $\begin{pmatrix} d_r \\ d_r^+ \end{pmatrix} d_r^+$ dimensional convex hulls is considerably smaller than computing a single d_r dimensional convex hull. This projection however only retains the $d_r^+ - 1$ order interactions between model responses. Herein, $d_r^+ = 1$ is the limit case where only the hyper-cubic approximation of $\tilde{\boldsymbol{z}}_m^r$ and $\tilde{\boldsymbol{z}}_e^r$ is retained. This however does not limit the accuracy of the method as long as $d_r^+ \ge 2$, 135 since by considering only the convex hulls all higher-order interactions between model responses are already linearised. Therefore, no further approximation of the dependence structure is made.

Finally, the multivariate interval uncertainty in θ^{I} is obtained by minimizing following objective function:

$$\delta(\boldsymbol{\theta}^{I}) = \sum_{i=1}^{\binom{d_{r}}{d_{r}^{+}}} \left(\Delta V_{m,i}^{2} + w_{o} \Delta V_{o,i}^{2} + \Delta c_{i}^{2} \right)$$
(12)

with:

$$\Delta V_{m,i} = 1 - \frac{\mathcal{V}_{m,i}(\boldsymbol{\theta}^I)}{\mathcal{V}_{e,i}}$$
(13a)

$$\Delta V_{o,i} = 1 - \frac{\mathcal{V}_{o,i}(\boldsymbol{\theta}^I)}{\mathcal{V}_{e,i}}$$
(13b)

$$\Delta c_i = \left\| \mathbf{c}_{e,i} - \mathbf{c}_{m,i}(\boldsymbol{\theta}^I) \right\|_2$$
(13c)

where $\mathcal{V}_{m,i}$ and $\mathcal{V}_{e,i}$ are the d_r^+ -dimensional volumes of respectively $\mathcal{C}_{\mathcal{B}_i^+}^m$ and $\mathcal{C}_{\mathcal{B}_i^+}^e$, $\mathcal{V}_{o,i}$ is the volume of the overlap of both convex hulls, and $\mathbf{c}_{e,i}$ and $\mathbf{c}_{m,i}$ are the centres of gravity of respectively $\mathcal{C}_{\mathcal{B}_i^+}^e$ and $\mathcal{C}_{\mathcal{B}_i^+}^m$. Note that for notational simplicity, the subscript \mathcal{B}_i^+ is simplified to i.

The advantage of this methodology for reducing the dimension over selecting a smaller value for d_r , is that more information on the dependence and nondeterminism that is present in the model quantities is retained in the analysis. ¹⁴⁵ Instead of allowing a larger ϵ , and hence more loss of information, this method decomposes the convex hull into lower-dimensional subsets, and quantifies the interval uncertainty based on a combinatorial combination of the discrepancy of those subsets. As a final note, it should be stressed that also other, non-linear, manifold learning techniques such as kernel principal component analysis [26],

- isomaps [27], or local linear embedding [28] have been presented in literature. These methods usually provide a better estimate of the "effective" manifold spanned by the measurement data, however at a higher computational cost. Nonetheless, since only the convex hull of the data is considered, application of these techniques is not expected to improve the quantification.
- Finally, due the construction of the objective function, it jointly tries to minimise the difference between the multidimensional volumes between the convex hull of the measurement data set and the convex hull resulting from the interval FE analysis, as to maximise the overlap between these convex hulls. As the set of measurement data generally does not contain the same extreme vertices
- as contained in the interval model, the volumes cannot be matched exactly. Therefore, a trade-off is found by the optimiser, where the volume of the convex hulls is matched at the price of losing some overlap. Increasing the factor w_o

alleviates this problem, however at the cost of increasing the over-conservatism of the prediction (see e.g., [14]) and making the optimisation problem harder to solve due to the barrier-function like behaviour of the term describing the overlap.

3. Bayesian model updating

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The use of Bayesian methods for uncertainty quantification is largely founded on the the pioneering work of Beck and Katafygiotis [6, 21] in the late 1990s. Following the Bayesian interpretation of probability, the probabilistic nature of an uncertain parameter is interpreted as the degree to which it is believed that each possible value of this parameter is consistent with the available information (e.g., the response of a high-fidelity model or measurement data). Following Bayes' rule, this degree of belief is adjusted using independent information. As such, the Bayesian methods translate this prior knowledge on the uncertainty corresponding to the parameter values to an updated posterior knowledge, based

on experimental data.

As a first step in the uncertainty characterisation, a prior probability distribution $p(\boldsymbol{\Theta}|\mathcal{M})$, conditioned upon a chosen mathematical model \mathcal{M} , is assigned to a set of uncertain parameters $\boldsymbol{\Theta}$. These distributions represent prior information on the uncertain parameters, stemming from e.g., expert opinions, lab-scale specimen testing, previous quantification of the uncertainty, etc. Then, experimental data \mathcal{D} are used to update this prior knowledge by means of Bayes' theorem, yielding the posterior distribution $p(\boldsymbol{\Theta}|\mathcal{D}, \mathcal{M})$:

$$p(\boldsymbol{\Theta}|\mathcal{D}, \mathcal{M}) = \frac{p(\mathcal{D}|\boldsymbol{\Theta}, \mathcal{M}) p(\boldsymbol{\Theta}|\mathcal{M})}{p(\mathcal{D}|\mathcal{M})}$$
(14)

where $p(\mathcal{D}|\Theta, \mathcal{M})$ is the likelihood of obtaining the data \mathcal{D} , given the value of the uncertain parameters Θ and a specific model \mathcal{M} . The denominator of eq. (14), also commonly referred to as evidence, ensures that the posterior distribution $p(\Theta|\mathcal{D}, \mathcal{M})$ integrates to one.

In the context of structural dynamics, the data \mathcal{D} usually consists of the residuals between experimental measurements and predictions of the model \mathcal{M}

as:

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$$\epsilon_i = z_i^e - z_i^m(\mathbf{\Theta}), \quad i = 1, \dots, n \tag{15}$$

where $z_i^e = \omega_i^2$ is the square of the *i*-th measured response, $z_i^m(\Theta) = \lambda_i(\Theta)$ is the *i*-th predicted response of a finite element model and *n* the number of considered responses. In the specific context of parameter uncertainty quantification using experimental modal analysis data, z_i^e and z_i^m correspond to measured and predicted eigenfrequencies of \mathcal{M} .

In practice, the likelihood function is often chosen to be a zero-mean multivariate normal distribution [29]:

$$p\left(\mathcal{D}|\boldsymbol{\Theta},\mathcal{M}\right) = \prod_{i=1}^{N} \frac{1}{(2\pi)^{n/2} |\Sigma|^{1/2}} \exp\left(-\frac{1}{2} \epsilon_{i}^{T}(\boldsymbol{\Theta}) \boldsymbol{\Sigma}^{-1} \epsilon_{i}(\boldsymbol{\Theta})\right)$$
(16)

where N denotes the number of data points in \mathcal{D} . The solution to eq. (14) is commonly approximated by sampling from a Markov Chain that is ergodic and stationary with respect to $p(\mathcal{D}|\Theta, \mathcal{M})$. In this paper, the transitional Markov Chain Monte Carlo (TMCMC) method, as introduced by Ching and Chen [30], is applied for the computation of the posterior distribution.

4. Conceptual comparison of Bayesian and interval methods for inverse UQ

The methods presented in section 2 and 3 can both be used for inverse un-¹⁹⁵ certainty quantification, but the underlying philosophy differs greatly. Similarly to forward uncertainty quantification, probabilistic and interval approaches are complementary, and the most appropriate method given an inverse UQ problem is highly case dependent. Important contrasts, next to the obvious philosophical differences, include the availability of objective and informative prior

information, and even more so, measurement data of the model responses and the desired information content on the uncertain model parameters. This section aims at guiding the analyst through the selection of the most appropriate method. Interval methods are based on the idea that the possible values of the un-²⁰⁵ certain parameter are located between two crisp boundaries, without making inference on the likelihood of each value within that interval. As such, interval UQ methods approach the problem from the *outside*, as they usually aim to bound all possible response of the model in an objective sense following a settheoretical approach. Consequently, only a worst-case inference is attainable ²¹⁰ based on the obtained information. Bayesian methods on the other hand start from assigning a degree of plausibility to each value of the uncertain parameters within a range and employ independent data to infer the most plausible parameter values based on Bayes' theorem. Therefore, Bayesian methods approach the uncertainty from the *inside* by searching the most probable point, and specifying how the plausibility of obtaining that specific value decreases for

4.1. Criticism on the Bayesian approach

other parameter values.

When applying Bayesian techniques, a full description of the joint degree of plausibility of each parameter over a range is obtained, even when such parameter is modelled by a random field. Also the selection of the most appropriate covariance function is possible in this context [31]. However, in order to obtain an objective quantification of the parameters and their correlation, a sufficiently accurate prior estimate of the uncertainty is needed in conjunction with sufficient and informative experimental data. When only scarce, limited or vague experimental data are available, as is usually the case in realistic experimental cases, the prior distribution influences the obtained quantification to a large extent. The obtained results are in this case highly subjective, which limits their credibility. This effect is further amplified when the prior is highly biased with respect to the actual parameter values [32]. The effect of adding more data to

the quantification method in order to minimise the effect of the prior can be understood from the following simple example. Consider the tossing of a biased dice that favours the numbers 4 and 5. The analyst, not knowing about the bias on the dice, has a prior assumption that the probability of rolling one of each number in the set $\{1, 2, 3, 4, 5, 6\}$ is equal. When the analyst starts rolling the

biased dice, more and more data come available showing the bias of the dice towards 4 and 5. As such, the effect of the prior assumption is counteracted. One can see that the posterior information will tend to the true biased distribution of the dice when more data becomes available.

In order to obtain a more objective prior, techniques such as the maximum entropy principle have been introduced. However objective and based on a solid mathematical foundation, it is not guaranteed that the maximum entropy principle also yields the physically most probable prior distribution. For example, when only information on the range of the uncertain parameter is available, the maximum entropy principle yields the uniform distribution as being most

²⁴⁵ appropriate as this distribution has the highest information entropy given the available data. However, given only information on the range of the measured responses, there exists no physical guarantee that each point between the predefined bounds is indeed equally probable. Therefore, while being objective, the obtained posterior distribution will not necessarily correspond to reality.

250 4.2. Criticism on the Interval approach

Interval quantification methods on the other hand deliver crisp bounds between which the uncertain parameter is believed to lie. The main advantage hereof is that each data-set can be uniquely described by an interval. Therefore, this method is inherently objective as no subjective estimates and approxima-

- tions on the underlying probabilistic nature of the non-determinism are made to steer the quantification process. However, it is not possible to asses the reliability of the designed structure using purely interval-based methods, as no information on the relative likelihood of different parameter or response values is given. Moreover, intervals are by definition not capable of describing depen-
- dency between different variables. As such, the identified non-determinism is inherently decoupled as well, and thus no information on the dependence can be obtained from the analysis. Finally, as long as the measurement data are obtained in an optimized way (i.e., such that they completely capture the needed

scatter in the responses that results from the uncertainty that is studied), interval methods provide sufficient information for worst-case analysis.

4.3. Bayesian or interval?

In case large data sets are available or the analyst has need for quantifying the relative likelihood of several parameter values being realised, including their (joint)-plausibility, correlation and multi-modal descriptors, Bayesian methods have the upper hand over interval approaches, but the analyst should ensure that sufficient informative data are available as discussed in section 4.1. On the other hand, when data are vague or scarce, interval methods are expected to provide a more objective and accurate quantification of the uncertainty, as less a priori assumptions on the underlying likelihood structure are needed. This however is achieved at the cost that only worst-case information is delivered to the analyst. This information is summarised in figure 2.



Figure 2: Selection flowchart to aid the analyst choosing the most appropriate method

5. Application: the DLR-AIRMOD test structure

5.1. Introduction

The first goal of this case study is to validate the novel reduction method that is presented in the context of the multivariate interval quantification. Secondly, it also aims at providing an in-depth discussion concerning the applicability of Bayesian and interval methods on a realistic dataset. The DLR-AIRMOD structure, with corresponding dataset (see [22]) is selected due to its challenging nature and the elaborate literature on the subject. Both presented techniques

- ²⁸⁵ are applied to this case, and the results are compared in terms of obtained information and accuracy. The inverse UQ is first performed using the complete experimental data set containing 86 measured eigenfrequencies. Then, to illustrate the performance of both methods under scarce data, the inverse UQ is also performed using only 5 measured replica that were randomly selected from the
- full dataset. Note that, since both methods are conceptually very different, a quantitative comparison of the accuracy of the obtained results is not possible. Therefore, the comparison is made qualitatively in terms of conservatism with respect to the measurement data set.

Finally, since both methods require a large number of function evaluations, Artificial Neural Networks are used as surrogate models. Specifically a set of 2-layer (18:16:14:1) Neural Networks that map each vector of uncertain model parameter to one eigenfrequency of the FE model is constructed using state-ofthe art ANN tools and ensuring proper performance (see also [33] or [34] for a discussion of the training and validation).

300 5.2. Model introduction

The DLR AIRMOD structure, as shown in figure 3, is a scaled replica of the GARTEUR SM-AG19 benchmark airplane model [22]. The physical AIRMOD structure is constructed from six aluminum beams that are connected by five bolted joints and weighs approximatly 40 kg to represent the fuselage, wings, winglets, vertical tail plate (VTP) and horizontal tail plate (HTP). It has a wing span of 2.0 m, the fuselage length is 1.5 m and the height is 0.46 m. The complete FE model, constructed in NX Nastran, consists of 1440 CHEXA, 6 CPENTA and 561 CELAS1, 55 CMASS1, 18 CONM2 and 3 CROD elements, and is constructed after [22].

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A set of 18 parameters including support and joint stiffness values, as well as

mass parameters are selected for the identification (see table 1), in correspondence with literature on the subject [12]. The locations of these parameters are indicated in figure 3.



Figure 3: Illustration of the AIRMOD test structure (adapted after [22])

Table 2 lists the eigenmodes and eigenfrequencies that are obtained by solving the deterministic model (denoted as $f_{FE,init}$). From this table, it can be noted that the model indeed exhibits some challenges for uncertainty quantification procedures such as asymmetric modal behaviour and closely spaced modes. To prevent that perturbations in the model parameters lead to eigenmode veering or cross-over, MAC-based mode tracking with respect to the reference solution is applied in each stage of the quantification. From this set, the $1^{st} - 8^{th}$, $10^{th} - 12^{th}$, 14^{th} , 19^{th} and 20^{th} mode are selected for the identification. These 14 modes are selected to be consistent with literature on the subject [12].

It can be shown that the eigenfrequencies predicted by a linear numerical model are a strict monotonous function of the uncertain model parameters [35]. In this case, the reduced transformation method [24] yields the exact bounds on the eigenfrequencies. However, the number of necessary function evaluations for the propagation scales exponentially with the number of uncertain model parameters. For the multivariate interval quantification method, it is therefore assumed that the masses at both wing-tips (i.e. θ_9 and θ_{10}) and the stiffness

	Type	Description	Orientation	Deterministic value	
θ_1	Stiffness	Support stiffness	y	$1.80 \cdot 10^{03} N/m$	
θ_2	Stiffness	Support stiffness	y	$7.50 \cdot 10^{03} N/m$	
θ_3	Stiffness	Cables	y	$1.30\cdot 10^{02}~N/m$	
$ heta_4$	Stiffness	Cables	y	$7.00 \cdot 10^{01} \ N/m$	
θ_5	Stiffness	Cables	y	$7.00 \cdot 10^{01} \ N/m$	
θ_6	Stiffness	Joint stiffness	x,y	$1.00\cdot 10^{07}~N/m$	
θ_7	Stiffness	Joint stiffness	z	$1.00 \cdot 10^{09} \ N/m$	
θ_8	Mass	Cables	/	$2.00 \cdot 10^{-01} \ kg$	
$ heta_9$	Mass	Screws	/	$1.86 \cdot 10^{-01} \ kg$	
θ_{10}	Mass	Screws	/	$1.86 \cdot 10^{-01} \ kg$	
θ_{11}	Mass	Cables	/	$1.50 \cdot 10^{-02} \ kg$	
θ_{12}	Mass	Cables	/	$1.50 \cdot 10^{-02} \ kg$	
θ_{13}	Mass	Cables	/	$1.50 \cdot 10^{-02} \ kg$	
θ_{14}	Stiffness	Joint stiffness	х	$2.00\cdot 10^{07}~N/m$	
θ_{15}	Stiffness	Joint stiffness	У	$2.00\cdot 10^{07}~N/m$	
θ_{16}	Stiffness	Joint stiffness	Z	$7.00 \cdot 10^{06} N/m$	
θ_{17}	Stiffness	Joint stiffness	х	$5.00 \cdot 10^{07} \ N/m$	
θ_{18}	Stiffness	Joint stiffness	У	$1.00\cdot 10^{07}~N/m$	

Table 1: Parameters that are used in the identification

- introduced by the cables at the top and the bottom of the structure (θ_4 and θ_5) are completely dependent, reducing the number of uncertain parameters to 16. Hence, the number of necessary function evaluations for a single interval computation reduces from 262144 to 65536 which makes it computationally more tractable. The validity of this assumption will be estimated based on
- the obtained results. This assumption is not made for the Bayesian method, as the Monte Carlo sampling that underlies the applied TMCMC approach is dimension-independent.

5.3. Measurement data set

A measurement data set containing 86 measurements of the 30 eigenmodes and -frequencies, as explained in detail in [22], is applied for the inverse UQ. For a complete explanation of the experimental campaign that was followed to construct this dataset, the reader is referred to [22]. The measured eigenfrequencies and their variability are shown in Table 2. In this table μ_f and σ_f denote respectively the mean and standard deviation of the measurement data.

To assess the robustness of both methods against the size of measurement data, both the full measurement data set and a subset of 5 arbitrarily drawn replica were employed for the UQ of the 18 parameters.

5.4. Multivariate interval quantification

5.4.1. Response set dimension reduction

- The number of considered model responses proves to be challenging in the application of the interval analysis due to the exponential time complexity of the applied QuickHull algorithm. The computation of a convex hull over all 14 pre-selected resonance frequencies, would therefore make the identification computationally intractable, as this computation has to be performed numerously during the quantification procedure. Therefore, this dimension is reduced following the method presented in section 2.2. In this context, application of equations (8) to (9) indicates that the effective dimensionality is equal to 13 when an approximation error ϵ below 0.1% is desired. This also can be seen in figure 4, which plots the reduction error ϵ as a function of the reduced dimension
- 360 d_r for the first 18 $\lambda_{e,i}$.

Based on the singular value decomposition of the covariance matrix (see eq. (8)) of the pre-selected 14 eigenmodes, an orthogonal basis $\mathcal{B} \in \mathbb{R}^{13}$ is constructed according to eq. (6), and these 14 eigenmodes are projected onto this basis. This dataset is further projected onto lower-dimensional bases, in accor-

dance with eq. (10). The computational gain of applying $d_r^+ = 2$ dimensional sub-bases of \mathcal{B} is illustrated in figure 5. This figure shows the wall-clock time that is needed to compute 1 d_r -dimensional convex hull, as well as the time needed

Table 2: eigenmodes and corresponding initial FE estimate of the corresponding eigenfrequency, as well as the statistics of the measured eigenfrequencies. The eigenmode naming convention is as follows. The letter before the dash – denotes the considered part, the letters after the dash the deformation mode. RB denotes the rigid body modes, W denotes the wing, Wl the winlets, F the fuselage, VTP the Vertical Tail Piece and HTP the Horizontal Tail Piece. B denotes horizontal bending, T torsion, FAB denotes fore-after bending, LB lateral bending, BL bending left, BR bending right and VB vertical bending.

No	Mode	$f_{FE,init}$	μ_f (Hz)	σ_f (Hz)	$\underline{f}(Hz)$	$\overline{f}(Hz)$	nsam
1	RB-Yaw	_	0.23	0.006	0.22	0.24	41
2	RB-Roll	0.56	0.65	0.019	0.6	0.68	81
3	RB-Pitch	0.82	0.83	0.017	0.8	0.88	83
4	RB-Heave	2.14	2.17	0.024	2.11	2.22	86
5	2^{nd} W-B	5.65	5.5	0.004	5.49	5.52	86
6	3^{rd} W-B	15.11	14.91	0.017	14.88	14.94	86
7	1^{st} as sym. W-T	31.31	31.96	0.02	31.92	32.01	86
8	1^{st} sym. W-T	33.62	32.33	0.017	32.29	32.38	86
9	1^{st} VTP-B	35.39	34.38	0.081	34.23	34.54	86
10	4^{th} W-B	44.66	43.89	0.015	43.85	43.92	86
11	1^{st} W-FAB	47.21	46.71	0.149	46.27	46.99	86
12	2^{nd} W-FAB	52.91	51.88	0.012	51.84	51.91	86
13	5^{th} W-B	60.59	58.59	0.075	58.33	58.76	86
14	1^{st} VTP-T	67.69	65.93	0.274	65.46	66.33	86
15	2^{nd} F-LB	102.59	100.05	0.28	99.38	100.48	86
16	2^{nd} VTP-B	128.62	124.56	0.356	123.85	125.1	86
17	6^{th} W-B	132.08	129.38	0.107	129.12	129.66	86
18	7^{th} W-B	145.91	141.47	0.347	140.79	142.76	85
19	2^{nd} HTP-B	206.73	205.59	1.023	203.24	206.87	86
20	1^{st} HTP-FAB	225.73	219.07	1.663	216.29	221.3	86
21	1^{st} W-BR	261.53	254.73	0.557	253.41	256.48	70
22	1^{st} W-BL	262.64	255.02	0.575	253.84	256.27	81
23	3^{rd} W-FAB	278.71	272.08	0.374	271.17	272.89	86
24	1^{st} Wl-BL	320.15	303.96	1.115	301.52	306.5	82
25	1^{st} Wl-BR	321.64	304.32	2.138	301.17	310.27	83
26	3^{rd} F-LB	324.12	313.68	1.218	309.38	314.76	86
27	2^{nd} sym. W-T	336.31	328.55	0.448	327.03	330.21	85
28	2^{nd} as sym. W-T	341.15	331.18	0.528	329.64	332.95	86
29	4^{th} W-FAB	343.55	336.21	0.647	330.88	337.41	86
30	2nd F-VB	359.54	348.68	1.141	346.75	350.3	86



Figure 4: Reduction error ϵ as a function of the reduced dimension d_r , applied to the AIRMOD data set.

to compute $\binom{d_r}{2}$ 2-dimensional convex hulls. The computations are made using the measurement data set with the 14 pre-selected Eigenmodes, containing a selection of either 25, 50 or 86 replica (i.e., the full dataset). All computations are made using a single-thread of an Intel Xeon E5-1620 @ 3.70 GHz with a total 32Gb of RAM as to prevent swapping. As can be seen, the projection method is very effective in reducing the computational burden of obtaining the convex hull. Specifically, when considering a 14-dimensional convex hull, the computational cost is reduced with a factor up to $1 \cdot 10^{05}$ when compared to the computation of the full-dimensional convex hull for $d_r = 14$. Finally, the influence of the measurement data size is considerably smaller when $d_r^+ = 2$, which is very relevant for reducing the result of the interval FE model, which contains $2^{16} = 65536$ responses.

Note that this does not limit the accuracy of the method since quadratic and higher order interactions between responses are already linearised by considering only the convex hull of the set. All interval quantification computations are therefore performed using 91 2-dimensional projections of the full 14-dimensional convex hull, where an equal weight is attributed to each projection.



Figure 5: Wall-clock time that is needed to compute 1 d_r -dimensional convex hull, as well as the time needed to compute $\binom{d_r}{d_r^+} d_r^+$ dimensional convex hulls

5.4.2. Uncertainty Quantification

The objective function presented in eq. (12) is in general not convex and should be solved in 32-dimensional space. Therefore, eq. (12) is solved using a hybrid Particle Swarm Algorithm (PSA). A swarm size of 100 particles is used, and the optimization is considered to be converged when it reached 15 stalling 390 iterations. These settings are found in a heuristic approach and based on prior experience with PSA. In a second step a sequential quadratic programming approach is followed until converged, starting from the estimate of the global minimum of eq. (12) obtained by the PSA. The optimality of this solution can be verified by qualitatively comparing cross-sections of the convex hulls of 395 the measurement data set and the result of the interval FE model. For the quantification of the multivariate interval uncertainty the weighting factor w_{o} in eq. (12) is set to unity, meaning that equal priority is given to maximising the overlap and matching the multidimensional volumes. Given the already challenging nature of the optimisation, this weight value prevents the encounter 400

of instabilities. These would inevitably appear in the case a large w_o is selected due to the then discontinuities behaviour of eq. (12) as a result of the barrierfunction like influence term $w_o \Delta V_{o,i}^2$.

5.5. Bayesian model updating

- The Bayesian uncertainty quantification was performed using 18 uncorrelated marginal uniform prior distributions. The range of the distribution of each parameter has been selected spanning an interval from 5% to 200% of the parameter nominal value. The likelihood, as introduced in eq. (16), is constructed under the assumption independence of the data.
- $_{410}$ 5.6. UQ with full data set

Bayesian result is illustrated in figure 7.

Figure 6 illustrates the marginal posterior distributions of the model parameters obtained via Bayesian updating, and the corresponding quantified interval bounds. For the sake of comparison, the parameters $\boldsymbol{\theta}$ are here normalised by their initial nominal values. The colour coding represents the normalised height of the histogram of each normalised θ_i value according to the obtained marginal posterior. On the other hand, the upward and downward facing triangles denote respectively the lower and upper interval bounds, as obtained by the multivariate interval quantification method. The correlation matrix of the

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First, a large conceptual difference between both methods can be noted. On the one hand, the intervals provide the analyst with crisp bounds for the possible parameter values, on the other hand, the marginal posterior distributions assign a non-zero plausibility to a large range of responses. Note that the range of illustrated values is limited due to the finite data set that is used to plot the

⁴²⁵ normalised histograms. Nevertheless, the most probable point in the Bayesian analysis is still clearly identifiable in most cases. It should be noted that the Bayesian analysis predicts a multi-modal plausibility distribution for θ_1 , θ_2 , θ_4 , θ_5 , θ_7 , θ_8 , θ_9 , θ_{10} , θ_{12} and θ_{15} . The multi-modality can be interpreted as slight changes in the test conditions that were not noted during the experimental



Figure 6: Normalised histograms of the posterior distribution samples, indicated as colour coded bar graphs, and interval bounds, indicated as triangles that are obtained by performing respectively Bayesian inference and Multivariate interval quantification on the full AIRMOD data set.

campaign (see e.g. [33]). This multi-modal behaviour cannot be captured by the interval UQ. By comparing the results for those parameters with multi-modal posteriors, the interval analysis either encompasses all modal peaks (e.g., θ_7), or just the most probable one (e.g., θ_{12} or θ_{15}).

Qualitatively, the predictions of both methods correspond rather well, as the ⁴³⁵ interval method usually bounds the same area in parameter space to which the Bayesian method assigns the highest degree of plausibility. Exceptions hereto exists for θ_6 , θ_8 , θ_{16} , θ_{17} and θ_{18} . These differences are due to a variety of factors, such as the difficulty in finding the exact global minimum in a 18dimensional uncertain space in case of the interval method, a highly peaked ⁴⁴⁰ marginal distributions in the case of Bayesian UQ, or a low-importance input that cannot be effectively updated with the available experiments. As a last

remark, the assumption that parameters 4-5 and 9-10 are dependent to

reduce the number of uncertain parameters in the case of the interval analysis is proven to be reasonable, as the marginal posteriors obtained by Bayesian UQ show similar plausibility distributions and they were found to be highly

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Figure 7: Correlation matrix, obtained by the Bayesian analysis on the full dataset.

From the point of view of the updated model outputs, Figure 8 shows all combinations of the pre-selected eigenfrequencies, obtained by propagating the quantified intervals and Bayesian posterior distribution through their corresponding ANN surrogates of the AIRMOD FE model. As an additional comparison, the measurement data set is also shown.

Marginally, the results obtained by the multivariate interval quantification method predict a wider range of possible responses for f_5 , f_7 , f_8 , f_{10} and f_{12} as compared to the samples from the Bayesian posterior distribution. Note that

these bounds, as opposed to the ones predicted by these Bayesian samples, are considered strict. For e.g., f_{19} and f_{20} , it is shown that some of the experimental responses are not encapsulated within the convex hull. This is a direct result of the construction of the objective function for the quantification, as elaborated in the final paragraph of section 2.2. This effect is clearly visible in this case since $w_o = 1$ to smooth the objective function. As concerns the combination of the 5-8, 7-8 and 8-12 eigenmodes, the mutual dependence is not predicted correctly by the quantified interval model. The interval model in fact predicts these responses to be highly dependent, whereas the measurement data show otherwise. Keeping in mind that the 7th, 8th and 12th eigenmode correspond to respectively anti-symmetric torsion, symmetric torsion and wing fore-after bending modes, the inaccurate prediction of their dependence in the interval model is a direct cause of the assumption that $\theta_4 - \theta_5$ and $\theta_9 - \theta_{10}$ are fully dependent.

Finally, although the Bayesian posterior distributions predict a non-zero
plausibility for a larger range of parameter values as compared to the interval model, tighter bounds on the prediction of the uncertain parameters as compared to the measurement data set are obtained. This observation stems from the low plausibility that is attributed to the tails of the marginal posterior distributions, combined with the peakedness of the distribution, such that the
moderately small dataset of 500 samples did not sample these tails.

These results indicate that, in case sufficient data are available, Bayesian inference outperforms the multivariate quantification method in terms of information content (e.g., correlation and multi-modality as compared to purely the crisp bounds) and accuracy (less over-conservative while still capturing almost all responses). Note that a part of the discrepancy between the results obtained

by both methods stems from different assumptions on parameter dependence that were made. This however does not invalidate this main conclusion.

5.7. UQ with small data set

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Figure 9 shows all combinations of considered eigenfrequencies, obtained by propagating the quantified intervals and posterior distributions through the AIRMOD FE model.

Concerning the marginal eigenfrequencies, the bounds predicted by the quantified interval method circumscribe the measurement data set tighter as compared to the Bayesian samples, which are shown to be more over-conservative. This is e.g., the case for f_1 , f_2 , f_3 , f_4 , f_5 , f_{11} , f_{14} and f_{19} . The only excep-



Figure 8: All combinations of considered eigenfrequencies, obtained by propagating the quantified intervals and posterior distributions through the AIRMOD FE model. The quantified results were obtained by using the full measurement data set.

tion hereto is the 8th eigenmode (i.e. the 1st symmetric wing torsion mode), which is completely missed by the quantified interval method, as it focussed on encompassing f_7 since both frequencies are equal as can be noted from figure 9. The main reason for this lack of overlap is explained by the assumed perfect dependence between parameters $\theta_4 - \theta_5$ and $\theta_9 - \theta_{10}$, combined with the low data availability. Ideally, the interval method would try to find a set of intervals that encompass both f_7 and f_8 perfectly, but due to the fact that w_0 is set to unity, this did not happen. This is a possible limitation of the method.

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Due to the scarcity of the dataset, none of the methods was able to give an accurate estimate of the dependence between the eigenfrequencies. However, the interval method, apart from the 8th eigenmode, was able to provide tighter bounds on the uncertainty. The ability to provide a better estimate is caused by the effect of the prior distribution for the Bayesian UQ on the identified posterior. To recall, a uniform distribution with very broad basis was selected as to correspond with no prior knowledge on the uncertainty. Since insufficient data are available to compensate for the effect of this prior, the Bayesian method provides a very broad estimate of possible responses.

Therefore, under scarce data, the interval method outperforms the Bayesian method since the lack of informative data makes the quantified posterior highly ⁵¹⁰ dependent on the prior knowledge. The interval method on the other hand does not need an initial estimate of the parameter uncertainty, since the global optimisation routine actively searches the space of input parameters for those intervals that best prescribe the available data.

6. Conclusions

In this paper, an improved version of a novel interval quantification method was presented in the context of high-dimensional models, which has been shown to be able to alleviate the curse of dimensionality that is attributed to the necessary convex hull computations. The latter is illustrated by the factor 10⁵ gain in computational efficiency for the evaluation of the discrepancy between the convex hulls.

The approach has been tested by analysing the DLR-AIRMOD test structure and compared with results obtained by Bayesian model updating. Concerning the interval method with the full dataset, the cross-sections of the resulting convex hull failed to encompass the dataset completely, which is a direct result

of some assumptions that were made in the modelling process. Moreover, propagation of the Bayesian posterior distribution provided a tighter estimation of the experimental dataset. The interval method however provides the analyst



Figure 9: All combinations of considered eigenfrequencies, obtained by propagating the quantified intervals and posterior distributions through the AIRMOD FE model. The quantified results were obtained by using only 5 measured replica.

with fixed bounds between the uncertainty is believed to be located in contrast to the possibly infinite support of the posterior distributions stemming from the Bayesian analysis. In case only limited data are available, the effect of the prior distribution dominates the obtained posterior distribution in the Bayesian case. In this case, propagation of the quantified intervals provides the analyst with a better representation of the measurement dataset in terms of over-conservatism and crispness of the model representation. However, the case study illustrates clearly that it is challenging to perfectly fit a numerically computed convex hull to a convex hull over experimental data. This is a direct result of the fact that in the context of inverse interval uncertainty quantification, a trade-off between conservatism and representing the data accurately exists, where some data points are missed to better represent the full dataset. This was specifically evident when only 5 data samples are considered, as one eigenfrequency was missed by the interval method.

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As such, using this challenging high-dimensional example and the presented theoretical discussion, it is shown that the choice between interval and Bayesian UQ boils down to some key questions an analyst has to consider:

• what is the desired information on the uncertainty?

If the analyst is only interested in bounds on the uncertain parameters, given a set of measurement data, interval analysis has been shown to be very robust with respect to the scarcity of the measurement data. In this case no prior, often subjective, knowledge is needed for the analysis. Should the analyst on the other hand be interested in a complete description of the (joint-)plausibility, including correlation and multi-modal descriptors, a Bayesian approach has to be applied.

• what is the availability of data?

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Since the prior distribution influences the posterior distribution in a Bayesian context under scarce data to a large extent, sufficient informative data should be available to perform a Bayesian analysis. If insufficient data are available, but the analyst requires the information provided by Bayesian analysis; (s)he should collect more data or take utmost care in constructing the prior distribution. Otherwise, the interval approach was proven to give more informative bounds in this case.

Acknowledgements

The authors would like to acknowledge the financial support of the Flemish Research Foundation (FWO) for travel grants K218117N and K217917N and the research project G0C2218N.

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