

Finite element model updating for structural applications

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Abstract

A novel method for performing model updating on finite element models is presented. The approach is particularly tailored to modal analyses of buildings, by which the lowest frequencies, obtained by using sensors and system identification approaches, need to be matched to the numerical ones predicted by the model. This is done by optimizing some unknown material parameters (such as mass density and Young’s modulus) of the materials and/or the boundary conditions, which are often known only approximately. In particular, this is the case when considering historical buildings.

The straightforward application of a general-purpose optimizer can be impractical, given the large size of the model involved. In the paper, we show that, by slightly modifying the projection scheme used to compute the eigenvalues at the lowest end of the spectrum one can obtain local parametric reduced order models that, embedded in a trust-region scheme, form the basis for a reliable and efficient specialized algorithm.

We describe an optimization strategy based on this approach, and we provide numerical experiments that confirm its effectiveness and accuracy.

Keywords: Model updating, Finite elements, Trust-region, Lanczos, Eigenvalue optimization

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1. Introduction

Finite element (FE) model updating is a procedure aimed at calibrating the FE model of a structure in order to match numerical and experimental results. Introduced in the 1980s, it turned out to play a crucial role in the design, analysis and maintenance of aerospace, mechanical and civil engineering structures [14, 19, 24, 31]. In structural mechanics, model updating techniques are used in conjunction with vibrations measurements to determine unknown system characteristics, such

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as the materials' properties, constraints, etc. The resulting updated FE model can then be used to obtain reliable predictions on the dynamic behavior of the structure subjected to time-dependent loads. A further important application of model updating, within the framework of structural health monitoring, is damage detection [41, 45]. Within this framework, damage can be identified based on the assumption that its presence is associated with a decrease in the stiffness of some elements, with consequent changes in the structure's modal characteristics.

Finite element model updating involves the solution of a constrained optimization problem, whose objective function is generally expressed as the discrepancy between experimental and numerical quantities, such as the structure's natural frequencies and mode shapes. Ill-posedness or ill-conditioning can affect model updating formulations and lead to numerical problems due to inaccuracy in the model and lack of information in the measurements. Among the methods aimed at quantifying uncertainties, the probabilistic Bayesian approach is one of the most adopted [41].

Application of FE model updating to ancient masonry buildings is relatively recent. In [4, 5, 12, 13, 15, 21, 23, 25, 29, 35, 39, 40, 46] a vibration-based model updating is conducted, and preliminary FE models are tuned by using the dynamic characteristics determined through system identification techniques. In the papers cited above the modal analysis of the FE models are conducted via commercial codes, and the model updating procedure is implemented separately. In this paper, on the contrary, FE model updating is integrated within a software package, the NOSA-ITACA code [11], able to manage the large-scale problems encountered in applications. In particular, the algorithms for the solution of the constrained optimization problem integrated in NOSA-ITACA exploit the structure of the stiffness and mass matrices and the fact that only a few of the smallest eigenvalues have to be calculated. This new procedure reduces both the total computation time of the numerical process and user's effort, thus providing the scientific and technical communities with efficient algorithms specific for FE model updating.

A simple form of FE model updating has been employed in combination with the NOSA-ITACA code in [6, 7, 34], to perform modal analyses of the San Frediano bell tower and the Clock tower in Lucca, Italy. In these works the optimal values of Young's modulus and the materials' mass density are determined by fitting the data measured by seismometric stations placed on the towers and running several simulation on a grid of feasible values. However, this approach becomes impractical if the number of free parameters or the size of the model is considerably increased.

Model reduction techniques have long been used to reduce the size of complex FE models to a more manageable order. This allows, for example, performing more demanding, and computationally costly, numerical tasks, such as optimization, simulation, and so on, in an efficient manner. In particular, reduced models are natural candidates for optimization of the model's free parameters, either to improve some engineering properties or to better match the empirically measured modes. In practice, this step often turns into an eigenvalue optimization, where part of the spectral structure is tuned following some prescribed criteria.

However, when the model depends on parameters, it is generally non-trivial to obtain a reduced parametric model that accurately reflects the behavior of the original one for all possible parameter values. This is precisely the aim of so-called parametric model reduction, which has recently been used in combination with optimization algorithms [3, 32, 38, 48], especially involving trust-region solvers [1, 47]. Very often, these strategies involve some form of low-rank approximation [42]; in this work we propose an efficient model reduction strategy derived by properly recycling the Lanczos projection. Our approach is tailored to the needs of structural FE analysis.

In fact, FE study of buildings has some peculiar features that justify trying to devise an adapted method. As discussed in the following, in this context the natural frequencies of interest are those

at the lowest end of the real spectrum. In order to compute them accurately, the natural choice is an (inverse) Lanczos method. When a parametric model is given, the Lanczos projection can be re-interpreted as a parameter dependent model reduction, whereby only the relevant part of the spectrum is matched. We show that this step can be performed efficiently, and that its combination with a trust-region method allows matching the measured frequencies with the ones predicted by the parametric model.

The paper is organized as follows. In Section 2 we formulate the optimization problem related to model updating. In Section 3 we describe the requirements for a reduced model to be used in a trust-region optimization method, in particular those required to guarantee convergence. Section 4 describes how to construct a model that satisfies these requirements by slightly modifying the Lanczos projection step used to compute the smallest frequencies, and finally Section 5 illustrates some tests of our implementation on some problems. In Section 6 we draw some conclusions and discuss future lines of research.

2. Modal analysis of masonry buildings and frequency matching

Although the constituent masonry materials of historical buildings exhibit different strengths under tension and compression and, therefore, behave nonlinearly, modal analysis, which is based on the assumption that the materials are linear elastic, is widely used in applications and provides important qualitative information on the dynamic behavior of masonry structures. Modal analysis consists in the solution of the constrained generalized eigenvalue problem

$$K v = \omega^2 M v, \quad \text{subject to } C v = 0. \quad (1)$$

The left part of (1) is derived from the differential equation

$$M \ddot{u} + K u = 0, \quad (2)$$

governing the undamped free vibrations of a linear elastic structure discretized into finite elements. In (2) u is the displacement vector, which belongs to $\mathbb{R}^n$ and depends on time t , $\ddot{u}$ is the second-derivative of u with respect to t , and K and $M \in \mathbb{R}^{n \times n}$ are the stiffness and mass matrices of the FE model. K is symmetric and positive-semidefinite, M is symmetric and positive-definite, and both are sparse and banded. Displacements u_i are also called degrees of freedom; the integer n is the total number of degrees of freedom of the system and is generally very large, since it depends on the level of discretization of the problem. By assuming that $u = v \cdot \sin(\omega t)$, with v a vector of $\mathbb{R}^n$ and ω a real scalar representing the circular frequency, and applying the mode superposition procedure [9], equation (2) is transformed into the generalized eigenvalue problem (1). The right part of equation (1) expresses the fixed constraints and the master-slave relations assigned to displacement u , written in terms of vector v , $C \in \mathbb{R}^{h \times n}$ and h is the number of boundary conditions, with $h \ll n$. Imposing the constraints and boundary conditions is equivalent to projecting the matrices K and M on a subspace where they are a symmetric positive definite pencil (the right kernel of the operator C). In the following, we assume that this projection has already been done — and we refer the reader to [37] for further details. Notice, in particular, that if K and M depend *linearly* on some parameters $\mathbf{x} := (x_1, \dots, x_\ell)$, the same holds true for their projection. More generally, the smooth dependency of K and M is preserved by the projection, since it is a linear operation. This will be exploited in the analysis of the parametric projection in Section 4.

The eigenvalues ω_i^2 of (1) are linked to the natural frequencies, or eigenfrequencies f_i of the structure via the relation $f_i = \omega_i/2\pi$, and the eigenvectors $v^{(i)}$ are the corresponding mode shape

vectors, or eigenmodes. Together with the natural frequencies, the mode shapes provide a good deal of qualitative information on the structure's deformations under dynamic loads.

Measuring the vibrations of masonry buildings is a common practice for assessing their dynamic behavior. Historic constructions are subjected to a number of sources of vibrations, such as traffic, micro-tremors, wind and earthquakes. The availability of sensitive instruments to detect buildings' movements makes it possible to conduct accurate, long-term monitoring campaigns. In fact, ambient vibration monitoring can provide important information on the structural health of old masonry constructions, as it is a non-destructive technique able to capture the most important features of their dynamic behavior, such as natural frequencies, damping ratios, mode shapes and wave propagation velocities. Once the influence of environmental factors has been accounted for, changes in these dynamic properties over time may represent effective structural damage indicators [7].

FE model updating is a procedure that enables calibrating a finite element model of a structure in order to match numerical and experimental results. Thus, model updating techniques are used in conjunction with vibrations measurements to determine a structure's characteristic, such as materials' properties, constraints, etc., which are generally unknown.

2.1. Formulation of the optimization problem

The model updating problem can be reformulated as an optimization problem by assuming that the (projected) stiffness and mass matrices K and M are functions of the parameter vector $\mathbf{x}$. We use the notation

$$K := K(\mathbf{x}), \quad M := M(\mathbf{x}), \quad \mathbf{x} \in \mathbb{R}^\ell,$$

to denote this dependency. The set of valid choices for the parameters is denoted by Ω , which we assume to be an ℓ -dimensional box, that is

$$\Omega = [a_1, b_1] \times \dots \times [a_\ell, b_\ell], \quad (3)$$

for certain values $a_i < b_i$, $i = 1, \dots, \ell$. We also assume that initial estimates for the parameter values are available, and we denote these by $\mathbf{x}^{(0)}$. When this is not the case, we choose $\mathbf{x}^{(0)}$ as the center of the box Ω (i.e., the vector with the midpoints of the intervals $[a_i, b_i]$ as coordinates). The intervals $[a_i, b_i]$ will be chosen accordingly to the physical quantities that they describe.

Our ultimate aim is to determine the optimal value of $\mathbf{x}$ that minimizes a certain cost functional $\phi(\mathbf{x})$ within the box Ω . This is an instance of the optimization problem

$$\min_{\mathbf{x} \in \Omega} \phi(\mathbf{x}). \quad (4)$$

The choice of the objective function $\phi(\mathbf{x})$ is related to the frequencies that we want to match. If we need to match s frequencies of the model, we choose a suitable weight vector $\mathbf{w} := [w_1, \dots, w_s]$, with $w_i \geq 0$, and define the functional $\phi(\mathbf{x})$ as follows:

$$\phi(\mathbf{x}) := \left\| \frac{\sqrt{\Lambda_s(K, M)}}{2\pi} - \mathbf{f} \right\|_{\mathbf{w}, 2}^2, \quad \|\mathbf{y}\|_{\mathbf{w}, 2} := \sqrt{\mathbf{y}^T D_{\mathbf{w}}^2 \mathbf{y}}, \quad (5)$$

where $\mathbf{f}$ is the vector of the measured frequencies, $D_{\mathbf{w}} = \text{diag}(w_1, \dots, w_s)$ and $\Lambda_s(K, M)$ is the vector containing the smallest s eigenvalues of the pencil $K - \lambda M$, ordered according to their magnitude.

The vector $\mathbf{w}$ encodes the weights that should be given to each frequency in the optimization scheme. If the aim is to minimize the distance between the vector of measured and computed frequencies in the usual Euclidean norm, then $\mathbf{w} = \mathbb{1}_s$, the vector of all ones, should be chosen. If, instead, a relative accuracy on the frequencies is desired, $w_i = f_i^{-1}$ is the natural choice. To avoid scaling issues in the objective function, we always normalize $\mathbf{w}$ to have Euclidean norm 1.

3. The optimization framework

In this section we describe the optimization framework that we use. We choose to rely on trust-region methods, well-known iterative methods for solving bound-constrained optimization problems, see e.g. [16, 17, 36]. Their main distinctive features are robustness and convergence to first-order critical points regardless of the choice of the initial estimate.

A trust-region method is defined by building a sequence of local models for the objective function. Construction of the local models is discussed in Section 4, where we show that such models satisfy the minimal requirements for a well-defined trust-region scheme. The practical choice of the objective function and the weights has been deferred to Section 5, where several options are discussed.

3.1. Trust region methods

We now review the basic steps in a trust-region algorithm for solving the bound-constrained optimization problem of Equation (4), where $\phi : \mathbb{R}^\ell \rightarrow \mathbb{R}$ is a smooth function and Ω is defined as in (3).

The key idea of a trust-region method is to define, at iteration k , a model $\phi_k^{\mathcal{R}}$ for the objective function ϕ around the current iterate $\mathbf{x}^{(k)}$, together with a region within such model can be trusted to provide an adequate representation of ϕ . The trial step $\mathbf{s}^{(k)}$ is then computed, either exactly or approximately, by minimizing this model within the trust-region. The trust-region is the set of all points

$$\mathcal{B}_k = \{\mathbf{x} \in \mathbb{R}^\ell \mid \|\mathbf{x} - \mathbf{x}^{(k)}\| \leq \Delta_k\},$$

where Δ_k is called the trust-region radius and $\|\cdot\|$ is a norm equivalent to the Euclidean norm.

Due to the presence of bound-constraints, all the generated iterates $\mathbf{x}^{(k)}$ must be ensured to be feasible, that is, $\mathbf{x}^{(k)} \in \Omega$ for all $k \geq 0$. One strategy consists in minimizing the model within the set $\mathcal{B}_k \cap \Omega$. When the feasible set Ω is a box, the trust-region is generally defined using the ∞ norm.

Having minimized the model on $\mathcal{B}_k \cap \Omega$, it must be decided whether to accept the trial step or to change the trust-region radius. Usually, the trust-region radius and the new point $\mathbf{x}^{(k)} + \mathbf{s}^{(k)}$ are tested simultaneously to assess the quality of the approximation yielded by the local model. This is measured using the ratio $\rho^{(k)}$ between the actual and the predicted reduction defined as follows:

$$\rho^{(k)} = \frac{\phi(\mathbf{x}^{(k)}) - \phi(\mathbf{x}^{(k)} + \mathbf{s}^{(k)})}{\phi_k^{\mathcal{R}}(\mathbf{x}^{(k)}) - \phi_k^{\mathcal{R}}(\mathbf{x}^{(k)} + \mathbf{s}^{(k)})} \quad (6)$$

If $\rho^{(k)}$ is close to 1, there is good agreement between the model $\phi^{\mathcal{R}}$ and the function ϕ over this step, so $\mathbf{x}^{(k)} + \mathbf{s}^{(k)}$ is accepted as the new iterate and it is safe to increase the radius of the trust-region for the next iteration. If $\rho^{(k)}$ is positive but not close to 1, then $\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \mathbf{s}^{(k)}$ but the trust-region radius is not altered. If $\rho^{(k)}$ is close to zero or negative, step $\mathbf{s}^{(k)}$ is rejected and the trust-region radius is shrunk.

Convergence of the iterative process is declared when a suitable criticality measure is sufficiently small. We consider the measure

$$\chi(\mathbf{x}^{(k)}) = \|P_\Omega(\mathbf{x}^{(k)} - \nabla\phi(\mathbf{x}^{(k)})) - \mathbf{x}^{(k)}\| \quad (7)$$

where P_Ω is the projection onto the feasible set, and $\chi(\mathbf{x}^{(k)})$ is the norm of the projected gradient onto the box, and reduces to $\|\nabla\phi(\mathbf{x}^{(k)})\|$ when $\mathbf{x}^{(k)}$ is in the interior of Ω . The approach is summarized in the pseudocode of Algorithm 1 where we assume that the model functions are minimized “exactly” (as it will be the case in our applications, see Section 5.1).

Algorithm 1 Basic trust-region algorithm

Require: The initial point $\mathbf{x}^{(0)} \in \Omega$, an initial trust-region radius Δ_0 , the constants $\eta_1, \eta_2, \gamma_1, \gamma_2$ such that

$$0 < \eta_1 \leq \eta_2 < 1 \text{ and } 0 < \gamma_1 \leq \gamma_2 < 1,$$

$k_{max} > 0$ and $\epsilon > 0$.

Compute $\phi(\mathbf{x}^{(0)})$

for $k = 1, \dots, k_{max}$ **do**

 Define a model $\phi_k^{\mathcal{R}}$ in $\mathcal{B}_k \cap \Omega$.

 Compute a step $\mathbf{s}^{(k)}$ that minimizes the model and such that $\mathbf{x}^{(k)} + \mathbf{s}^{(k)} \in \mathcal{B}_k \cap \Omega$.

 Compute $\phi(\mathbf{x}^{(k)} + \mathbf{s}^{(k)})$ and define the ratio (6).

if $\rho^{(k)} \geq 1$ **then**

 define $\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \mathbf{s}^{(k)}$;

else

 define $\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)}$.

end if

 Update the trust-region radius setting

$$\Delta_{k+1} \in \begin{cases} [\Delta_k, \infty) & \text{if } \rho^{(k)} \geq \eta_2 \\ [\gamma_2 \Delta_k, \Delta_k] & \text{if } \rho^{(k)} \in [\eta_1, \eta_2] \\ [\gamma_1 \Delta_k, \gamma_2 \Delta_k] & \text{if } \rho^{(k)} < \eta_1 \end{cases}$$

 Compute the criticality measure (7).

if $\chi(\mathbf{x}^{(k+1)}) \leq \epsilon$ **then**

return An approximate first-order critical point $\mathbf{x}^{(k+1)}$.

end if

end for

In order to obtain a robust and globally convergent trust-region algorithm involving approximate models, the following assumptions should hold [17] for the objective ϕ and the model $\phi_k^{\mathcal{R}}$ functions:

(AF.1) $\phi : \mathbb{R}^\ell \rightarrow \mathbb{R}$ is twice-continuously differentiable on Ω .

(AF.2) The function ϕ is bounded below for all $\mathbf{x} \in \Omega$.

(AF.3) The second derivatives of ϕ are uniformly bounded for all $\mathbf{x} \in \Omega$.

(AM.1) For all k , $\phi_k^{\mathcal{R}}$ is twice differentiable on $\mathcal{B}_k$.

(AM.2) The values of the objective and the model function coincide at the current iterate, i.e., for all k ,

$$\phi(\mathbf{x}^{(k)}) = \phi_k^{\mathcal{R}}(\mathbf{x}^{(k)}).$$

(AM.3) The gradients of the objective and the model function coincide at the current iterate, i.e., for all k ,

$$\nabla \phi(\mathbf{x}^{(k)}) = \nabla \phi_k^{\mathcal{R}}(\mathbf{x}^{(k)}).$$

(AM.4) The second derivatives of $\phi^{\mathcal{R}}(\mathbf{x}^{(k)})$ remain bounded within the trust-region $\mathcal{B}_k$ for all k .

Theorem 3.1. [17, Section 12.2.2] *Under Assumptions (AF.1)-(AF.3) and (AM.1)-(AM.4), every limit point of the sequence $\{\mathbf{x}^{(k)}\}$ generated by Algorithm 1 is first-order critical, i.e.*

$$\lim_{k \rightarrow \infty} \|P_{\Omega}(\mathbf{x}^{(k)} - \nabla \phi(\mathbf{x}^{(k)})) - \mathbf{x}^{(k)}\| = 0.$$

where P_{Ω} is the projection on the feasible set.

3.2. Enforcing first-order matching

Here we use the trust-region framework described in the previous section to solve problem (4) with objective function (5). Note that conditions (AF.1) – (AF.3) are automatically satisfied by our definition of objective function. As discussed in Section 4.1, the local model $\tilde{\phi}_k^{\mathcal{R}}(\mathbf{x})$ that we build does not necessarily agree at the first order with the original function $\phi(\mathbf{x})$, which apparently breaks assumption (AM.3) — and only satisfies (AM.2). However, if the gradient $\nabla \phi_k^{\mathcal{R}}(\mathbf{x})$ is known at the expansion point $\mathbf{x}^{(k)}$, this can be easily fixed by defining

$$\phi_k^{\mathcal{R}}(\mathbf{x}) := \tilde{\phi}_k^{\mathcal{R}}(\mathbf{x}) + (\phi(\mathbf{x}^{(k)}) - \tilde{\phi}_k^{\mathcal{R}}(\mathbf{x}^{(k)})) + (\nabla \phi(\mathbf{x}^{(k)}) - \nabla \tilde{\phi}_k^{\mathcal{R}}(\mathbf{x}^{(k)}))^T (\mathbf{x} - \mathbf{x}^{(k)}), \quad (8)$$

see, e.g., [1, 2, 26].

4. Building the reduced model by Lanczos projection

Our aim in this section is to develop a low-dimensional model of the frequencies as a function of the parameters. The model must be efficient to evaluate, and needs to satisfy the requirements for use as a local model in the trust-region approach described in Section 3.

This task can be viewed as a special case of parametric model order reduction. Several approaches have been developed in this field that allow obtaining low-dimensional models for accurately approximating the behavior of complex systems. For instance, when controlling large-scale dynamical systems, it is possible to employ projection approaches when the observations and inputs have low-rank properties (see [10] and references therein).

Our aim is slightly different; we are interested only in a particular region on the spectrum (the lowest frequencies), but we have no particular assumptions on the input (which is typically induced by the environment).

In the next sections, we first briefly summarize the standard (inverse) Lanczos iteration for matrix pencils, mainly to fix the notation, and then we show that we can perform some modifications to make it parameter dependent. We refer to [8] for a discussion of the relation between the Lanczos process and (nonparametric) model reduction.

4.1. The Lanczos method

The Lanczos iteration, for a symmetric $n \times n$ matrix A , can be summarized as follows. We construct an orthogonal basis W_m for the subspace $\mathcal{K}_m(A, v) \subseteq \mathbb{R}^n$ defined as:

$$\mathcal{K}_m(A, v) := \text{span}(v, Av, \dots, A^{m-1}v).$$

Unless a breakdown occurs, this space has dimension m , and the basis W_m is represented as a $n \times m$ orthogonal matrix (that is, $W_m^T W_m = I_m$). Given the basis, one then computes the projection $T_m = W_m^T A W_m$. The matrix T_m is tridiagonal, and its eigenvalues (known as Ritz values) are good approximations of the extremal eigenvalues of A even when $m \ll n$. This procedure is appealing for several reasons.

- (i) Computation of the space W_m only requires matrix-vector multiplications, and therefore it is easy to exploit the sparsity or, more generally, the structure of A .
- (ii) The orthogonalization of the basis requires only scalar products, which can be computed very efficiently if m does not increase too much.
- (iii) The algorithm can be carried out iteratively, i.e., the optimal value of m does not need to be known a priori, and T_m can be built incrementally. In our framework, we stop the iteration when the error on the computed eigenvalues is guaranteed to be smaller than a certain relative threshold τ (see [18] for a classical reference on the stopping criterion of Lanczos for computing eigenvalues). The actual value of τ is reported in Section 5.

When the matrix A arises from a FE discretization of a partial differential equation (PDE) involving only local operators and elements with compact support and few overlappings, the number of nonzero elements is only $\mathcal{O}(n)$; hence, the matrix vector multiplication can be carried out with linear complexity. Moreover, the matrix usual has a band or multiband structure, and using good sparse solvers or appropriate preconditioners we can solve linear systems in $\mathcal{O}(n)$ time. Indeed, this is easy to establish for 1D problems, where the stiffness matrix is usually banded and a system solution can be computed by a banded LU and backsubstitution in linear time. For 2D and 3D problems, linear-time complexity approaches have been developed by relying, for instance, on optimal multigrid solver, see [20, 27]. Therefore, the smallest eigenvalue of A can be extracted applying Lanczos to A^{-1} . Since we are interested in modal analysis of structures discretized by FE, we need to approximate the smallest eigenvalues of a pencil $K - \lambda M$, with K and M symmetric positive definite; the same technique to extract the smallest eigenvalues can be applied (implicitly) to the symmetric matrix $M^{\frac{1}{2}} K^{-1} M^{\frac{1}{2}}$, obtaining:

$$T_m = W_m^T M^{\frac{1}{2}} K^{-1} M^{\frac{1}{2}} W_m = U_m^T M K^{-1} M U_m, \quad U_m = M^{-\frac{1}{2}} W_m.$$

It is immediately verifiable that U_m contains a basis of the same subspace spanned by W_m , but this matrix is M -orthogonal (that is, orthogonal with respect to the inner product defined by M), so $U_m^T M U_m = I_m$ [22].

Remark 4.1. In the above framework, there is no need to explicitly compute the square root of the positive definite matrix M , which would be a very expensive operation. The matrix T_m is implicitly obtained using the matrix-vector product with $M K^{-1} M$ and re-orthogonalization by the scalar-product induced by M , computing the basis U_m directly.

However, using this procedure as a black-box to evaluate the objective function in an optimization method can easily become too onerous: even if the complexity is only linear, the size of the problem can make this operation unpractical (we often deal with $n \approx 10^6$ or even more in finite element models).

Therefore, we wish to investigate the use of the Lanczos process to build a local parametric reduced model for the case when K and M are not constant matrices, but depend on ℓ parameters $x_1, \dots, x_\ell$.

4.2. Parametric finite element models

Let us assume that a finite element model depending on a variable $\mathbf{x} \in \mathbb{R}^\ell$ is given. These parameters $\mathbf{x}$ may encode different properties of the system (our method does not make any assumption about what they mean — but only on their algebraic properties). Most of the time, these will be some materials' physical characteristics, such as Young's modulus, or mass density. In almost all the cases of interest, the dependency is *linear* in $\mathbf{x}$. However, for our discussion we just assume it to be smooth, i.e., at least $\mathcal{C}^2$.

4.3. A first-order approximation to the updated Lanczos projection

In this section we develop a strategy that, given a certain initial value $\mathbf{x}^{(0)}$ for the model parameters, builds a local model to be used in the trust-region scheme presented in Section 3.

Let us denote with $F(\mathbf{x})$ the matrix valued function that, for a given choice of the parameters $\mathbf{x}$, returns the tridiagonal matrix obtained running the Lanczos process on $M^{\frac{1}{2}}(\mathbf{x})K(\mathbf{x})^{-1}M^{\frac{1}{2}}(\mathbf{x})$.

We shall approximate $F(\mathbf{x})$ in a neighborhood of $\mathbf{x}^{(0)}$. This approximation calls for two steps: first, we fix the subspace used for the projection, as described in Lemma 4.3, obtaining an approximation $F_0(\mathbf{x})$ of $F(\mathbf{x})$. Then, we show how to further approximate this function in order to make it cheap to evaluate. Lemma 4.7 describes how to combine these approximations to obtain a first-order correct local model for the trust-region scheme.

Remark 4.2. The outcome of the Lanczos process depends on the initial vector chosen for the scheme. Here and in the following, we assume that this vector is fixed whenever we keep $\mathbf{x}^{(0)}$ unchanged. This makes the subspace generated by the Lanczos projection unique.

Lemma 4.3. *Let $F(\mathbf{x}) = U_m(\mathbf{x})M(\mathbf{x})K^{-1}(\mathbf{x})M(\mathbf{x})U_m(\mathbf{x})$ be the matrix-valued function that associates the point $\mathbf{x}$ with the Lanczos projection of the pencil $K(\mathbf{x}) - \lambda M(\mathbf{x})$ previously defined. Then, there exists a matrix-valued function $U_{m,0}(\mathbf{x})$ such that, in a neighborhood $\mathcal{N}$ of $\mathbf{x}^{(0)}$,*

1. *For any $\mathbf{x} \in \mathcal{N}$ the matrix-valued function $U_{m,0}(\mathbf{x})$ is an $M(\mathbf{x})$ -orthogonal basis for the column-span of $U_m(\mathbf{x}^{(0)})$.*
2. *The function $F_0(\mathbf{x})$ defined as*

$$F_0(\mathbf{x}) := U_{m,0}(\mathbf{x})^T M(\mathbf{x}) K(\mathbf{x})^{-1} M(\mathbf{x}) U_{m,0}(\mathbf{x})$$

is $\mathcal{C}^2(\mathcal{N})$ and $F_0(\mathbf{x}^{(0)}) = F(\mathbf{x}^{(0)})$, that is it is a zeroth-order approximation of $F(\mathbf{x})$.

Proof. By construction, the basis generated by the Lanczos method at $\mathbf{x}^{(0)}$ returns an $M(\mathbf{x}^{(0)})$ -orthogonal basis $U_m(\mathbf{x}^{(0)})$. We have

$$U_m(\mathbf{x}^{(0)})^T M(\mathbf{x}) U_m(\mathbf{x}^{(0)}) = Z(\mathbf{x}), \quad Z(\mathbf{x}^{(0)}) = I,$$

and since $M(\mathbf{x})$ is of class $\mathcal{C}^2$, the same can be said of the symmetric matrix $Z(\mathbf{x})$. Therefore, we can define $U_{m,0}(\mathbf{x}) := U_m(\mathbf{x}^{(0)})Z(\mathbf{x})^{-\frac{1}{2}}$. For $\mathbf{x}$ close enough to $\mathbf{x}^{(0)}$, the value of $Z(\mathbf{x})$ is bounded away from singular matrices (recall that $Z(\mathbf{x}^{(0)}) = I$), so the inverse square-root is locally analytic and therefore $U_{m,0}(\mathbf{x})$ is at least $\mathcal{C}^2(\mathcal{N})$ in a neighborhood $\mathcal{N}$. Direct substitution yields $U_{m,0}^T(\mathbf{x})M(\mathbf{x})U_{m,0}(\mathbf{x}) = I$ and $F_0(\mathbf{x}^{(0)}) = F(\mathbf{x}^{(0)})$, as requested. $\square$

The idea behind the approximation of Lemma 4.3 is to obtain the eigenvalues of $F_0(\mathbf{x})$ by a subspace projection method, where the subspace is approximated by choosing the one obtained by running Lanczos for the parameters at $\mathbf{x}^{(0)}$ instead of the one at $\mathbf{x}$. If the two values are close, this subspace is still “good enough” to provide accurate approximations of the spectrum. In fact a similar strategy underlies several methods known as *Krylov subspace recycling* [33, 43]. These techniques find applications in solving sequences of shifted linear systems, which is a step required, for instance, in model reduction algorithms.

The definition of $F_0(\mathbf{x})$, as is, does not make it particularly easier to evaluate with respect to $F(\mathbf{x})$. In Section 4.3.1, we will show how to approximate $F_0(\mathbf{x})$ at the first-order in a way that makes its computation very efficient. The next result justifies this approach.

Lemma 4.4. *Let $\hat{F}_0(\mathbf{x})$ be any locally $\mathcal{C}^2$ first-order symmetric positive definite approximation of $F_0(\mathbf{x})$, and let $\hat{\Lambda}_s(\cdot)$ be the function that associates a symmetric matrix with the inverse of its largest s eigenvalues. If the eigenvalues at $\mathbf{x}^{(0)}$ of $F(\mathbf{x}^{(0)})$ have multiplicities that do not change in an open neighborhood of $\mathbf{x}^{(0)}$, then*

$$\hat{\Lambda}_s(F(\mathbf{x})) = \hat{\Lambda}_s(\hat{F}_0(\mathbf{x})) + O(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2), \quad \mathbf{x} \in \mathcal{N},$$

where $\mathcal{N}$ is an appropriate neighborhood of $\mathbf{x}^{(0)}$.

Proof. When the eigenvalues of a symmetric matrix are all distinct, they locally depend analytically on the entries of the matrix (see, for instance, [28, 44]). This statement is valid also in the more general case where the eigenvalues are not simple, but their multiplicity is locally constant (i.e., there exists an open neighborhood where the eigenvalues do not change the multiplicity). Since the matrices involved are all positive definite, the same holds for their inverses λ_i^{-1} . Therefore, the result follows by composing $\hat{\Lambda}_s(\cdot)$ with the functions $F(\mathbf{x})$ and $\hat{F}_0(\mathbf{x})$, that is considering $\hat{\Lambda}_s(F(\mathbf{x}))$ and $\hat{\Lambda}_s(\hat{F}_0(\mathbf{x}))$, since the composition of analytic functions is again analytic. $\square$

Remark 4.5. In the event two or more eigenvalues coincide the situation is slightly more problematic, since the dependency of the eigenvalues is only continuous in general, and although partial derivatives exist, a higher-order smoothness in the eigenvalue functions cannot be claimed [44]. We note that the objective function is still well-defined in the case of multiple eigenvalues; the only limitation is a possible loss of regularity at the eigenvalues crossing. This could slow down the convergence if the minimum is at that point. The example reported in Section 5 shows that the case of multiple eigenvalues can be managed in a satisfactory way by the proposed algorithm, even when the hypothesis of Lemma 4.4 is not satisfied.

4.3.1. First-order expansion of the projection

In view of Lemma 4.4, in this Section we aim at constructing an approximation to $F_0(\mathbf{x})$ that matches at the first-order, and that is numerically undemanding to evaluate. More precisely, we

aim at complexity $\mathcal{O}(m^3)$, where m is the dimension of the Krylov subspace. Let us write the value of $K(\mathbf{x})$ and $M(\mathbf{x})$ as small perturbations of their values at $\mathbf{x}^{(0)}$:

$$K(\mathbf{x}) = K(\mathbf{x}^{(0)}) + \delta K(\mathbf{x}), \quad M(\mathbf{x}) = M(\mathbf{x}^{(0)}) + \delta M(\mathbf{x}).$$

We can expand the above expressions to the first-order, which yields

$$K(\mathbf{x}) = K(\mathbf{x}^{(0)}) + \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) \frac{\partial K(\mathbf{x}^{(0)})}{\partial x_j} + O(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2), \quad (9)$$

and the analogous formula for $M(\mathbf{x})$. Let us assume that we are given an $M(\mathbf{x})$ -orthogonal basis $U_{m,0}(\mathbf{x})$ spanning the same subspace of $U_m(\mathbf{x}^{(0)})$ that we have computed with the Krylov projection method at $\mathbf{x}^{(0)}$. In this case, we can compute the new projected counterpart of $M^{\frac{1}{2}}(\mathbf{x})K(\mathbf{x})^{-1}M^{\frac{1}{2}}(\mathbf{x})$ as follows:

$$F_0(\mathbf{x}) = U_{m,0}^T(\mathbf{x})M(\mathbf{x})K(\mathbf{x})^{-1}M(\mathbf{x})U_{m,0}(\mathbf{x}). \quad (10)$$

In order to derive a first-order expansion for the above formula we need first-order expansions for all the terms involved. However, we still do not have such an expression for the inverse of $K(\mathbf{x})$. For $\mathbf{x}$ sufficiently close to $\mathbf{x}^{(0)}$ we can write the Neumann expansion:

$$K(\mathbf{x})^{-1} = (I + K(\mathbf{x}^{(0)})^{-1}\delta K(\mathbf{x}))^{-1}K(\mathbf{x}^{(0)})^{-1} = \sum_{j=0}^{\infty} (-1)^j \left[K(\mathbf{x}^{(0)})^{-1}\delta K(\mathbf{x}) \right]^j K(\mathbf{x}^{(0)})^{-1}.$$

Relying on Equation (9), and dropping the second-order terms yields

$$K(\mathbf{x})^{-1} = K(\mathbf{x}^{(0)})^{-1} - \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) K(\mathbf{x}^{(0)})^{-1} \frac{\partial K(\mathbf{x}^{(0)})}{\partial x_j} K(\mathbf{x}^{(0)})^{-1} + O(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2),$$

which in turn can be rewritten in the following form:

$$K(\mathbf{x})^{-1} = K(\mathbf{x}^{(0)})^{-1} \left(I - \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) \frac{\partial K(\mathbf{x}^{(0)})}{\partial x_j} K(\mathbf{x}^{(0)})^{-1} \right) + O(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2). \quad (11)$$

Equation (11) has the desirable property that all the inverses of $K(\mathbf{x})$ appear evaluated at $\mathbf{x}^{(0)}$, and the dependency on $\mathbf{x}$ is linear.

We now have all the ingredients to derive a cheap formula for computing $F_0(\mathbf{x})$, by substituting the expansions in Equation (10). This leads to:

$$\begin{aligned} F_0(\mathbf{x}) &= U_{m,0}^T(\mathbf{x})M(\mathbf{x}^{(0)})K(\mathbf{x}^{(0)})^{-1}M(\mathbf{x}^{(0)})U_{m,0}(\mathbf{x}) \\ &\quad + \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) U_{m,0}(\mathbf{x})^T G_j(\mathbf{x}^{(0)}) U_{m,0}(\mathbf{x}) + O(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2), \end{aligned} \quad (12)$$

where $G_j(\mathbf{x}^{(0)})$ is obtained by grouping together all the first-order contributions, that is,

$$\begin{aligned} G_j(\mathbf{x}^{(0)}) &= \frac{\partial M(\mathbf{x}^{(0)})}{\partial x_j} K(\mathbf{x}^{(0)})^{-1} M(\mathbf{x}^{(0)}) - M(\mathbf{x}^{(0)}) K(\mathbf{x}^{(0)})^{-1} \frac{\partial K(\mathbf{x}^{(0)})}{\partial x_j} K(\mathbf{x}^{(0)})^{-1} M(\mathbf{x}^{(0)}) \\ &\quad + M(\mathbf{x}^{(0)}) K(\mathbf{x}^{(0)})^{-1} \frac{\partial M(\mathbf{x}^{(0)})}{\partial x_j}. \end{aligned}$$

The terms $G_j(\mathbf{x}^{(0)})$ are large matrices, and computing them explicitly would be unfeasible. However, by observing Equation (12) closely, we immediately notice that we do not need the complete matrices, but just their projected counterparts $\hat{G}_j(\mathbf{x}) := U_{m,0}(\mathbf{x})^T G_j(\mathbf{x}^{(0)}) U_{m,0}(\mathbf{x})$, which are much smaller. Moreover, in the definition of $G_j(\mathbf{x}^{(0)})$ the only matrix that appears as an inverse is $K(\mathbf{x}^{(0)})$. In our Lanczos implementation we use a direct sparse solver so, after the projection, we already have a sparse Cholesky factorization of $K(\mathbf{x}^{(0)})$, and we compute $\hat{G}_j(\mathbf{x}^{(0)})$ at the cost of some extra back-substitutions and sparse matrix-vector products (which have a comparable cost). Similar savings could be obtained recycling the preconditioner when using an iterative method.

However, the same computational cost would be needed to calculate $\hat{G}_j$ at $\mathbf{x} \neq \mathbf{x}^{(0)}$. We would like to avoid this extra computation. To this end, observe that according to the proof of Lemma 4.3, we have $U_{m,0}(\mathbf{x}) = U_m(\mathbf{x}^{(0)}) Z(\mathbf{x})^{-\frac{1}{2}}$. Therefore, we can write the formulas to compute $\hat{G}_j(\mathbf{x})$ as follows:

$$\hat{G}_j(\mathbf{x}) = Z(\mathbf{x})^{-\frac{1}{2}} U_m(\mathbf{x}^{(0)})^T G_j(\mathbf{x}^{(0)}) U_m(\mathbf{x}^{(0)}) Z(\mathbf{x})^{-\frac{1}{2}},$$

which shifts the problem to that of computing $Z(\mathbf{x}) = U_m(\mathbf{x}^{(0)})^T M(\mathbf{x}) U_m(\mathbf{x}^{(0)})$. As in the previous equations, we can obtain a first-order approximation of $Z(\mathbf{x})$ at any $\mathbf{x}$ with a first-order truncation by precomputing some small projected matrices at $\mathbf{x}^{(0)}$:

$$\begin{aligned} \hat{Z}(\mathbf{x}) &:= U_m(\mathbf{x}^{(0)})^T \left(M(\mathbf{x}^{(0)}) + \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) \frac{\partial M(\mathbf{x}^{(0)})}{\partial x_j} \right) U_m(\mathbf{x}^{(0)}) \\ &= I + \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) U_m(\mathbf{x}^{(0)})^T \frac{\partial M(\mathbf{x}^{(0)})}{\partial x_j} U_m(\mathbf{x}^{(0)}). \end{aligned}$$

We can now derive a formula for an approximation $F_{m,0}(\mathbf{x})$ of $F_0(\mathbf{x})$ by combining all these considerations:

$$F_{m,0}(\mathbf{x}) = \hat{Z}(\mathbf{x})^T T_m(\mathbf{x}_0) \hat{Z}(\mathbf{x}) + \sum_{j=1}^{\ell} (x_j - x_j^{(0)}) \hat{G}_j(\mathbf{x}) \quad (13)$$

According to the previous remarks, $\hat{Z}(\mathbf{x})$, and $\hat{G}_j(\mathbf{x})$, can be computed cheaply in $\mathcal{O}(m^3)$ flops, so with a computational cost that is *independent* of the size of the original problem. In view of this analysis, we arrive at the following result.

Lemma 4.6. *The matrix-valued function $F_{m,0}(\mathbf{x})$ defined in (13) is a first-order approximation of $F_0(\mathbf{x})$ at $\mathbf{x} = \mathbf{x}^{(0)}$. More precisely, there exists a neighborhood $\mathcal{N}$ containing $\mathbf{x}^{(0)}$ such that*

$$F_0(\mathbf{x}) = F_{m,0}(\mathbf{x}) + \mathcal{O}(\|\mathbf{x} - \mathbf{x}^{(0)}\|^2), \quad \mathbf{x} \in \mathcal{N}.$$

4.3.2. Restarted Lanczos and other subspace iterations

At this stage, at no time we have used the property that $W_m(\mathbf{x})$ spans a Krylov subspace generated by $K(\mathbf{x})^{-1}$. In fact, this is not a requirement at all. Any subspace $W_m(\mathbf{x})$ generated through a subspace iteration method will work equally well in this framework.

In particular, we could rely on well-known large-scale eigenvalue solver such as ARPACK [30], which provides a tried-and-tested implementation of a restarted Lanczos method to compute the smaller end of the spectrum. This is what we currently use in the code.

Other approaches can be considered as well. The only requirement is that the projected matrix have the form

$$W_m(\mathbf{x})^T M(\mathbf{x})^{\frac{1}{2}} K(\mathbf{x})^{-1} M(\mathbf{x})^{\frac{1}{2}} W_m(\mathbf{x}),$$

which in turn implies that we are interested in the lowest end of the spectrum. For this reason, and in order to keep the explanation simple and self-contained, we decided to consider only the Lanczos method herein. Even if we never mention restarted variants explicitly, their use does not require any change in the proposed strategy.

4.3.3. Computing derivatives

In order to apply Algorithm 1 and, in particular, to perform the first-order correction (8), we need to evaluate $\nabla\phi(\mathbf{x}^{(k)})$ and $\nabla\phi_k^{\mathcal{R}}(\mathbf{x}^{(k)})$. In both cases, the problem can be rephrased in terms of the computation of a derivative of the eigenvalues of a pencil $K - \lambda M$, in view of

$$\frac{\partial}{\partial x_j} \left\| \frac{\sqrt{\Lambda_s(K(\mathbf{x}), M(\mathbf{x}))}}{2\pi} - \mathbf{f} \right\|_{\mathbf{w},2}^2 = \left(\frac{\sqrt{\Lambda_s(K(\mathbf{x}), M(\mathbf{x}))}}{2\pi} - \mathbf{f} \right)^T D_{\mathbf{w}}^2 \frac{\frac{\partial}{\partial x_j} \Lambda_s(K(\mathbf{x}), M(\mathbf{x}))}{2\pi \sqrt{\Lambda_s(K(\mathbf{x}), M(\mathbf{x}))}}, \quad (14)$$

where $D_{\mathbf{w}} = \text{diag}(w_1, \dots, w_s)$, as in (5). Classical perturbation theory for eigenvalues yields the following, for $i = 1, \dots, s$:

$$\left[\frac{\partial}{\partial x_j} \Lambda_s(K(\mathbf{x}), M(\mathbf{x})) \right]_i = \left(v_i^T \left(\frac{\partial K(\mathbf{x})}{\partial x_j} - \lambda_i \frac{\partial M(\mathbf{x})}{\partial x_j} \right) v_i \right) \cdot (v_i^T M(\mathbf{x}) v_i)^{-1},$$

where v_i is an eigenvector relative to the eigenvalues λ_i .

4.4. Constructing a local model for the trust-region scheme

Now that we have an approximation $F_{m,0}(\mathbf{x})$ of $F(\mathbf{x})$, we can obtain a local model for the trust-region scheme by replacing $\hat{\Lambda}_s(F(\mathbf{x}))$ with $\hat{\Lambda}_s(F_{m,0}(\mathbf{x}))$.

In view of Lemma 4.3, the approximation of $F(\mathbf{x})$ with $F_{m,0}(\mathbf{x})$ is correct only up to the zeroth order, and the same holds when composing it with $\hat{\Lambda}_s(\cdot)$, and so for $\phi(\mathbf{x})$ and $\phi^{\mathcal{R}}(\mathbf{x})$ as well. However, using the strategy described in Section 4.3.1, we can modify $\phi^{\mathcal{R}}(\mathbf{x})$ a posteriori to match at the first-order.

Since both $\phi(\mathbf{x})$ and $\phi^{\mathcal{R}}(\mathbf{x})$ are of the form described in equation (14), and in both cases the eigenvectors are available while computing the eigenvalues, the same formula can be used directly.

Lemma 4.7. *Let $\phi^{\mathcal{R}}(\mathbf{x})$ be the function defined as follows*

$$\phi^{\mathcal{R}}(\mathbf{x}) = \widetilde{\phi}^{\mathcal{R}}(\mathbf{x}) + (\nabla\phi(\mathbf{x}^{(0)}) - \nabla\widetilde{\phi}^{\mathcal{R}}(\mathbf{x}^{(0)}))^T (\mathbf{x} - \mathbf{x}^{(0)}), \quad \widetilde{\phi}^{\mathcal{R}}(\mathbf{x}) = \left\| \frac{\sqrt{\hat{\Lambda}_s(F_{m,0}(\mathbf{x}))}}{2\pi} - \mathbf{f} \right\|_{\mathbf{w},2}^2. \quad (15)$$

Then, $\phi^{\mathcal{R}}(\mathbf{x})$ is a first-order approximation of $\phi(\mathbf{x})$ at $\mathbf{x} = \mathbf{x}^{(0)}$.

Proof. In view of Lemma 4.3, 4.4, 4.6 we have $\phi(\mathbf{x}^{(0)}) = \widetilde{\phi}^{\mathcal{R}}(\mathbf{x}^{(0)})$. Given this condition, the first-order matching is ensured by the gradient correction done as described with Section 4.3.1. $\square$

Remark 4.8. Although the formulas for $\nabla\phi(\mathbf{x})$ and $\nabla\phi^{\mathcal{R}}(\mathbf{x})$ are not reported explicitly in Lemma 4.7, they are directly implementable in view of Equation (14), and require almost no computational effort when the eigenvectors are obtained through the Lanczos process.

5. Numerical experiments and real examples

In this section we verify the performance of the proposed approach on both artificial and real examples. To this end, we will compare the number of iterations needed to achieve convergence, which is reported in terms of outer iterations, that is to say, the number of reduced models computed in order to complete the procedure.

The number of iterations of the inner optimizer, used within a single trust-region, are ignored because they are largely irrelevant in determining the final computational cost. Typical tests show that the total time spent optimizing the models within the trust-region is less than 1% of the total running time of the algorithm.

5.1. Some implementation details

The tests were run on a computer with an Intel Core i7-920 CPU running at 2.67 GHz, with 18 GB of RAM clocked at 1066 MHz. We used the single-core version of MUMPS 4.10, and the Intel MKL BLAS shipped with MATLAB R2017b.

In all experiments the tolerance τ for the accuracy of the Lanczos method was set to 10^{-5} while the trust-region procedure was stopped when the norm of the projected gradient fell below 10^{-4} (i.e. $\epsilon = 10^{-4}$ in Algorithm 1). All the other parameters in Algorithm 1 were set to standard values as suggested in [17, Chapter 17]. The trust-region radius update follows [17, Chapter 17, page 782] as well. Also, in Algorithm 1 the trust-region step $\mathbf{s}^{(k)}$ is computed “exactly”, that is, we minimize the reduced model (15) in $\mathcal{B}_k \cap \Omega$ to high accuracy. This is a reasonable choice due to the low dimension of the problem. To this end, we used the function `fmincon` included in the MATLAB’s Optimization toolbox, setting the built-in `sqp` solver and using the default parameter settings.

Furthermore, the parameter space is always preliminary scaled, so that the initial point $\mathbf{x}^{(0)}$ is the vector of all ones. This ensures that checking norm-wise conditions on the eigenvalues and on the gradient yields relative accuracy on the parameters, independently of their scaling.

The weight vector $\mathbf{w}$ is always chosen as $w_i = f_i^{-1}$, which ensures relative accuracy on the recovered frequency, except where otherwise stated. In particular, in the clock tower example, we emphasize the consequences of a different choice of $\mathbf{w}$.

5.2. Arch on piers

In this section we consider a simple example for demonstration purposes. We built a 2D discretization of the arch on piers, shown in Figure 1. The arch spans 4 meters, and rests on two 4-meters-high lateral piers. The structure is modeled by means of 336 finite elements, and clamped at the base of the two piers. This corresponds to 851 total degrees of freedom in the structure. We have used plane strain 4-node quadrilateral elements.

The three materials composing the structure are depicted with different colors in Figure 1. Material 1 is used for the arch (green in the figure), and materials 2 and 3 for the piers (red and blue, respectively). Poisson’s ratio ν is set to 0.2 for all the materials, and assumed to be known a priori.

For the three materials we assume the following values:

$$\begin{array}{lll} E_1 = 3250 \text{ MPa} & E_2 = 5000 \text{ MPa} & E_3 = 4800 \text{ MPa} \\ \rho_1 = 1800 \frac{\text{kg}}{\text{m}^3} & \rho_2 = 2200 \frac{\text{kg}}{\text{m}^3} & \rho_3 = 2100 \frac{\text{kg}}{\text{m}^3}, \end{array}$$

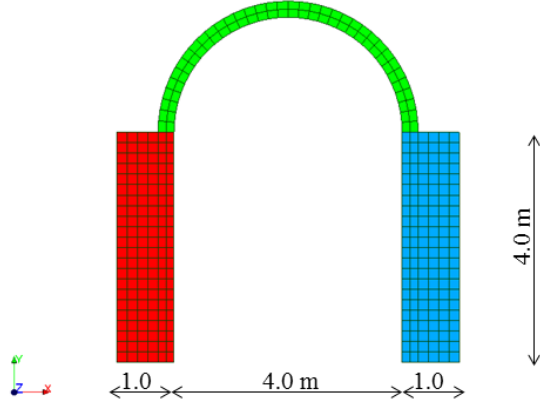


Figure 1: The arch on piers modeled in the experiments. Each differently colored region corresponds to a different material.

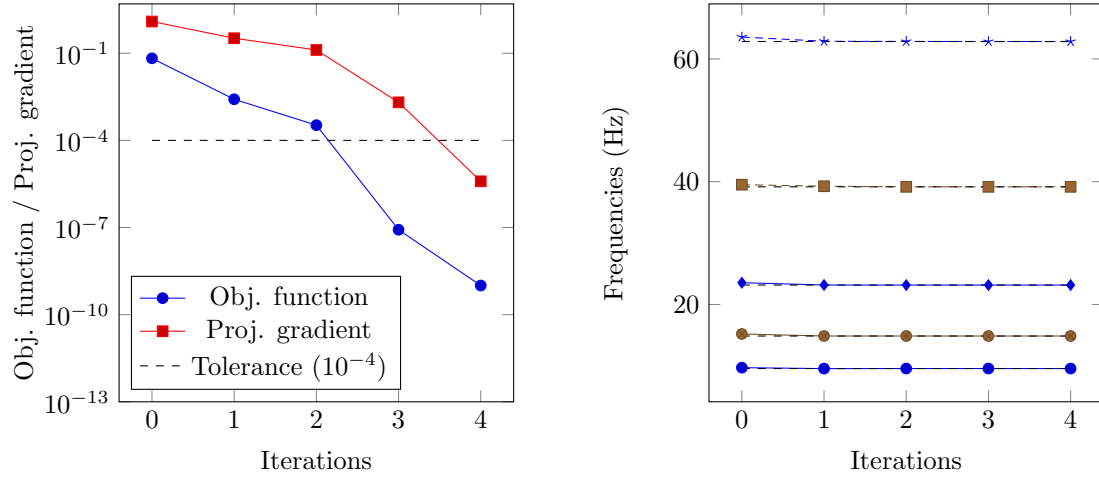


Figure 2: The left graph shows the convergence of the objective function to the minimum for each new reduced model generated in the optimization procedure. The dashed line is the tolerance set for the projected gradient in the optimization scheme. The right graph plots the convergence of the frequencies during the process.

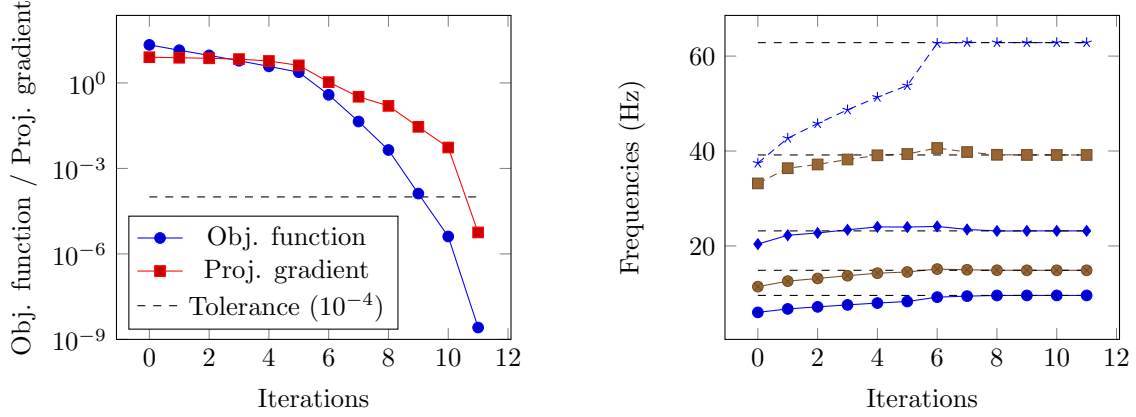


Figure 3: The left graph shows the convergence of the objective function to the minimum for each new reduced model generated in the optimization procedure. The dashed line is the tolerance set for the projected gradient in the optimization scheme. The right graph plots the convergence of the frequencies during the process. In this case, the initial points were chosen quite far from the original ones, selecting $E_2 = 2000 \text{ MPa}$, $\rho_2 = 1100 \text{ kg} \cdot \text{m}^{-3}$, $E_3 = 1100 \text{ MPa}$.

where the index $j \in \{1, 2, 3\}$ indicates the material under consideration. We computed the 5 leading frequencies by running the NOSA-ITACA code (with these fixed parameters), to obtain:

$$\mathbf{f} \approx [9.575 \quad 14.87 \quad 23.17 \quad 39.17 \quad 62.84] .$$

Our aim is to validate the optimization method by recovering the original parameters matching these frequencies, assuming E_2 , ρ_2 , and E_3 are unknown. We consider the following bounds of realistic parameters:

$$1000 \text{ MPa} \leq E_2 \leq 9000 \text{ MPa}, \quad 1000 \frac{\text{kg}}{\text{m}^3} \leq \rho_2 \leq 3000 \frac{\text{kg}}{\text{m}^3}, \quad 1000 \text{ MPa} \leq E_3 \leq 9000 \text{ MPa},$$

which define the corresponding box Ω in (3)

In the default implementation, the starting points have been chosen the midpoint of the intervals, which are quite good estimates of the true values. Convergence in this case is displayed in Figure 2. The initial points are sufficiently close to the correct ones for the frequencies to almost match already, and the distance to the optimum during convergence is not easily distinguishable. However, the left plot in Figure 2, which shows the convergence of the objective function in log-scale, clearly demonstrates that we are recovering the right solution.

To test the robustness of the approach, we modified the starting points to be relatively far from the correct ones by making the following choices:

$$E_2^{(0)} = 2000 \text{ MPa}, \quad \rho_2^{(0)} = 1100 \frac{\text{kg}}{\text{m}^3}, \quad E_3^{(0)} = 1100 \text{ MPa}.$$

Convergence of the method is displayed in the graphs in Figure 3, which clearly shows that the initial frequencies are quite far from the correct ones; yet the method reaches convergence using only 12 reduced models. Moreover, a rough estimate (though probably sufficient for most engineering

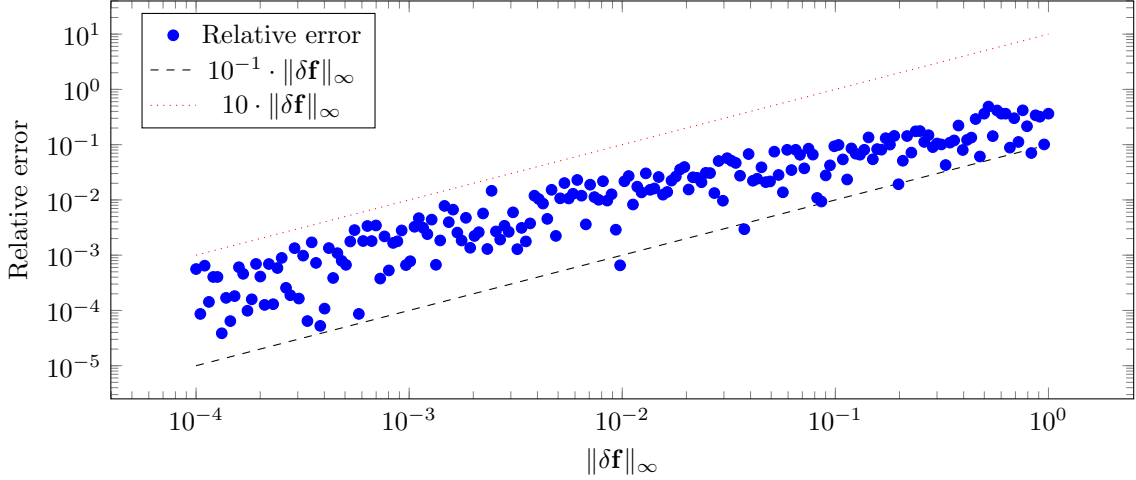


Figure 4: This plot shows the relation between the noise affecting the frequency measurements and the component-wise relative error on the retrieved optimal parameters. The graph shows the linear dependency of the error on the estimated parameters and the noise level of the input data.

purposes) is already obtained after about 6 steps. Moreover, the Young’s modulus and the mass density of the initial model are recovered exactly up to 5 digits in this example.

Another important feature for a method that must deal with experimentally measured frequencies is its robustness when the input is subject to noise. In fact, we would expect the frequencies to be accurate only up to a certain relative threshold.

To simulate the behavior of the method in a predictable environment, we perturbed the frequency vector $\hat{\mathbf{f}} = \mathbf{f} + \delta\mathbf{f}$, by imposing $|\delta f_i| \leq |f_i| \cdot \delta$, with δ the prescribed noise level. We have run tests for δ ranging from 0.01% to 100%: the corresponding error in the retrieved frequencies is plotted in Figure 4.

The error is measured in relative norm, that is we have plotted the infinity norm of the vector with components $(x_i - \hat{x}_i)/x_i$, where x_i is the vector with the actual model parameters, and $\hat{x}_i$ those estimated by the optimization process.

The linear behavior of the relative error with respect to the noise level is clearly visible. This behavior suggests that, even if the frequency measurements are contaminated with noise, it is still possible to retrieve meaningful parameters if enough information is provided. Note that, in general, correct (or accurate) recovery of the parameters might not be possible, and a more detailed study of the condition number of this inverse problem will be object of further in-depth study in future research.

5.3. A dome

Here we provide a more complex example, represented by a dome supported by four 14-meters-high pillars. The dome consists of an octagonal shaped cloister vault 5 meters high, resting on a drum inscribed on a 10×11 meters rectangle. We assume the structure to be composed of 4 different materials: material 1 for the vault (red in Figure 5), material 2 for the top of the drum (blue), material 3 for the lower part of the drum (green) and material 4 for the pillars (gray).

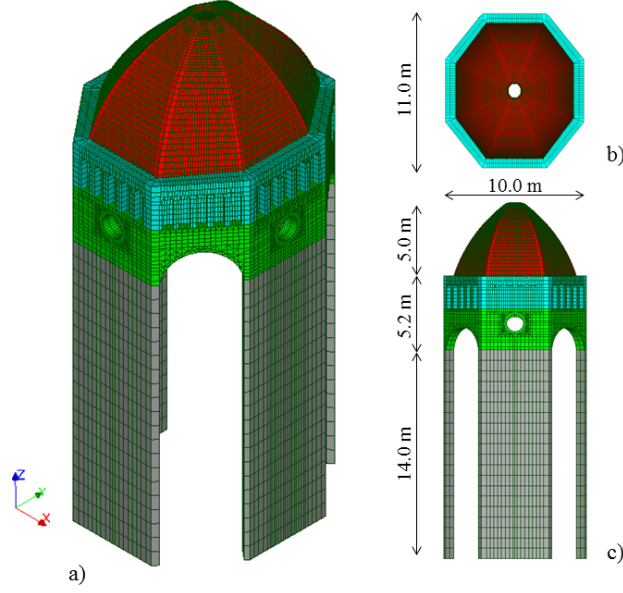


Figure 5: Domed temple

The finite element model consists of 31,052 elements and 41,245 nodes. It is modeled using 3D 8-node hexahedron brick elements, which gives rise to a (projected) model with 122.853 degrees of freedom.

We performed a test similar to that carried out for the arch on piers: we computed the frequencies for a certain choice of parameters, some of which were assumed to be unknown, and ran the optimization scheme trying to recover the original parameters.

In order to obtain the reference frequencies the characteristics of the 4 materials (Young's modulus and mass density) were set as follows:

$$\begin{aligned}
 E_1 &= 3000 \text{ MPa} & \rho_1 &= 1800 \frac{\text{kg}}{\text{m}^3} & E_2 &= 4000 \text{ MPa} & \rho_2 &= 2000 \frac{\text{kg}}{\text{m}^3} \\
 E_3 &= 3500 \text{ MPa} & \rho_3 &= 1900 \frac{\text{kg}}{\text{m}^3} & E_4 &= 5000 \text{ MPa} & \rho_4 &= 2200 \frac{\text{kg}}{\text{m}^3} \\
 \nu_j &= 0.25, & j &= 1, \dots, 4.
 \end{aligned}$$

The leading 10 frequencies have been computed using the NOSA-ITACA code. The result, expressed in Hertz, are the frequencies in the vector:

$$\mathbf{f} \approx [2.19 \quad 2.23 \quad 3.76 \quad 3.83 \quad 4.32 \quad 4.60 \quad 4.72 \quad 8.26 \quad 8.30 \quad 9.21].$$

The optimization code was run setting the bounds as

$$2000 \text{ MPa} \leq E_j \leq 6000 \text{ MPa} \quad 1600 \frac{\text{kg}}{\text{m}^3} \leq \rho_j \leq 2400 \frac{\text{kg}}{\text{m}^3} \quad j = 1, \dots, 4,$$

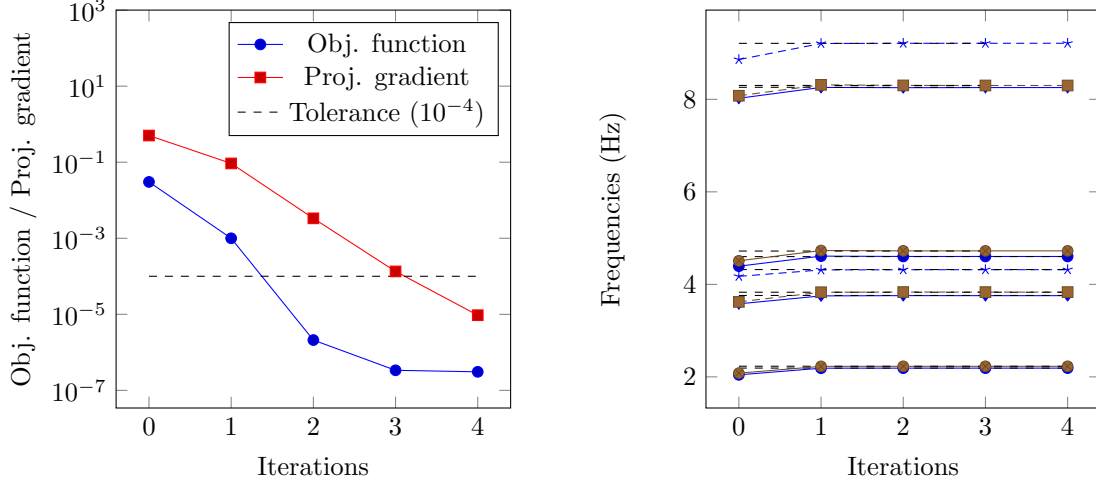


Figure 6: Convergence history for the dome model, matching 10 frequencies with 7 free parameters. The left plot shows the convergence of the objective function and the norm of the projected gradient, and in the right plot the convergence of frequencies to the exact ones is reported.

with the sole exception of ρ_3 which was set to the correct value. This leaves 7 parameters to be optimized. No particular starting conditions were specified. In this case, the algorithm chooses the midpoint of the intervals. The convergence history is reported in Figure 6, which reveals that high accuracy is attained with only 4 outer iterations — with the residual of the objective function being very small even after two steps. The weights have been chosen in order to reach relative accuracy³ (i.e., $w_i = f_i^{-1}$) on the target frequencies, with the usual tolerance $\epsilon = 10^{-4}$.

The estimated mechanical parameters obtained through the optimization are as follows:

$$\begin{aligned}
 E_1 &= 2953 \text{ MPa} & \rho_1 &= 1769 \frac{\text{kg}}{\text{m}^3} & E_2 &= 3932 \text{ MPa} & \rho_2 &= 1960 \frac{\text{kg}}{\text{m}^3} \\
 E_3 &= 3426 \text{ MPa} & E_4 &= 4951 \text{ MPa} & \rho_4 &= 2174 \frac{\text{kg}}{\text{m}^3}.
 \end{aligned}$$

This corresponds to a relative error between 0.97% and 2.1%, with an average of about 1.6%. Although such accuracy is quite good for practical purposes, even more precise estimates can be obtained by reducing the tolerance to below 10^{-4} that we have used in these tests.

5.4. Examples with multiple eigenvalues

In this section we test the code on two examples that involve multiple eigenvalues. In the first example, we consider a model with a structural symmetry that is preserved when the materials parameters change. This corresponds to an eigenvalue that has constant multiplicity for all the parameters, and therefore the convergence is guaranteed by Lemma 4.4.

³A detailed discussion of the influence of the weights on the accuracy of the recovered frequencies is presented in the next section, where its relevance is discussed for a real example.

In the second case, we consider a model where the symmetry is present only for a suitable choice of materials parameters. In this case, the multiplicity of the eigenvalue is not constant in a neighborhood of the optimal point, and therefore we cannot ensure the convergence of the approach. Nevertheless, we verify that the method works well in practice.

For both cases, the model is a 45 meters high masonry tower having a square cross section of 8 meters long side and constant wall thickness of about 2 meters. The tower is discretized into 1440 thick shell elements with 8832 degrees of freedom. The structure is assumed to be clamped at the base.

In the first test (Test 1), we suppose that the bottom part (34m high) is made of material 1 with Young's modulus E_1 , while the upper part made of material 2 with Young's modulus E_2 .

We choose the following values for the material properties:

$$E_1 = 3000 \text{ MPa} \quad E_2 = 1000 \text{ MPa} \quad \rho_1 = \rho_2 = 1800 \frac{\text{kg}}{\text{m}^3} \quad \nu_1 = \nu_2 = 0.2.$$

The corresponding frequencies to match are:

$$\mathbf{f} := [1.08729 \ 1.08729 \ 4.035].$$

The interval of valid parameters is chosen as

$$10^2 \text{ MPa} \leq E_1, E_2 \leq 9 \cdot 10^3 \text{ MPa},$$

for both materials. The convergence histories are reported in Figure 7 for the starting point $E_1^{(0)} = 1600 \text{ MPa}$ and $E_2^{(0)} = 200 \text{ MPa}$, and in Figure 8 for the starting points with Young's moduli $E_1^{(0)} = 6000 \text{ MPa}$ and $E_2^{(0)} = 500 \text{ MPa}$. The points have been chosen quite far from the true one to create a more difficult test for the method. We have also tested the starting points $E_1^{(0)} = 3500 \text{ MPa}$ and $E_2^{(0)} = 1500 \text{ MPa}$, and the convergence is reached in 4 steps.

As a second test (Test 2), we use the same model but now assuming a different materials distribution: the west and south facades are made of material 1, and the north and east of material 2. We assume that

$$E_1 = E_2 = 3000 \text{ MPa}.$$

For this choice of the Young's moduli, due to the symmetry, the structure has a double eigenvalue that, unlike the former example, does not maintain its multiplicity in a neighborhood (if the Young's moduli do not coincide). The first three frequencies have been computed numerically and are equal to

$$\mathbf{f} := [1.0945 \ 1.0945 \ 4.1360].$$

The theory cannot ensure convergence in this case, but nevertheless the proposed method works quite well in practice. Using the same interval as before

$$10^2 \text{ MPa} \leq E_1, E_2 \leq 9 \cdot 10^3 \text{ MPa},$$

for the Young's moduli, we have tested different starting points, listed in Table 1. In all cases the recovered solution provided the correct frequencies up to 4 digits. In addition, when starting from an initial point where the symmetry is present (that is, $E_1^{(0)} = E_2^{(0)}$) the method converges to a critical point which is closer to the original one. The convergence history for the test case 2 is reported in Figure 9.

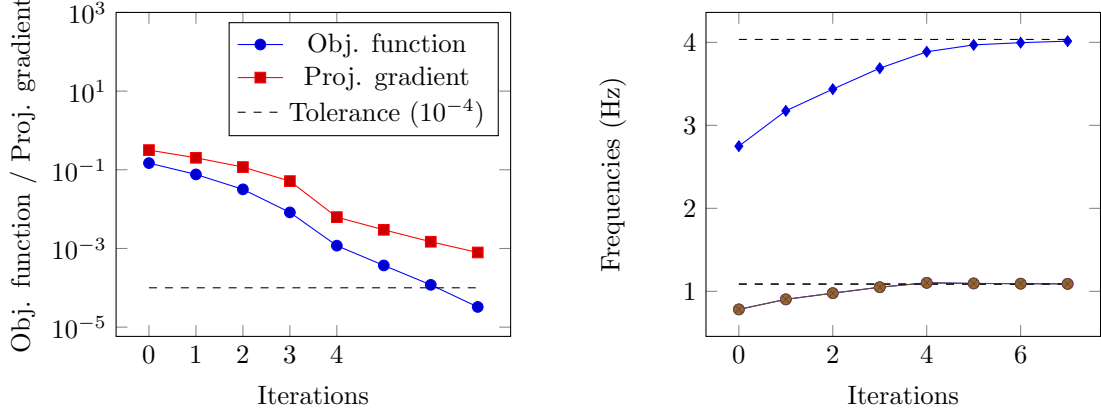


Figure 7: Test 1. Convergence history for the tower model with a double structural eigenvalue at about 1.08 Hz, matching 3 frequencies with 2 free parameters. The starting point is $E_1^{(0)} = 1.6 \cdot 10^9 \text{ Pa}$ and $E_2^{(0)} = 2 \cdot 10^8 \text{ Pa}$. The left plot shows the convergence of the objective function and the norm of the projected gradient, and in the right plot the convergence of frequencies to the exact ones is reported.

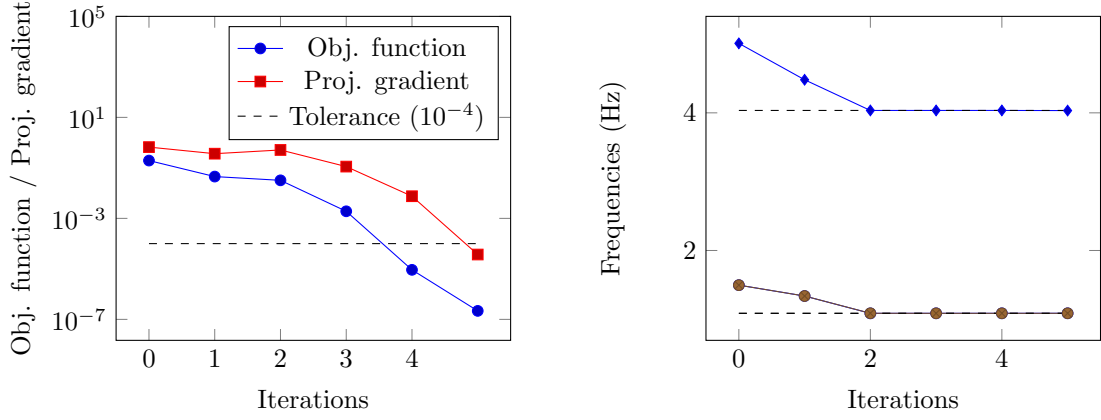


Figure 8: Test 1. Convergence history for the tower model with a double structural eigenvalue at about 1.08 Hz, matching 3 frequencies with 2 free parameters. The starting point is $E_1^{(0)} = 6 \cdot 10^9 \text{ Pa}$ and $E_2^{(0)} = 5 \cdot 10^8 \text{ Pa}$. The left plot shows the convergence of the objective function and the norm of the projected gradient, and in the right plot the convergence of frequencies to the exact ones is reported.

Case	Starting $E_1^{(0)}$	Starting $E_2^{(0)}$	Recovered E_1	Recovered E_2	Iterations
1	$3.5 \cdot 10^3 \text{ MPa}$	$3.5 \cdot 10^3 \text{ MPa}$	$3.0 \cdot 10^3 \text{ MPa}$	$3.0 \cdot 10^3 \text{ MPa}$	2
2	$1.2 \cdot 10^3 \text{ MPa}$	$6 \cdot 10^3 \text{ MPa}$	$2.92 \cdot 10^3 \text{ MPa}$	$3.07 \cdot 10^3 \text{ MPa}$	7
3	$1.0 \cdot 10^3 \text{ MPa}$	$2.0 \cdot 10^3 \text{ MPa}$	$2.87 \cdot 10^3 \text{ MPa}$	$3.12 \cdot 10^3 \text{ MPa}$	7
4	$5 \cdot 10^2 \text{ MPa}$	$5 \cdot 10^2 \text{ MPa}$	$3.0 \cdot 10^3 \text{ MPa}$	$3.0 \cdot 10^3 \text{ MPa}$	13

Table 1: Test 2. Number of iterations and recovered parameters for the model with multiple eigenvalues at the solution, but not in a neighborhood. The results for several different cases are reported. The optimal Young's moduli for the given frequencies are $3.0 \cdot 10^3 \text{ MPa}$.

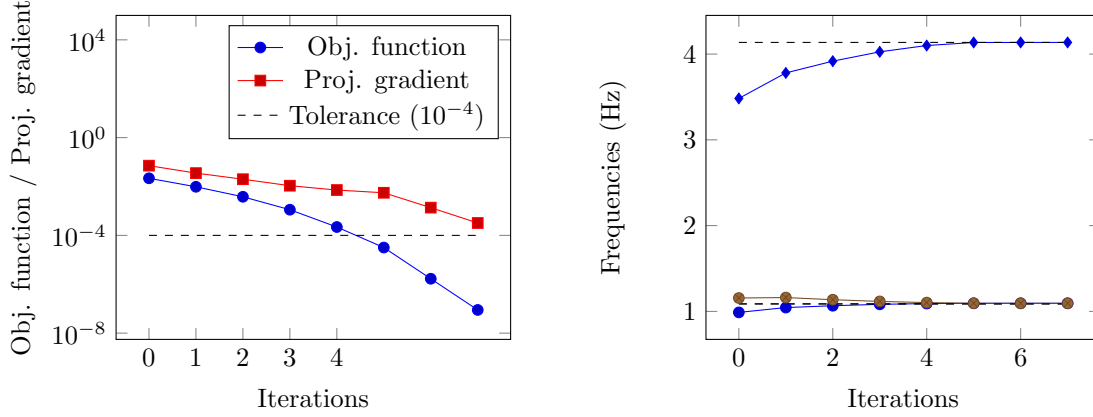


Figure 9: Test 2. Convergence history for the tower model with a double structural eigenvalue at about 1.0945 Hz, matching 3 frequencies with 2 free parameters. The starting points are Young’s moduli $E_1^{(0)} = 1.2 \cdot 10^3$ MPa and $E_2^{(0)} = 6 \cdot 10^3$ MPa. The left plot shows the convergence of the objective function and the norm of the projected gradient, and in the right plot the convergence of frequencies to the exact ones is reported.

5.5. The Clock tower

As a real example, we consider the case study of the Clock tower (“Torre delle Ore”) in Lucca, Italy. The 48.4 m high masonry structure has a rectangular cross section of about 5.1×7.1 meters and walls of thickness varying from 1.77 m at the base to 0.85 m at the top. Two barrel vaults are set at heights of about 12.5 m and 42.3 m. The bell chamber is covered by a pavillon roof made up of wooden trusses and rafters. At a height of about 33 m the walls have been instrumented with 4 steel tie rods of 30×30 mm rectangular section. The adjacent buildings about the tower on two sides for a height of about 13 m and constitute asymmetric boundary conditions. The modal behavior of the tower has already been analyzed using the NOSA-ITACA code and through experimental measurements conducted in November 2016, when the tower was fitted with 4 triaxial seismometric stations. The two procedures yielded similar, albeit not perfectly matching results [34]. Figure 10 shows a picture of the tower, together with the mesh used for its discretization.

The frequencies obtained by analyzing data from the fitted instruments via Operational Modal Analysis techniques, are as follows:

$$\mathbf{f} \approx [1.05 \quad 1.3 \quad 4.19 \quad 4.50].$$

The frequencies obtained with the finite element model in [34] are, instead, $[0.98 \quad 1.24 \quad 4.32 \quad 4.38]$. The materials constituting the tower are wood, steel, and masonry, with the latter being different in the bell chamber and the lower part of the tower. The mechanical properties used in the finite element model are

$$\begin{array}{lll} E_w = 10000 \text{ MPa} & \rho_w = 800 \frac{\text{kg}}{\text{m}^3} & \nu_w = 0.35 \\ \rho_{m_l} = 2100 \frac{\text{kg}}{\text{m}^3} & \rho_{m_c} = 1700 \frac{\text{kg}}{\text{m}^3} & \nu_{m_l} = \nu_{m_c} = 0.2 \\ E_s = 210,000 \text{ MPa} & \rho_s = 7850 \frac{\text{kg}}{\text{m}^3} & \nu_s = 0.3, \end{array}$$



Figure 10: On the left, a picture of the clock tower in Lucca. On the right, the mesh used for the finite element discretization, where the different materials for the bell chamber and the lower part of the tower are reported in different colors (respectively red and light blue).

and we leave the Young's moduli E_{m_c} and E_{m_l} as free parameters, since they are not known explicitly due to the lack of experimental information. Here, the subscript m_l identifies the lower part of the tower, and m_c the bell chamber.

We can reformulate the problem as an optimization one, trying to match the frequencies of the system to the measured ones, allowing the parameters to vary in a reasonable range. In our case we can set

$$2500 \text{ MPa} \leq E_{m_l} \leq 5500 \text{ MPa} \quad 1000 \text{ MPa} \leq E_{m_c} \leq 5500 \text{ MPa}.$$

As a first test, we choose the weights as the vector of all ones. The convergence of the optimization scheme, obtained with these initial parameters, is reported in Figure 11. Note that the method optimizes the third and fourth frequencies to match quite closely, but at the price of decreasing the quality of the match for the first two (see Table 2).

This is connected to the choice of $\mathbf{w}$, which ensures norm-wise accuracy, but not a relative one. Often, we prefer to have a low relative error on the frequencies, instead of an absolute one. Therefore, it is natural to choose $w_i = f_i^{-1}$ (in fact, this is the default choice in the implementation). This yields different results, and different approximations for both the parameters and the frequencies. More specifically, we expect a better match on the lowest end of the spectrum, and a worse one on the highest frequencies.

The convergence history with this choice of $\mathbf{w}$ is reported in Figure 12, and the results are shown in Table 2. It appears immediately evident that the relative errors are balanced with $w_i = f_i^{-1}$, whereas they tend to increase on the small frequencies when $w_i = 1$.

The parameters identified by the model updating phase give $E_{m_l} \approx 3182 \text{ MPa}$ and $E_{m_c} \approx 1873 \text{ MPa}$ when using weights equal to one, and $E_{m_l} \approx 3076 \text{ MPa}$ and $E_{m_c} \approx 1950 \text{ MPa}$ when aiming for relative accuracy.

We conclude this section by comparing the overall CPU time of our implementation with those of alternative solvers, which in various ways fail to exploit the problem structure. Table 3 compares

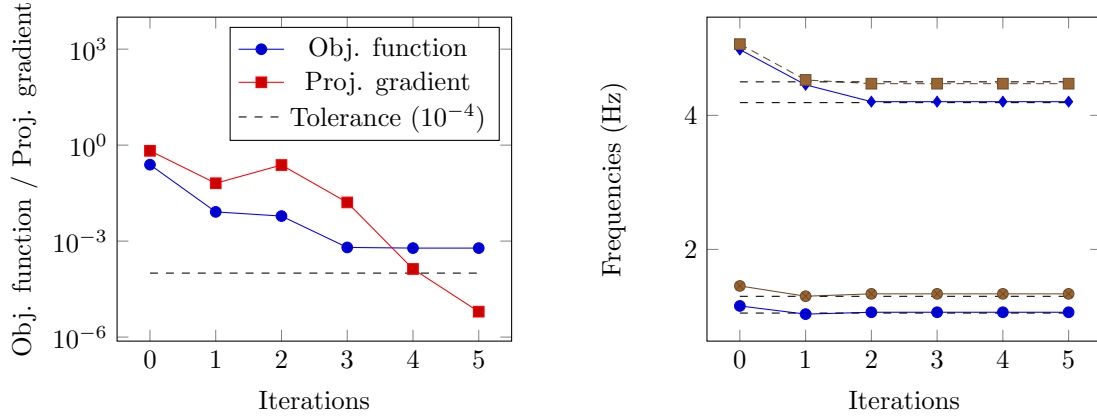


Figure 11: The left graph reports the convergence history of the objective function to the minimum for each new reduced model generated in the optimization procedure. The norm of the projected gradient is reported as well, and the dashed line is the tolerance of the optimization scheme. The right plot represents the convergence of the frequencies during the process (the dashed lines are the experimental frequencies). Iteration 0 is the evaluation function at the starting point. In this example, the free parameters are E_{m_l} and E_{m_c} .

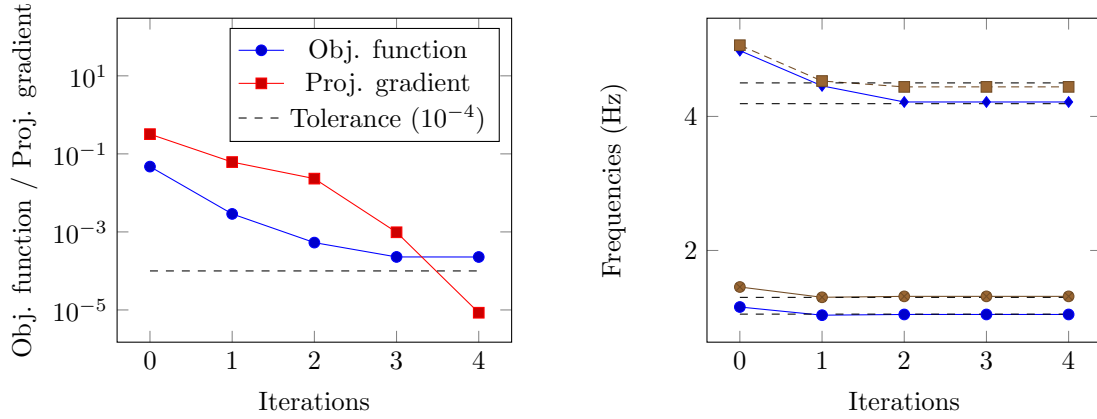


Figure 12: The left graph reports the convergence history of the objective function to the minimum for each new reduced model generated in the optimization procedure. The norm of the projected gradient is reported as well, and the dashed line is the tolerance set for the projected gradient in the optimization scheme. The right plot represents the convergence of the frequencies during the process (the dashed line are the experimental frequencies). Iteration 0 is the evaluation function at the starting point. In this example, the free parameters are E_{m_l} and E_{m_c} , and the objective function is defined using the weights $w_i = f_i^{-1}$.

Exp. freq.	Freq. ($w_i = 1$)	Rel. error ($w_i = 1$)	Freq. ($w_i = f_i^{-1}$)	Rel. error ($w_i = f_i^{-1}$)
1.05 Hz	1.0621 Hz	1.15%	1.0449 Hz	0.49%
1.30 Hz	1.3366 Hz	2.82%	1.315 Hz	1.15%
4.19 Hz	4.2041 Hz	0.33%	4.2154 Hz	0.61%
4.50 Hz	4.4729 Hz	0.60%	4.4409 Hz	1.31%

Table 2: Frequencies and corresponding relative errors computed using the weights $w_i = 1$ and $w_i = f_i^{-1}$ for the problem of the clock tower in Lucca, Italy.

4 different implementations:

RM This implementation corresponds to the current proposal, that is, a solver based on the trust-region and projection approach described in the previous section and used in the tests presented herein. This strategy is labeled as “RM” in the table — highlighted with the bold font.

BB Here we interpret the finite element code as a black-box function and solve the optimization problem using a general purpose optimizer. This is done to reflect a scenario in which the user has no access to the internals of the finite element code — and requires an assembly phase at every function evaluation. Derivatives are computed here by finite differences. This is column “BB” in the table.

A In this approach, we separately assemble the parametric generalized eigenvalue problem $K(\mathbf{x}) - \lambda M(\mathbf{x})$ and then pass to a general optimizer a function that evaluates the frequencies for a given value of $\mathbf{x}$. This function corresponds to the Lanczos projection routine used internally in RM. Derivatives are computed here by finite differences. This is approach “A” in the table.

AD The same as A, but derivatives are evaluated analytically as described to Section 4.3.3. This is “AD” in the table.

The general optimizer chosen for the BB, A and AD implementations is again `fmincon` with the `sqp` option available in the Optimization toolbox of MATLAB. In our experience, this choice turns out to be the most efficient among those available in the toolbox. This same function has been used for computation of the step in the trust-region scheme of Algorithm 1.

In Table 3, the time for assembly (which needs to be added to the costs in columns A, AD, and RM) is reported separately, as well as the problem’s degrees of freedom.

It is immediately apparent that the proposed RM method is the fastest in all cases except the arch on piers, where given the small dimension incurs in some overhead. Moreover, it is worth stressing that our approach exhibits a complexity that seems to grow nicely with the dimension. The general optimization approach BB, instead, becomes costly when the number of degrees of freedom increases.

For instance, in the dome example, the timings range from about 7 hours for BB to slightly more than 3 minutes with our approach. The case where we supply derivatives to the optimizer (AD) still takes more than 30 minutes.

The timings suggest the unsurprising conclusion that better knowledge of the underlying optimization problem leads to more efficient and faster optimization routines. It should be noted that the stopping criterion for all the approaches is set to ensure that the first order optimality condition (given by the norm of the projected gradient) is smaller than $\epsilon = 10^{-4}$ used in the experiments. All the approaches provide solutions with comparable accuracy.

	# dof	BB	A	AD	RM	Assembly
Arch on piers	851	16.40s	1.17s	0.62s	0.68s	0.42s
Dome	122,853	25,032s	7,662.4s	1,990.5s	226.02s	134.32s
Clock tower	45,511	476.9s	305.39s	202.87s	44.10s	32.22s

Table 3: Timings for different optimization approaches. The timings reported for the approaches “Assembled” (A), “Assembled + derivatives” (AD), and “Reduced model” (RM) do not include the time needed to assembly the parametric finite element model, which is reported in the last column. The Portal model has been run with the default starting points.

6. Concluding remarks

In this work we have presented a model updating scheme focusing on the optimization of finite element models for structural analysis purposes. We have shown how the subspace projection method used to compute the eigenvalues at the lowest end of the spectrum can be used directly as a parametric model order reduction step, and that embedding it in a trust-region scheme can be very effective.

Embedding the model updating step directly in the finite element code naturally leads to more efficient procedures, compared to applying an optimizer fed with a black-box finite element code. In the proposed approach, we benefit from the fact that the FE code at our disposal not only provides the mass and stiffness matrices for certain values of the parameters, but also allows building the parametric model in a single step, instead of assembling the matrices for different parameter values over and over.

Several examples have been reported on, both to test the theoretical properties of the scheme as well as demonstrate its practical applicability and its good performance with respect to alternative solvers. We have also presented a preliminary numerical evidence of its robustness against perturbations in the data, an aspect which will be further investigated and characterized in future work.

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