A CONVERGENT ADAPTIVE METHOD FOR ELLIPTIC EIGENVALUE PROBLEMS

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Abstract. We prove the convergence of an adaptive linear finite element method for computing eigenvalues and eigenfunctions of second-order symmetric elliptic partial differential operators. The weak form is assumed to yield a bilinear form which is bounded and coercive in $H^1$. Each step of the adaptive procedure refines elements in which a standard a posteriori error estimator is large and also refines elements in which the computed eigenfunction has high oscillation. The error analysis extends the theory of convergence of adaptive methods for linear elliptic source problems to elliptic eigenvalue problems, and in particular deals with various complications which arise essentially from the nonlinearity of the eigenvalue problem. Because of this nonlinearity, the convergence result holds under the assumption that the initial finite element mesh is sufficiently fine.

Key words. second-order elliptic problems, eigenvalues, adaptive finite element methods, convergence

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1. Introduction. In the last decades, mesh adaptivity has been widely used to improve the accuracy of numerical solutions to many scientific problems. The basic idea is to refine the mesh only where the error is high, with the aim of achieving an accurate solution using an optimal number of degrees of freedom. There is a large amount of numerical analysis literature on adaptivity, in particular on reliable and efficient a posteriori error estimates (e.g., [1]). Recently, the question of convergence of adaptive methods has received intensive interest and a number of convergence results for the adaptive solution of boundary value problems have appeared (e.g., [8, 18, 19, 7, 6, 23]).

We prove here the convergence of an adaptive linear finite element algorithm for computing eigenvalues and eigenvectors of scalar symmetric elliptic partial differential operators in bounded polygonal or polyhedral domains, subject to Dirichlet boundary data. Such problems arise in many applications, e.g., resonance problems, nuclear reactor criticality, and the modelling of photonic band gap materials, to name but three.

Our refinement procedure is based on two locally defined quantities, firstly, a standard a posteriori error estimator and secondly a measure of the variability (or “oscillation”) of the computed eigenfunction. (Measures of “data oscillation” appear in the theory of adaptivity for boundary value problems, e.g., [18]. In the eigenvalue problem the computed eigenvalue and eigenfunction on the present mesh plays the role of “data” for the next iteration of the adaptive procedure.) Our algorithm performs local refinement on all elements on which the minimum of these two local quantities is sufficiently large. We prove that the adaptive method converges provided the initial mesh is sufficiently fine. The latter condition, while absent for adaptive methods for

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linear symmetric elliptic boundary value problems, commonly appears for nonlinear problems and can be thought of as a manifestation of the nonlinearity of the eigenvalue problem.

We believe that the present paper is the first contribution to the topic of convergence of adaptive methods for eigenvalue problems. Since writing this paper, substantial improvements in the theory have been made in [5], where the need to adapt on the oscillations of the eigenvalue is removed and, in addition, the general convergence of the adaptive scheme to a nonspurious eigenvalue of the continuous problem is established.

The outline of the paper is as follows. In section 2 we briefly describe the model elliptic eigenvalue problem and the numerical method and in section 3 we describe a priori estimates, most of which are classical. Section 4 describes the a posteriori estimates and the adaptive algorithm. Section 5 proves that proceeding from one mesh to another ensures error reduction (up to oscillation of the computed eigenfunction) while the convergence result is presented in section 6. Numerical experiments illustrating the theory are presented in section 7.

2. Eigenvalue problem and numerical method. Throughout, $\Omega$ will denote a bounded domain in $\mathbb{R}^d$ ($d = 2$ or 3). In fact, $\Omega$ will be assumed to be a polygon ($d = 2$) or polyhedron ($d = 3$). We will be concerned with the problem of finding an eigenvalue $\lambda \in \mathbb{R}$ and eigenfunction $0 \neq u \in H^1_0(\Omega)$ satisfying

\begin{equation}
(2.1) \quad a(u, v) := \lambda b(u, v), \quad \text{for all} \quad v \in H^1_0(\Omega),
\end{equation}

where, for real valued functions $u$ and $v$,

\begin{equation}
(2.2) \quad a(u, v) = \int_{\Omega} \nabla u(x)^T A(x) \nabla v(x) \, dx \quad \text{and} \quad b(u, v) = \int_{\Omega} B(x) u(x) v(x) \, dx.
\end{equation}

Here, the matrix-valued function $A$ is required to be uniformly positive definite, i.e.,

\begin{equation}
(2.3) \quad 0 < a \leq \xi^T A(x) \xi \leq \overline{a} \quad \text{for all} \quad \xi \in \mathbb{R}^d \quad \text{with} \quad |\xi| = 1 \quad \text{and all} \quad x \in \Omega.
\end{equation}

The scalar function $B$ is required to be bounded above and below by positive constants for all $x \in \Omega$, i.e.,

\begin{equation}
(2.4) \quad 0 < \underline{b} \leq B(x) \leq \overline{b} \quad \text{for all} \quad x \in \Omega.
\end{equation}

We will assume that $A$ and $B$ are both piecewise constant on $\Omega$ and that any jumps in $A$ and $B$ are aligned with the meshes $T_n$ (introduced below), for all $n$.

Throughout the paper, for any polygonal (polyhedral) subdomain of $D \subset \Omega$, and any $s \in [0, 1]$, $\| \cdot \|_{s,D}$ and $| \cdot |_{s,D}$ will denote the standard norm and seminorm in the Sobolev space $H^s(D)$. Also $(\cdot, \cdot)_{0,D}$ denotes the $L^2(D)$ inner product. We also define the energy norm induced by the bilinear form $a$:

$$
\| u \|_{a,\Omega}^2 := a(u, u) \quad \text{for all} \quad u \in H^1_0(\Omega),
$$

which, by (2.3), is equivalent to the $H^1(\Omega)$ seminorm. (The equivalence constant depends on the contrast $\overline{a}/\underline{a}$, but we are not concerned with this dependence in the present paper.) We also introduce the weighted $L^2$ norm:

$$
\| u \|_{0,B,\Omega}^2 = b(u, u) = \int_{\Omega} B(x)|u(x)|^2 \, dx,
$$

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\[ a(u, v) = \lambda b(u, v), \quad \text{for all } v \in H^1_0(\Omega) \]

By the continuity of \( a \) and \( b \) and the coercivity of \( a \) on \( H^1_0(\Omega) \) it is a standard result that (2.6) has a countable sequence of nondecreasing positive eigenvalues \( \lambda_j \), \( j = 1, 2, \ldots \) with corresponding eigenfunctions \( u_j \in H^1_0(\Omega) \) [3, 12, 24].

In this paper we will need some additional regularity for the eigenfunctions \( u_j \), which will be achieved by making the following regularity assumption for the elliptic problem induced by \( a \).

**Assumption 2.1.** We assume that there exists a constant \( C_{\text{ell}} > 0 \) and \( s \in [0, 1] \) with the following property. For \( f \in L^2(\Omega) \), if \( v \in H^1_0(\Omega) \) solves the problem \( a(v, w) = (f, w)_{\Omega, \Omega} \) for all \( w \in H^1_0(\Omega) \), then \( ||v||_{1+s, \Omega} \leq C_{\text{ell}} ||f||_{0, \Omega} \).

Assumption 2.1 is satisfied with \( s = 1 \) when \( A \) is constant (or smooth) and \( \Omega \) is a smooth boundary or is a convex polygon. In a range of other practical cases \( s \in (0, 1) \), for example, \( \Omega \) nonconvex (see [4]), or \( A \) having a discontinuity across an interior interface (see [2]). Under Assumption 2.1 it follows that the eigenfunctions \( u_j \) of the problem (2.6) satisfy \( ||u_j||_{1+s, \Omega} \leq C_{\text{ell}} \lambda_j \sqrt{\pi} \).

To approximate problem (2.6) we use the piecewise linear finite element method. Accordingly, let \( T_n, n = 1, 2, \ldots \) denote a family of conforming triangular \((d = 2)\) or tetrahedral \((d = 3)\) meshes on \( \Omega \). Each mesh consists of elements denoted \( \tau \in T_n \). We assume that for each \( n \), \( T_{n+1} \) is a refinement of \( T_n \). For a typical element \( \tau \) of any mesh, its diameter is denoted \( H_\tau \) and the diameter of its largest inscribed ball is denoted \( \rho_\tau \). For each \( n \), let \( H_n \) denote the piecewise constant mesh function on \( \Omega \), whose value on each element \( \tau \in T_n \) is \( H_\tau \) and let \( H_n^{\text{max}} = \max_{\tau \in T_n} H_\tau \). Throughout we will assume that the family of meshes \( T_n \) is shape regular; i.e., there exists a constant \( C_{\text{reg}} \) such that

\[ H_\tau \leq C_{\text{reg}} \rho_\tau, \quad \text{for all } \tau \in T_n \quad \text{and all } \quad n = 1, 2, \ldots \]

In the later sections of the paper, the \( T_n \) will be produced by an adaptive process which ensures shape regularity.

We let \( V_n \) denote the usual finite dimensional subspace of \( H^1_0(\Omega) \), consisting of all continuous piecewise linear functions with respect to the mesh \( T_n \). Then the discrete formulation of problem (2.6) is to seek the eigenpairs \( (\lambda_n, u_n) \in \mathbb{R} \times V_n \) such that

\[ a(u_n, v_n) = \lambda_n b(u_n, v_n), \quad \text{for all } v_n \in V_n \]

The problem (2.8) has \( N = \dim V_n \) positive eigenvalues (counted according to multiplicity) which we denote in nondecreasing order as \( \lambda_{n,1} \leq \lambda_{n,2} \leq \cdots \leq \lambda_{n,N} \). It is well-known (see [24, section 6.3]) that for any \( j \), \( \lambda_{n,j} \to \lambda_j \) as \( H_n^{\text{max}} \to 0 \) and (by the minimax principle—see, e.g., [24, section 6.1]) the convergence of the \( \lambda_{n,j} \) is monotone decreasing, i.e.,

\[ \lambda_{n,j} \geq \lambda_{m,j} \geq \lambda_j, \quad \text{for all } \quad j = 1, \ldots, N, \quad \text{and all } m \geq n. \]
Thus, it is clear that there exists a separation constant \( \rho > 0 \) (depending on the spectrum of (2.6)) with the following property: If \( \lambda_j = \lambda_{j+1} = \cdots = \lambda_{j+R-1} \) is any eigenvalue of (2.6) of multiplicity \( R \geq 1 \), then

\[
\lambda_j \quad \frac{1}{|\lambda_{n,\ell} - \lambda_j|} \leq \rho, \quad \ell \neq j, j+1, \ldots, j+R-1,
\]

provided \( H_n^{\text{max}} \) is sufficiently small. (Note that for \( \ell \neq j, j+1, \ldots j+R-1 \), \( \lambda_n,\ell \rightarrow \lambda \neq \lambda_j \).

The a priori error analysis for our eigenvalue problem is classical (see, e.g., [3], [12], and [24]). In the next section, we briefly recall some of the main known results and also prove a nonclassical result (Theorem 3.2) which is essential to the proof of convergence of our adaptive scheme.

3. A priori analysis. In this section we shall assume that \( \lambda_j \) is an eigenvalue of (2.6) and \( \lambda_{n,j} \) is its approximation as described above. Let \( u_j \) and \( u_{n,j} \) be any corresponding normalized eigenvectors as defined in (2.6) and (2.8). From these we obtain the important basic identity:

\[
a(u_j - u_{n,j}, u_j - u_{n,j}) = a(u_j, u_j) + a(u_{n,j}, u_{n,j}) - 2a(u_j, u_{n,j})
= \lambda_j + \lambda_{n,j} - 2\lambda_j b(u_j, u_{n,j})
= \lambda_{n,j} - \lambda_j + \lambda_j (2 - 2b(u_j, u_{n,j}))
= \lambda_{n,j} - \lambda_j + \lambda_j b(u_j - u_{n,j}, u_j - u_{n,j}).
\]

Using this and (2.9), we obtain

\[
|||u_j - u_{n,j}|||_\Omega^2 = |\lambda_j - \lambda_{n,j}| + \lambda_j \|u_j - u_{n,j}\|_{L^2,\Omega}^2.
\]

The following theorem investigates the convergence of discrete eigenpairs. Although parts of it are very well-known, we do not know a suitable reference for all the results given below, so a brief proof is given for completeness. In the proof we make use of the orthogonal projection \( Q_n \) of \( H_0^1(\Omega) \) onto \( V_n \) with respect to the inner product induced by \( a(\cdot, \cdot) \), which has the property:

\[
a(Q_n u, v_n) = \lambda b(u, v_n) \quad \text{for all} \quad v_n \in V_n.
\]

In the main result of this paper we prove convergence for adaptive approximations to eigenvalues and eigenvectors assuming for simplicity a simple eigenvalue. The following preliminary theorem is stated for a simple eigenvalue. However, this result is known for multiple eigenvalues (see, e.g., [24]). More details are given in [10].

**Theorem 3.1.** Let \( \lambda_j \) be a simple eigenvalue of (2.6), let \( \lambda_{n,j} \) be its associated approximation from solving (2.8), and let \( u_j \) and \( u_{n,j} \) be any corresponding normalized eigenvectors. Then for all \( 1 \leq j \leq N \),

\[
(i) \quad |\lambda_j - \lambda_{n,j}| \leq |||u_j - u_{n,j}|||_\Omega^2;
\]

\[
(ii) \quad \text{There are constants } C_1, C_2 > 0 \text{ and scalars } \alpha_{n,j} \in \{\pm 1\} \text{ such that}
\]

\[
|||u_j - \alpha_{n,j} u_{n,j}|||_{L^2,\Omega} \leq C_1 (H_n^{\text{max}})^s |||u_j - Q_n u_j|||_\Omega
\]

\[
\leq C_1 (H_n^{\text{max}})^s |||u_j - \alpha_{n,j} u_{n,j}|||_\Omega,
\]

where \( s \) is as in Assumption 2.1.
(iii) For sufficiently small $H_{n}^{\text{max}}$ there is a constant $C_{2}$ such that

$$
\|u_{j} - \alpha_{n,j}u_{n,j}\|_{\Omega} \leq C_{2}(H_{n}^{\text{max}})^{s}.
$$

The constants $C_{1}, C_{2}$ depend on the spectral information $\lambda_{\ell}$, $u_{\ell}$, $\ell = 1, \ldots, j$, the separation constant $\rho$, the constants $C_{\text{cll}}, C_{\text{reg}}$ in Assumption 2.1 and in (2.7) and on the bounds $\|u_{n}\|_{B}, \|u_{j}\|_{B}$ in (2.3), (2.4).

Proof. The estimate (3.4) follows directly from (3.2). Note that (3.4) holds even if $u_{n,j}$ is not close to $u$, which may occur due to the nonuniqueness of the eigenvectors.

The proof of (3.5) is obtained by a reworking of the results in [24]. By the symmetry of $a$ and $b$ there exists a basis $\{u_{n,\ell} : \ell = 1, \ldots, N\}$ of $V_{n}$ (containing $u_{n,j}$) which is orthonormal with respect to inner product $b$, and each $u_{n,\ell}$ is an eigenvector of (2.8) corresponding to eigenvalue $\lambda_{n,\ell}$. Then with $\beta_{n,j} := b(Q_{n}u_{j}, u_{n,j})$, Parseval’s equality yields

$$
\|Q_{n}u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega}^{2} = \sum_{\ell \neq j}^{N} b(Q_{n}u_{j}, u_{n,\ell})^{2}.
$$

Then, since

$$
\lambda_{n,\ell} b(Q_{n}u_{j}, u_{n,\ell}) = a(Q_{n}u_{j}, u_{n,\ell}) = a(u_{j}, u_{n,\ell}) = \lambda_{j} b(u_{j}, u_{n,\ell}),
$$

we have $(\lambda_{n,\ell} - \lambda_{j}) b(Q_{n}u_{j}, u_{n,\ell}) = \lambda_{j} b(u_{j} - Q_{n}u_{j}, u_{n,\ell})$, and so

$$
\|Q_{n}u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega}^{2} = \sum_{\ell \neq j}^{N} \left( \frac{\lambda_{j}}{\lambda_{n,\ell} - \lambda_{j}} \right)^{2} b(u_{j} - Q_{n}u_{j}, u_{n,\ell})^{2}
$$

$$
\leq \rho^{2} \sum_{\ell \neq j}^{N} b(u_{j} - Q_{n}u_{j}, u_{n,\ell})^{2} \leq \rho^{2} \|u_{j} - Q_{n}u_{j}\|_{0,B,\Omega}^{2},
$$

with the last step again by Parseval’s equality. Hence,

$$
\|u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega} \leq (1 + \rho) \|u_{j} - Q_{n}u_{j}\|_{0,B,\Omega}.
$$

Moreover,

$$
\|u_{j}\|_{0,B,\Omega} - \|u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega} \leq \|\beta_{n,j}u_{n,j}\|_{0,B,\Omega} \leq \|u_{j}\|_{0,B,\Omega} + \|u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega}.
$$

Since the $u_{j}$ and the $u_{n,j}$ are normalized, this implies

$$
1 - \|u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega} \leq |\beta_{n,j}| \leq 1 + \|u_{j} - \beta_{n,j}u_{n,j}\|_{0,B,\Omega}
$$

and, combining these with (3.8), we have

$$
|\beta_{n,j} - 1| \leq (1 + \rho) \|u_{j} - Q_{n}u_{j}\|_{0,B,\Omega}.
$$

Thus, with $\alpha_{n,j} := \text{sign}(\beta_{n,j})$, we have $|\beta_{n,j} - \alpha_{n,j}| \leq (1 + \rho) \|u_{j} - Q_{n}u_{j}\|_{0,B,\Omega}$, and

$$
\|u_{j} - \alpha_{n,j}u_{n,j}\|_{0,B,\Omega} \leq 2(1 + \rho) \|u_{j} - Q_{n}u_{j}\|_{0,B,\Omega}.
$$
The first inequality in (3.5) now follows from an application of the standard Aubin–Nitsche duality argument, while the second is just the best approximation of $Q_n$ in the energy norm.

The proof of (3.6) is a slight modification of that given in [24, Theorem 6.2]. The argument consists of obtaining an $O((H^\text{max}_n)^{2s})$ estimate for the eigenvalue error $|\lambda_j - \lambda_{n,j}|$ and then combining this with (3.2) and (3.5). □

The next theorem is a generalization to eigenvalue problems of the standard monotone convergence property for linear symmetric elliptic PDEs, namely, that if one enriches the finite dimensional space, then the error is bound to decrease. This result fails to hold for eigenvalue problems (even for symmetric elliptic partial differential operators), because of the nonlinearity of such problems. The best that we can do is to show that if the finite dimensional space is enriched, then the error will not increase very much. This is the subject of Theorem 3.2.

**Theorem 3.2.** For any $1 \leq j \leq N$, there exists a constant $q > 1$ such that, for $m \geq n$, the corresponding computed eigenpair $(\lambda_{m,j}, u_{m,j})$ satisfies:

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{\Omega} \leq q \|u_j - \alpha_{n,j}u_{n,j}\|_{\Omega}.
\]

**Proof.** From Theorem 3.1 (ii), we obtain

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{0,\Omega} \leq C_1(H^\text{max}_m)^s \|u_j - Q_m u_j\|_{\Omega}.
\]

Since $T_m$ is a refinement of $T_n$, it follows that $V_n \subset V_m$ and so the best approximation property of $Q_m$ ensures that

\[
\|u_j - Q_m u_j\|_{\Omega} \leq \|u_j - Q_n u_j\|_{\Omega}.
\]

Hence, from (3.10) and using the fact that $H^\text{max}_m \leq H^\text{max}_n$, we have

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{0,\Omega} \leq C_1(H^\text{max}_n)^s \|u_j - Q_n u_j\|_{\Omega}.
\]

Recalling that (3.2) holds for all eigenfunctions, and using (3.11) and then (2.9), we obtain

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{\Omega}^2 \leq |\lambda_j - \lambda_{m,j}| + \lambda_j \|u_j - \alpha_{m,j}u_{m,j}\|_{0,\Omega}^2 \\
\leq |\lambda_j - \lambda_{m,j}| + \lambda_j C_1^2(H^\text{max}_n)^{2s} \|u_j - Q_n u_j\|_{\Omega}^2 \\
\leq |\lambda_j - \lambda_{n,j}| + \lambda_j C_1^2(H^\text{max}_n)^{2s} \|u_j - Q_n u_j\|_{\Omega}^2.
\]

Hence, from (3.4) we obtain

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{\Omega}^2 \leq \|u_j - \alpha_{n,j}u_{n,j}\|_{\Omega}^2 + \lambda_j C_1^2(H^\text{max}_n)^{2s} \|u_j - Q_n u_j\|_{\Omega}^2.
\]

But, since $Q_n$ yields the best approximation from $V_n$ in the energy norm, we have

\[
\|u_j - \alpha_{m,j}u_{m,j}\|_{\Omega}^2 \leq (1 + \lambda_j C_1^2(H^\text{max}_n)^{2s}) \|u_j - \alpha_{n,j}u_{n,j}\|_{\Omega}^2,
\]

which is in the required form. □

**Remark 3.3.** From now on we will be concerned with a true eigenpair $(\lambda_j, u_j)$ and its computed approximation $(\lambda_{j,n}, u_{j,n})$ on the mesh $T_n$. Theorem 3.1 tells us that a priori $\lambda_{n,j}$ is “close” to $\lambda_j$ and that the spaces spanned by $u_j$ and $u_{n,j}$ are close. From now on we drop the subscript $j$ and we simply write $(\lambda, u)$ for the eigenpair of (2.6) $(\lambda_n, u_n)$ for a corresponding eigenpair of (2.8) and the scalar $\alpha_{n,j}$ is abbreviated $\alpha_n$. 

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4. A posteriori analysis. This section contains our a posteriori error estimator and the definition of the adaptive algorithm for which convergence will be proved in the following sections.

Recalling the mesh sequence $T_n$ defined above, we let $S_n$ denote the set of all the interior edges (or the set of interior faces in 3D) of the elements of the mesh $T_n$. For each $S \in S_n$, we denote by $\tau_1(S)$ and $\tau_2(S)$ the elements sharing $S$ (i.e., $\tau_1(S) \cap \tau_2(S) = S$) and we write $\Omega(S) = \tau_1(S) \cup \tau_2(S)$. We let $\bar{n}_S$ denote the unit normal vector to $S$, orientated from $\tau_1(S)$ to $\tau_2(S)$. All elements, faces, and edges are considered to be closed sets. Furthermore, we denote the diameter of $S$ by $H_S$. Note that, by mesh regularity, $\text{diam}(\Omega(S)) \sim H_{\tau_i(S)}$, $i = 1, 2$.

**Notation 4.1.** We write $A \lesssim B$ when $A/B$ is bounded by a constant which may depend on the functions $A$ and $B$ in (2.2), on $\underline{a}$, $\overline{a}$, $\underline{b}$, and $\overline{b}$, on $C_{\text{cell}}$ in Assumption 2.1, $C_{\text{reg}}$ in (2.7). The notation $A \equiv B$ means $A \lesssim B$ and $A \gtrsim B$.

All the constants depending on the spectrum, namely, $\rho$ in (2.10), $q$ in (3.9), $C_1$ and $C_2$ in (3.5) and (3.6), are handled explicitly. Similarly all mesh size dependencies are explicit. Note that all eigenvalues of (2.8) satisfy $\lambda_n \gtrsim 1$, since $\lambda_n \geq \lambda_1 = a(u_1, u_1) \geq \|u_1\|_{0, \Omega}^2 \gtrsim \|u_1\|_{0, \Omega}^2 \geq \|u_1\|_{0, B, \Omega}^2 = 1$.

Our error estimator is obtained by adapting standard estimates for source problems to the eigenvalue problem. Analogous eigenvalue estimates can be found in [9] (for the Laplace problem) and [25] (for linear elasticity) and related results are in [14].

For a function $g$, which is piecewise continuous on the mesh $T_n$, we introduce its jump across an edge (face) $S \in S_n$ by:

$$[g]_S(x) := \left( \lim_{\tilde{x} \to x} g(\tilde{x}) - \lim_{\tilde{x} \to x} g(\tilde{x}) \right), \quad \text{for } x \in \text{int}(S).$$

Then for any function $v$ with piecewise continuous gradient on $T_n$ we define, for $S \in S_n$,

$$J_S(v)(x) := [\bar{n}_S \cdot \nabla v]_S(x), \quad \text{for } x \in \text{int}(S).$$

The error estimator $\eta_n$ on the mesh $T_n$ is defined as

$$\eta_n^2 := \sum_{S \in S_n} \eta_{S,n}^2,$$

where, for each $S \in S_n$,

$$\eta_{S,n}^2 := \|H_n \lambda_n u_n\|_{0, B, \Omega(S)}^2 + \left|\left|\left| H_n^{1/2} J_S(u_n) \right|\right|_{0, S}^2.\right.$$  \hspace{1cm} (4.2) (reliability).

The following lemma is proved, in a standard way, by adapting the usual arguments for linear source problems. Note again that $\lambda$ is an eigenvalue of (2.6), $\lambda_n$ is a nearby eigenvalue of (2.8), and $u$, $u_n$ are any corresponding normalized eigenfunctions which are only “near” in the sense of Theorem 3.1.

**Lemma 4.2** (reliability).

$$\|u - u_n\|_{\Omega} \lesssim \eta_n + G_n,$$

and

$$G_n := \frac{1}{2} (\lambda + \lambda_n) \frac{\|u - u_n\|_{0, B, \Omega}^2}{\|u - u_n\|_{\Omega}^2}.$$
Remark 4.3. Recalling Remark 3.3, $u_n$ in Lemma 4.2 is any normalized eigenvector of (2.8) corresponding to the simple eigenvalue $\lambda$; i.e., its sign is not unique. However, the error estimators $\eta_{S,n}$ are independent of the sign of $u_n$. This is not a contradiction: we shall see that only one choice of eigenfunction will guarantee that the second term on the right-hand side of (4.3) is small, and only in this case is the left-hand side also guaranteed to be small.

A similar result to Lemma 4.2 was proved in [25, Proposition 5].

Proof. To ease readability we set $e_n = u - u_n$ in the proof. Note first that, since $(\lambda, u)$ and $(\lambda_n, u_n)$, respectively, solve the eigenvalue problems (2.1) and (2.8), we have, for all $w \in V_n$,

$$\|w\|_{H^1(\Omega)}^2 = a(e_n, e_n) = a(e_n, e_n - w_n) + a(e_n, w_n)$$

$$= a(e_n, e_n - w_n) + a(u, w_n) - a(u_n, w_n)$$

$$= a(e_n, e_n - w_n) + b(\lambda u - \lambda_n u_n, w_n)$$

$$= a(e_n, e_n - w_n) - b(\lambda u - \lambda_n u_n, e_n - w_n) + b(\lambda u - \lambda_n u_n, e_n).$$

(4.5)

To estimate the first two terms on the right-hand side of (4.5), first note that, for all $v \in H^1_0(\Omega)$,

$$a(e_n, v) - b(\lambda u - \lambda_n u_n, v) = -a(u_n, v) + \lambda_n b(u_n, v).$$

Hence, using elementwise integration by parts (and the fact that $A \nabla u_n$ is constant on each element and $v$ vanishes on $\partial\Omega$), we obtain

$$a(e_n, v) - b(\lambda u - \lambda_n u_n, v) = -\sum_{\tau \in T_n} \int_{\tau} (A \nabla u_n) \cdot \nabla v + \lambda_n b(u_n, v)$$

(4.6)

$$= -\sum_{S \in S_n} \int_S J_S(u_n) v + \lambda_n b(u_n, v),$$

and hence, for all $w_n \in V_n$,

$$a(e_n, e_n - w_n) - b(\lambda u - \lambda_n u_n, e_n - w_n) = -\sum_{S \in S_n} \int_S J_S(u_n)(e_n - w_n) + \lambda_n b(u_n, e_n - w_n).$$

(4.7)

Now recall the Scott–Zhang quasi-interpolation operator ([22]) which has the property that, for all $v \in H^1_0(\Omega)$, $I_n v \in V_n$ and

$$\|v - I_n v\|_{0,\tau} \lesssim H_\tau |v|_{1,\omega(\tau)}, \quad \|v - I_n v\|_{0,S} \lesssim H^1_S |v|_{1,\omega(S)},$$

(4.8)

where $\omega(\tau)$ is the union of all elements sharing at least a point with $\tau$, and $\omega(S)$ is the union of all elements sharing at least a point with $S$. (Note $\Omega(S) \subseteq \omega(S)$.) Substituting $w_n = I_n e_n$ in (4.7) and using the Cauchy–Schwarz inequality, together with estimates (4.8), we obtain

$$a(e_n, e_n - w_n) - b(\lambda u - \lambda_n u_n, e_n - w_n) \lesssim \eta_n \|e_n\|_{\Omega}^2.$$  

(4.9)

To estimate the third term on the right-hand side of (4.5), we simply observe that due to the normalization in each of the eigenvalue problems (2.1) and (2.8) we have

$$b(\lambda u - \lambda_n u_n, e_n) = (\lambda + \lambda_n)(1 - b(u, u_n)) = \frac{1}{2} (\lambda + \lambda_n) \|e_n\|_{0,B,\Omega}^2.$$  

(4.10)
Now, combine (4.9) and (4.10) with (4.5) and divide by $|||e_n|||_{\Omega}$ to obtain the result. \[ \square \]

Remark 4.4. We shall see below that $G_n$ defined above constitutes a “higher order term”.

For mesh refinement based on the local contributions to $\eta_n$, we use the same marking strategy as in [8] and [18]. The idea is to refine a subset of the elements of $T_n$ whose side residuals sum up to a fixed proportion of the total residual $\eta_n$.

Definition 4.5 (marking strategy 1). Given a parameter $0 < \theta < 1$, the procedure is: mark the sides in a minimal subset $\hat{S}_n$ of $S_n$ such that

\[
\left( \sum_{S \in \hat{S}_n} \eta_{S,n}^2 \right)^{1/2} \geq \theta \eta_n.
\]

To compute $\hat{S}_n$, we compute all the “local residuals” $\eta_{S,n}$, then insert edges (faces) into $\hat{S}_n$ in order of nonincreasing magnitude of $\eta_{S,n}$, until (4.11) is satisfied. A minimal subset $\hat{S}_n$ may not be unique. After this is done, we construct another set $\hat{T}_n$, containing all the elements of $T_n$, which contain at least one edge (face) belonging to $\hat{S}_n$.

In order to prove our convergence theory, we require an additional marking strategy based on oscillations (Definition 4.7 below). This also appears in some theories of adaptivity for source problems, e.g., [8], [18], [16], [7], and [6]), but to our knowledge has not yet been used in connection with eigenvalue problems.

The concept of “oscillation” is just a measure of how well a function may be approximated by piecewise constants on a particular mesh. For any function $v \in L^2(\Omega)$, and any mesh $T_n$, we introduce its orthogonal projection $P_n v$ onto piecewise constants defined by

\[
(P_n v)|_{\tau} = \frac{1}{|\tau|} \int_{\tau} v_n, \quad \text{for all} \quad \tau \in T_n.
\]

Then we make the definition:

Definition 4.6 (oscillations). On a mesh $T_n$, we define

\[
\text{osc}(v, T_n) := \|H_n(v - P_n v)||_0, B, \Omega.
\]

Note that

\[
\text{osc}(v, T_n) = \left( \sum_{\tau \in T_n} H^2_{\tau}\|v - P_n v\|^2_{0, B, \tau} \right)^{1/2},
\]

and that (by standard approximation theory and the ellipticity of $a(\cdot, \cdot)$),

\[
\text{osc}(v, T_n) \lesssim (H_n^{\text{max}})^2||v||_{\Omega}, \quad \text{for all} \quad v \in H^1_0(\Omega).
\]

The second marking strategy (introduced below) aims to reduce the oscillations corresponding to a particular approximate eigenfunction $u_n$.

Definition 4.7 (marking strategy 2). Given a parameter $0 < \tilde{\theta} < 1$: mark the elements in a minimal subset $\tilde{T}_n$ of $T_n$ such that

\[
\text{osc}(u_n, T_n) \geq \tilde{\theta} \text{osc}(u_n, T_n).
\]
Fig. 4.1. The refinement procedure applied to an element of the mesh. In (a) the element before the refinement, in (b) after the three sides have been refined, and in (c) after the bisection of one of the three new segments.

Analogously to (4.11), we compute $\tilde{T}_n$ by inserting elements $\tau$ into $\tilde{T}_n$ according to nonincreasing order of their local contributions $H_\tau^2 \| (u_n - P_n u_n) \|_{0, B, \tau}^2$ until (4.15) is satisfied.

Our adaptive algorithm can then be stated:

**Algorithm 1** Converging algorithm

**Require:** $0 < \theta < 1$

**Require:** $0 < \tilde{\theta} < 1$

**Loop**

- Solve the Problem (2.8) for $(\lambda_n, u_n)$
- Mark the elements using the first marking strategy (Definition 4.5)
- Mark any additional unmarked elements using the second marking strategy (Definition 4.7)
- Refine the mesh $T_n$ and construct $T_{n+1}$

**End Loop**

In 2D at the $n$th iteration in Algorithm 1 each element in the set $\tilde{T}_n \cup \tilde{T}_n$ is refined using the algorithm illustrated in Figure 4.1. This consists of three recursive applications of the newest node algorithm [17] to each marked triangle, first creating two sons, then four grandsons, and finally bisecting two of the grandsons. This well-known algorithm is stated without name in [18, section 5.1]), is called “bisection5” in [7] and is called “full refinement” in [23]. This technique creates a new node in the middle of each marked side in $\tilde{S}_n$ and also a new node in the interior of each marked element. It follows from [17] that this algorithm yields shape regular conforming meshes in 2D.

In the 3D case we use a suitable refinement that creates a new node on each marked face in $S_n$ and a node in the interior of each marked element.

In [18] and [16] it has been shown for linear source problems that the reduction of the error, as the mesh is refined, is triggered by the decay of oscillations of the source on the sequence of constructed meshes. For the eigenvalue problem (2.1) the quantity $\lambda_n$ plays the role of data and in principle we have to ensure that oscillations of this quantity (or, more precisely, of its finite element approximation $\lambda_n u_n$) are sufficiently small. However, $\lambda_n u_n$ may change if the mesh changes and so the proof of error reduction for eigenvalue problems is not as simple as it is for linear source problems. This is the essence of the theoretical difficulty dealt with in this paper.

**5. Error reduction.** In this section we give the proof of error reduction for Algorithm 1. The proof has been inspired by the corresponding theory for source problems in [18]. However, the nonlinearity of the eigenvalue problem introduces new complications, and there are several lemmas before the main theorem (Theorem 5.6).

For the rest of the section let $(\lambda_n, u_n)$ be an approximate eigenpair on a mesh $T_n$, let...
THE n+1 be the mesh obtained by one iteration of Algorithm 1, and let \((\lambda_{n+1}, u_{n+1})\) be the corresponding eigenpair in the sense made precise in Remark 3.3.

The first lemma uses ideas from [18, Lemma 4.2] for the 2D case. The extension of this lemma to the 3D case is treated in Remark 5.2.

**Lemma 5.1.** Consider the

\[
\lambda_n \int_{\Omega(S)} B(P_n u_n) \Phi - \int_S J_S(u_n) \Phi_S = \|H_n \lambda_n P_n u_n\|_{0, B,\Omega(S)}^2 + \|H_S^{1/2} J_S(u_n)\|_{0, S}^2,
\]

and

\[
|||\Phi_S|||_{0, S}^2 \leq \|H_n \lambda_n P_n u_n\|_{0, B,\Omega(S)}^2 + \|H_S^{1/2} J_S(u_n)\|_{0, S}^2,
\]

where \(||v||_{\Omega(S)}^2 := \int_{\Omega(S)} v^T A \nabla v\).

**Proof.** Figure 5.1 illustrates two possible configurations of the domain \(\Omega(S)\).

We then define

\[
\Phi_S := \alpha_S \varphi_S + \beta_1 \varphi_1 + \beta_2 \varphi_2,
\]

where \(\varphi_S\) and \(\varphi_i\) are the nodal basis functions associated with the points \(x_S\) and \(x_i\) on \(T_{n+1}\), and \(\alpha_S\), \(\beta_i\) are defined by

\[
\alpha_S = \begin{cases} \frac{\|H_S^{1/2} J_S(u_n)\|_{0, S}^2}{\int_S J_S(u_n) \varphi_S} & \text{if } J_S(u_n) \neq 0, \\ 0 & \text{otherwise,} \end{cases}
\]

and

\[
\beta_i = \begin{cases} \frac{\|H_n \lambda_n P_n u_n\|_{0, B,\tau_i(S)}^2 - \alpha_S \int_{\tau_i(S)} B \lambda_n (P_n u_n) \varphi_S}{\int_{\tau_i(S)} B \lambda_n (P_n u_n) \varphi_i} & \text{if } P_n u_n|_{\tau_i(S)} \neq 0, \\ 0 & \text{otherwise,} \end{cases}
\]

for \(i = 1, 2\).
Note that $J_S(u_n)$ and $P_n u_n$ are constant on each element $\tau$. Using the fact that $\text{supp}(\varphi_i) = \tau_i(S)$, for $i = 1, 2$ we can easily see that the above formulae imply
\begin{equation}
\alpha_S \int_S J_S(u_n) \varphi_S = - \left\| H_S^{1/2} J_S(u_n) \right\|_{0,S},
\end{equation}
\begin{equation}
\int_{\Omega(S)} B \lambda_n(P_n u_n)(\alpha_S \varphi_S + \beta_1 \varphi_1 + \beta_2 \varphi_2) = \left\| H_n \lambda_n P_n u_n \right\|_{0,\beta,\Omega(S)}^2
\end{equation}
(and that these formulae remain true even if $J_S(u_n)$ or $P_n u_n|_{\tau_i(S)}$ vanish). Hence,
\begin{align*}
\lambda_n \int_{\Omega(S)} N(P_n u_n) \Phi_S - \int_S J_S(u_n) \Phi_S &= \lambda_n \int_{\Omega(S)} B(P_n u_n)(\alpha_S \varphi_S + \beta_1 \varphi_1 + \beta_2 \varphi_2) - \alpha_S \int_S J_S(u_n) \varphi_S
\end{align*}
and (5.1) follows immediately on using (5.6) and (5.7).

To proceed from here note that by the shape-regularity of the mesh and the standard inverse estimate,
\begin{equation}
\left\| \varphi_S \right\|_{0,\Omega(S)} \lesssim H_S^{-1} \left\| \varphi_S \right\|_{0,\Omega(S)}.
\end{equation}
Also, for all elements $\tau \in T_{n+1}$ with $\tau \subset \text{supp} \varphi_S$, there exists an affine map $\chi : \hat{\tau} \to \tau$, where $\hat{\tau}$ is the unit simplex in $\mathbb{R}^2$ and $\hat{\varphi}_S := \varphi_S \circ \chi$ is a nodal basis function on $\hat{\tau}$. The Jacobian $J_\chi$ of $\chi$ is constant and is proportional to the area of $\tau$. Hence,
\begin{equation}
\left\| \hat{\varphi}_S \right\|^2_{0,\hat{\tau}} = \int_{\tau} \left\| \varphi_S \right\|^2 = \int_{\hat{\tau}} \left\| \hat{\varphi}_S \right\|^2 J_\chi \sim \frac{1}{H_S^2},
\end{equation}
which ensures at $\left\| \varphi_S \right\|_{0,\Omega(S)} \lesssim 1$ and, similarly, $\left\| \varphi_i \right\|_{0,\Omega(S)} \lesssim 1$. Combining these with (5.3), we obtain
\begin{equation}
\left\| \Phi_S \right\|^2_{0,\Omega(S)} \lesssim |\alpha_S|^2 + |\beta_1|^2 + |\beta_2|^2.
\end{equation}

Now, note that by a simple change of variable, $\int_S \varphi_S$ is the integral over $[-H_S/2, H_S/2]$ of the one-dimensional hat function centered on 0 and so $\int_S \varphi_S \sim H_S$. Since $J_S(u_n)$ is constant on $S$, we have
\begin{equation}
|\alpha_S| \lesssim \frac{\left\| J_S(u_n) \right\|_{0,\Omega(S)}^2 H_S^{1/2}}{H_S} \lesssim \left\| J_S(u_n) \right\|_{0,\Omega(S)} \sim \left\| H_S^{1/2} J_S(u_n) \right\|_{0,\Omega(S)}.
\end{equation}
Also, since $P_n u_n$ is constant on each $\tau_i(S)$ and, since $\int_{\tau_i(S)} B \phi_i \sim H_{\tau_i(S)}^2$, we have
\begin{align*}
|\beta_i| &\lesssim \frac{\lambda_n (P_n u_n|_{\tau_i(S)}) \left\| H_{\tau_i(S)} \right\|^2_{0,\beta,\tau_i(S)} + |\alpha_S| H_{\tau_i(S)}^2}{H_{\tau_i(S)}^2} \\
&\lesssim \lambda_n (P_n u_n|_{\tau_i(S)}) \left\| H_{\tau_i(S)}^2 \right\|_{0,\beta,\tau_i(S)} + |\alpha_S| \sim \lambda_n \left\| H_n P_n u_n \right\|_{0,\beta,\tau_i(S)} + |\alpha_S|.
\end{align*}
This implies
\begin{equation}
|\beta_i|^2 \lesssim \lambda_n \left\| H_n P_n u_n \right\|^2_{0,\beta,\tau_i(S)} + |\alpha_S|^2 \lesssim \lambda_n \left\| H_n P_n u_n \right\|^2_{0,\beta,\tau_i(S)} + \left\| H_S^{1/2} J_S(u_n) \right\|^2_{0,\Omega(S)},
\end{equation}
and the proof is completed by combining (5.8) with (5.9) and (5.10).
Remark 5.2. To extend the results in Lemma 5.1 to the 3D case we need to use a refinement procedure for tetrahedra that creates a new node on each marked face in $\bar{S}_n$ and a node in the interior of each marked element. The proof in the 3D case is similar to the proof in the 2D case: for each couple of refined elements we define

$$\Phi_S := \alpha_S \varphi_S + \beta_1 \varphi_1 + \beta_2 \varphi_2,$$

where $\varphi_S$ is the nodal basis function associated to the new node on the shared face and $\varphi_i$ are the nodal basis functions associated to the new nodes in the interior of the elements. The coefficients $\alpha_S$, $\beta_1$, and $\beta_2$ can be chosen in the same way as in Lemma 5.1, and the rest of the proof proceeds similarly.

In the next lemma, we bound the local error estimator above by the local difference of two discrete solutions coming from consecutive meshes, plus higher order terms. This kind of result is called “discrete local efficiency” by many authors.

Recall that $T_{n+1}$ is the refinement of $T_n$ obtained by applying Algorithm 1.

**Lemma 5.3.** For any $S \in \bar{S}_n$, we have

$$\eta_{S,n}^2 \lesssim \|u_{n+1} - u_n\|_{\Omega(S)}^2 + \|H_n(\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n)\|_{0, B, \Omega(S)}^2$$

(5.11)

$$+ \|H_n \lambda_n (u_n - P_n u_n)\|_{0, B, \Omega(S)}^2.$$

**Proof.** Since the function $\Phi_S$ defined in Lemma 5.1 is in $V_{n+1}$ and $\text{supp}(\Phi_S) = \Omega(S)$, we have

$$a(u_{n+1} - u_n, \Phi_S) = a(u_{n+1}, \Phi_S) - a(u_n, \Phi_S) = \lambda_{n+1} \int_{\Omega(S)} B u_{n+1} \Phi_S - a(u_n, \Phi_S).$$

(5.12)

Now applying integration by parts to the last term on the right-hand side of (5.12), we obtain

$$a(u_{n+1} - u_n, \Phi_S) = \lambda_{n+1} \int_{\Omega(S)} B u_{n+1} \Phi_S - \int_S J_S(u_n) \Phi_S.$$

(5.13)

Rewriting (5.13) and combining with (5.1), we obtain

$$a(u_{n+1} - u_n, \Phi_S) - \int_{\Omega(S)} B(\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n) \Phi_S$$

$$= \lambda_n \int_{\Omega(S)} B(P_n u_n) \Phi_S - \int_S J_S(u_n) \Phi_S$$

$$= \|H_n \lambda_n P_n u_n\|_{0, B, \Omega(S)}^2 + \left\|H_n^{1/2} J_S(u_n)\right\|^2_{0, S}.$$

(5.14)

Rearranging this, and then applying the triangle and Cauchy–Schwarz inequalities, we obtain

$$\|H_n \lambda_n P_n u_n\|_{0, B, \Omega(S)}^2 + \left\|H_n^{1/2} J_S(u_n)\right\|^2_{0, S}$$

$$\leq \|a(u_{n+1} - u_n, \Phi_S)\| + \left\|\int_{\Omega(S)} B(\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n) \Phi_S\right\|$$

$$\leq \|u_{n+1} - u_n\|_{\Omega(S)} \|\Phi_S\|_{\Omega(S)} + \|\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n\|_{0, B, \Omega(S)} \|\Phi_S\|_{0, B, \Omega(S)}$$

$$\left(\|u_{n+1} - u_n\|_{\Omega(S)} + \|H_n(\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n)\|_{0, B, \Omega(S)}\right) \|\Phi_S\|_{\Omega(S)}.$$
In the final step of (5.15) we made use of the Poincaré inequality \(\|\Phi_S\|_{0,B,\Omega(S)} \lesssim H_S\|\Phi_S\|_{\Omega(S)}\) and also the shape-regularity of the meshes. In view of (5.2), we have

\[
\|H_n\lambda_n P_n u_n\|_{0,B,\Omega(S)}^2 + \left\| H_n^{1/2} J_S(u_n) \right\|_{0,S}^2 \\
\lesssim \left( \|u_{n+1} - u_n\|_{\Omega(S)} + \|H_n(\lambda_{n+1} u_{n+1} - \lambda_n P_n u_n)\|_{0,B,\Omega(S)} \right)^2
\]

(5.16)

Now, from the definition of \(\eta_{S,n}\) in (4.2), and the triangle inequality, we have

\[
\eta_{S,n}^2 \lesssim \|H_n\lambda_n P_n u_n\|_{0,B,\Omega(S)}^2 + \left\| H_n^{1/2} J_S(u_n) \right\|_{0,S}^2 + \|H_n\lambda_n(u_n - P_n u_n)\|_{0,B,\Omega(S)}^2.
\]

(5.17)

The required inequality (5.11) now follows from (5.16) and (5.17).

In the main result of this section, Theorem 5.6 below, we will be interested in achieving an error reduction result of the form \(\|u - \alpha_{n+1} u_{n+1}\|_{\Omega} \leq \rho\|u - \alpha_n u_n\|_{\Omega}\) for some \(\rho < 1\). Note that we need to introduce the scalar \(\alpha_n\) here to ensure nearness of the approximate eigenfunction to the true one.

To prove error reduction we exploit the identity

\[
\|u - \alpha_n u_n\|_{\Omega}^2 = \|u - \alpha_{n+1} u_{n+1} + \alpha_{n+1} u_{n+1} - \alpha_n u_n\|_{\Omega}^2 \\
= \|u - \alpha_{n+1} u_{n+1}\|_{\Omega}^2 + \|\alpha_{n+1} u_{n+1} - \alpha_n u_n\|_{\Omega}^2 \\
+ 2\alpha(u - \alpha_{n+1} u_{n+1}, \alpha_{n+1} u_{n+1} - \alpha_n u_n).
\]

(5.18)

In the case of source problems (e.g., [18, 19]), the \(\alpha_n\) is not needed and the last term on the right-hand side vanishes due to Galerkin orthogonality. However, this approach is not available to us in the eigenvalue problem. Therefore, a more technical approach is needed to bound the last two terms on the right-hand side of (5.18) from below. The main technical result is in the following lemma. Recall the convention in Notation 4.1.

**Lemma 5.4.** With \(u, u_n, \alpha_n\) as in Remark 3.3,

(5.19) \[\|\alpha_{n+1} u_{n+1} - \alpha_n u_n\|_{\Omega}^2 \gtrsim \theta^2 \|\alpha_{n+1} u_{n+1} - \alpha_n u_n\|_{\Omega}^2 - \text{osc}(\lambda_n u_n, T_n)^2 - L_n^2,\]

where \(\theta\) is defined in the marking strategy in Definition 4.5 and \(L_n\) satisfies the estimate:

(5.20) \[L_n \leq \hat{C}(H_n^{max})^q\|u - \alpha_n u_n\|_{\Omega},\]

where \(\hat{C}\) depends on \(\theta, \lambda, C_1, C_2\), and \(q\).

**Remark 5.5.** Note that the oscillation term in (5.19) is unaffected if we replace \(\alpha_n u_n\) by \(u_n\).

**Proof.** By Definition 4.5 and Lemma 5.3, we have

\[
\theta^2 \eta_n^2 \leq \sum_{S \in \mathcal{S}_n} \eta_{S,n}^2 \\
\lesssim \|\alpha_{n+1} u_{n+1} - \alpha_n u_n\|_{\Omega}^2 \\
+ \|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n P_n \alpha_n u_n)\|_{0,B,\Omega}^2 + \text{osc}(\lambda_n u_n T_n)^2.
\]
Hence, rearranging and making use of Lemma 4.2 and Remark 4.3, we have

\[
\|\alpha_{n+1} u_{n+1} - \alpha_n u_n\|_\Omega^2 \geq \theta^2 \eta_n^2 - \|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n P_n \alpha_n u_n)\|^2_{0,B,\Omega}
- \text{osc}(\lambda_n u_n T_n)^2
\geq \theta^2 \|u - \alpha_n u_n\|_\Omega^2 - \text{osc}(\lambda_n u_n T_n)^2
\geq \theta^2 \tilde{G}_n^2 - \|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n P_n \alpha_n u_n)\|^2_{0,B,\Omega},
\]

(5.21)

where \( \tilde{G}_n \) is the same as \( G_n \) in Lemma 4.2, but with \( u_n \) replaced by \( \alpha_n u_n \).

Note that (5.21) is of the required form (5.19) with

\[
L_n := \left( \theta^2 \tilde{G}_n^2 + \|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n P_n \alpha_n u_n)\|^2_{0,B,\Omega} \right)^{1/2}
\]

We now estimate the last two terms in (5.21) to obtain (5.20). To estimate \( \tilde{G}_n \), we use Theorem 3.1(ii) to obtain

\[
\tilde{G}_n \leq \frac{1}{2}(\lambda + \lambda_n) C_1^2 (H_n^{\text{max}})^{2s} \frac{\|u - Q_n u\|_\Omega^2}{\|u - \alpha_n u_n\||u|_\Omega}
\leq \frac{1}{2}(\lambda + \lambda_n) C_1^2 (H_n^{\text{max}})^{2s} \|u - \alpha_n u_n\|_\Omega.
\]

(5.22)

To estimate the last term in (5.21), we first use the triangle inequality to obtain

\[
\|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n P_n \alpha_n u_n)\|_{0,B,\Omega} \leq \|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n \alpha_n u_n)\|_{0,B,\Omega} + \text{osc}(\lambda_n u_n, T_n).
\]

(5.23)

For the first term on the right-hand side of (5.23), we have

\[
\|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n \alpha_n u_n)\|_{0,B,\Omega} \leq H_n^{\text{max}} (\|\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}\|_{0,B,\Omega} + \|\lambda u - \lambda_n \alpha_n u_n\|_{0,B,\Omega}).
\]

(5.24)

Then, recalling (2.6) and Theorem 3.1, we obtain

\[
\|\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}\|_{0,B,\Omega} \leq |\lambda - \lambda_{n+1}| \|u\|_{0,B,\Omega}
+ \lambda_{n+1} \|u - \alpha_{n+1} u_{n+1}\|_{0,B,\Omega}
\leq |\|u - \alpha_{n+1} u_{n+1}\|_\Omega^2
+ \lambda_{n+1} C_1 (H_n^{\text{max}})^{s}\|u - \alpha_{n+1} u_{n+1}\|_\Omega.
\]

(5.25)

Using Theorem 3.1 (iii) and then Theorem 3.2, this implies

\[
\|\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}\|_{0,B,\Omega} \leq (C_2 + \lambda_{n+1} C_1) (H_n^{\text{max}})^s\|u - \alpha_{n+1} u_{n+1}\|_\Omega
\leq q(C_2 + \lambda_{n+1} C_1) (H_n^{\text{max}})^s\|u - \alpha_n u_n\|_\Omega.
\]

(5.26)

An identical argument shows

\[
\|\lambda u - \lambda_n \alpha_n u_n\|_{0,B,\Omega} \leq (C_2 + \lambda_n C_1) (H_n^{\text{max}})^s\|u - \alpha_n u_n\|_\Omega.
\]

(5.27)

Combining (5.26) and (5.27) with (5.24), and using (2.9), we obtain

\[
\|H_n(\lambda_{n+1} \alpha_{n+1} u_{n+1} - \lambda_n \alpha_n u_n)\|_{0,B,\Omega} \leq (1 + q)(C_2 + \lambda_n C_1) (H_n^{\text{max}})^{s+1}\|u - \alpha_n u_n\|_\Omega.
\]

(5.28)
Now combining (5.28) with (5.21), (5.22), and (5.23) we obtain the result.

The next theorem contains the main result of this section. It shows that, provided we start with a “fine enough” mesh $T_n$, the mesh adaptivity algorithm will reduce the error in the energy norm.

**Theorem 5.6 (error reduction).** For each $\theta \in (0,1)$, there exists a sufficiently fine mesh threshold $H_n^{\max}$ and constants $\mu > 0$ and $\rho \in (0,1)$ (all of which may depend on $\theta$ and on the eigenvalue $\lambda$), with the following property. For any $\varepsilon > 0$ the inequality

\begin{equation}
\text{osc}(\lambda_n u_n, T_n) \leq \mu \varepsilon
\end{equation}

implies either $|||u - \alpha_n u_n|||_{\Omega} \leq \varepsilon$ or

\begin{equation}
|||u - \alpha_{n+1} u_{n+1}|||_{\Omega} \leq \rho |||u - \alpha_n u_n|||_{\Omega}.
\end{equation}

**Proof.** In view of (5.18) and remembering that $\alpha_{n+1} u_{n+1} - \alpha_n u_n \in V_{n+1}$ we have

\begin{equation}
|||u - \alpha_n u_n|||_{\Omega}^2 = |||u - \alpha_{n+1} u_{n+1}|||_{\Omega}^2 + 2a(u - \alpha_{n+1} u_{n+1}, \alpha_{n+1} u_{n+1} - \alpha_n u_n)
\end{equation}

(5.30)

Before proceeding further, recall that by the assumptions (2.3) and (2.4), and the Poincaré inequality, there exists a constant $C_P$ (depending on $A, B$ and $\Omega$) such that

\begin{equation}
||v||_{0, \theta, \Omega} \leq C_P |||v|||_{\Omega}, \quad \text{for all} \quad v \in H_0^1(\Omega).
\end{equation}

Now using Cauchy–Schwarz and then the Young inequality $2ab \leq \frac{1}{4C_p^2}a^2 + 4C_p^2b^2$ on the second term on the right-hand side of (5.30), we get

\begin{equation}
|||u - \alpha_n u_n|||_{\Omega}^2 - |||u - \alpha_{n+1} u_{n+1}|||_{\Omega}^2
\geq |||\alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{\Omega}^2 - 2\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}|||_{0, \theta, \Omega} \alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{0, \theta, \Omega}
\geq |||\alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{\Omega}^2 - \frac{1}{4C_p^2}|||\alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{0, \theta, \Omega}^2
- 4C_p^2||\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}|||_{0, \theta, \Omega}^2
\geq \frac{3}{4}|||\alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{\Omega}^2 - 4C_p^2||\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}|||_{0, \theta, \Omega}^2.
\end{equation}

Hence

\begin{equation}
|||u - \alpha_{n+1} u_{n+1}|||_{\Omega}^2 \leq |||u - \alpha_n u_n|||_{\Omega}^2 - \frac{3}{4}|||\alpha_{n+1} u_{n+1} - \alpha_n u_n|||_{\Omega}^2
+ 4C_p^2||\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}|||_{0, \theta, \Omega}^2.
\end{equation}

Applying Lemma 5.4, we see that there exist constants $C, \hat{C}$ such that

\begin{equation}
|||u - \alpha_{n+1} u_{n+1}|||_{\Omega}^2 \leq \left(1 - \frac{3}{4}C\theta^2 + \frac{3}{4}C\hat{C}^2(H_n^{\max})^2\right)|||u - \alpha_n u_n|||_{\Omega}^2
+ 4C_p^2||\lambda u - \lambda_{n+1} \alpha_{n+1} u_{n+1}|||_{0, \theta, \Omega}^2
+ \frac{3}{4}C \text{osc}(\lambda_n u_n, T_n)^2.
\end{equation}
Then, making use of (5.26) we have
\begin{equation}
|||u - \alpha_{n+1} u_{n+1}|||_\Omega^2 \leq \gamma_n |||u - \alpha_n u_n|||_\Omega^2 + \frac{3}{4} C \text{osc}(\lambda_n u_n, T_n)^2
\end{equation}
with
\begin{equation}
\gamma_n := \left[ 1 - \frac{3}{4} C \theta^2 + C'(H_n^{\text{max}})^{2s} \right],
\end{equation}
where $C'$ is another constant independent of $n$. Note that $H_n^{\text{max}}$ can be chosen sufficiently small so that $\gamma_m \leq \gamma$ for some $\gamma \in (0, 1)$ and all $m \geq n$.
Consider now the consequences of the inequality (5.29). If $|||u - \alpha_n u_n|||_\Omega > \varepsilon$, then (5.32) implies
\begin{equation}
|||u - \alpha_{n+1} u_{n+1}|||_\Omega^2 \leq \left[ \gamma + \frac{3}{4} C \mu^2 \right] |||u - \alpha_n u_n|||_\Omega^2.
\end{equation}
Now choose $\mu$ small enough so that
\begin{equation}
\rho := \left( \gamma + \frac{3}{4} C \mu^2 \right)^{1/2} < 1
\end{equation}
to complete the proof. \hfill \Box

6. Proof of convergence. The main result of this paper is Theorem 6.2 below, which proves convergence of the adaptive method and also demonstrates the decay of oscillations of the sequence of approximate eigenfunctions. Before proving this result we need a final lemma.

Lemma 6.1. There exists a constant $\hat{\rho} \in (0, 1)$ such that
\begin{equation}
\text{osc}(u_{n+1}, T_{n+1}) \leq \hat{\rho} \text{osc}(u_n, T_n) + (1 + q)(H_n^{\text{max}})^2 |||u - \alpha_n u_n|||_\Omega.
\end{equation}

Proof. First, recall that one of the key results in [18], namely, [18, Lemma 3.8], is the proof that the oscillations of any fixed function $v \in H_0^1(\Omega)$ are reduced by applying one refinement based on Marking Strategy 2 (Definition 4.7). Thus, we have (in view of Algorithm 1):
\begin{equation}
\text{osc}(u_n, T_{n+1}) \leq \hat{\rho} \text{osc}(u_n, T_n),
\end{equation}
where $0 < \hat{\rho} < 1$ is independent of $u_n$. Thus, a simple application of the triangle inequality combined with (6.2) yields
\begin{equation}
\text{osc}(u_{n+1}, T_{n+1}) \leq \text{osc}(u_n, T_{n+1}) + \text{osc}(\alpha_{n+1} u_{n+1} - \alpha_n u_n, T_{n+1})
\leq \hat{\rho} \text{osc}(u_n, T_n) + \text{osc}(\alpha_{n+1} u_{n+1} - \alpha_n u_n, T_{n+1}).
\end{equation}
(Recall, again, that $\text{osc}(u_n, T_n) = \text{osc}(\alpha_n u_n, T_n)$.) A further application of the triangle inequality and then (4.14) yields
\begin{equation}
\text{osc}(\alpha_{n+1} u_{n+1} - \alpha_n u_n, T_{n+1}) \leq \text{osc}(u - \alpha_{n+1} u_{n+1}, T_{n+1}) + \text{osc}(u - \alpha_n u_n, T_{n+1})
\leq (H_n^{\text{max}})^2 (|||u - \alpha_{n+1} u_{n+1}|||_\Omega + |||u - \alpha_n u_n|||_\Omega),
\end{equation}
and then combining (6.3) and (6.4) and applying Theorem 3.2 completes the proof.

**Theorem 6.2.** Provided the initial mesh $T_0$ is chosen so that $H_0^{\text{max}}$ is small enough, there exists a constant $p \in (0,1)$, such that the recursive application of Algorithm 1 yields a convergent sequence of approximate eigenvalues and eigenvectors, with the property:

(6.5) \[ \| u - \alpha_n u_n \|_\Omega \leq B_0 q^p n, \]

and

(6.6) \[ \lambda_n \text{osc}(u_n, T_n) \leq B_1 p^n, \]

where $B_0$ and $B_1$ are constants and $q$ is the constant defined in Theorem 3.2.

**Remark 6.3.** The initial mesh convergence threshold and the constants $B_0$ and $B_1$ may depend on $\theta$, $\tilde{\theta}$, and $\lambda$.

**Proof.** The proof of this theorem is by induction and the induction step contains an application of Theorem 5.6. In order to ensure the reduction of the error, we have to assume that the starting mesh $T_0$ is fine enough and $\mu$ in Theorem 5.6 is small enough such that, for the chosen value of $\theta$, the quantity $\rho$ in (5.34) satisfies $\rho < 1$.

Then with $\tilde{\rho}$ as in Lemma 6.1, choose $p$ in the range

$$\max\{\rho, \tilde{\rho}\} < p < 1.$$  

We also set

$$B_1 = \text{osc}(\lambda_0 u_0, T_0) \quad \text{and} \quad B_0 = \max\{\mu^{-1} p^{-1} B_1, \| u - \alpha_0 u_0 \|_\Omega \}.$$  

To perform the inductive proof, first note that by the definition of $B_0$ and Theorem 3.2,

$$\| u - \alpha_0 u_0 \|_\Omega \leq B_0 \leq B_0 q,$$

since $q > 1$. Combined with the definition of $B_1$ we have shown the result for $n = 0$.

Now, suppose that, for some $n > 0$, the inequalities (6.5) and (6.6) hold.

Now let us consider the outcomes, depending on whether the inequality

(6.7) \[ \| u - \alpha_n u_n \|_\Omega \leq B_0 p^{n+1} \]

holds or not. If (6.7) holds, then we can apply Theorem 3.2 to conclude that

$$\| u - \alpha_{n+1} u_{n+1} \|_\Omega \leq q \| u - \alpha_n u_n \|_\Omega \leq q B_0 p^{n+1},$$

which proves (6.5) for $n + 1$.

On the other hand, if (6.7) does not hold, then, by definition of $B_0$,

(6.8) \[ \| u - \alpha_n u_n \|_\Omega > B_0 p^{n+1} \geq \mu^{-1} B_1 p^n. \]
Also, since we have assumed (6.6) for \( n \), we have

\[(6.9) \quad \lambda_n \text{osc}(u_n, T_n) \leq \mu \varepsilon \quad \text{with} \quad \varepsilon := \mu^{-1}B_1p^n.\]

Then (6.8) and (6.9) combined with Theorem 5.6 yields

\[|||u - \alpha_{n+1}u_{n+1}|||_\Omega \leq \rho|||u - \alpha_n u|||_\Omega,\]

and so, using the inductive hypothesis (6.5) combined with the definition of \( p \), we have

\[|||u - \alpha_{n+1}u_{n+1}|||_\Omega \leq \rho B_0 q p^n \leq q B_0 p^{n+1},\]

which, again, proves (6.5) for \( n + 1 \).

To conclude the proof, we have to show that also (6.6) holds for \( n + 1 \). Using Lemma 6.1, (2.9), and the inductive hypothesis, we have

\[(6.10) \quad \lambda_{n+1} \text{osc}(u_{n+1}, T_{n+1}) \leq \hat{\rho} B_1p^n + (1 + q)(H_n^{\text{max}})^2 \lambda_n B_0 q p^n \leq (\hat{\rho} B_1 + (1 + q)(H_0^{\text{max}})^2 \lambda_0 B_0 q) p^n.\]

Now, (recalling that \( \hat{\rho} < \rho \)), in addition to the condition already imposed on \( H_0^{\text{max}} \), we can further require that

\[\hat{\rho} B_1 + (1 + q)(H_0^{\text{max}})^2 \lambda_0 B_0 q \leq p B_1.\]

This ensures that

\[\lambda_{n+1} \text{osc}(u_{n+1}, T_{n+1}) \leq B_1 p^{n+1},\]

thus concluding the proof. \( \Box \)

7. Numerical experiments. We present numerical experiments to illustrate the convergence theory. Algorithm 1 has been implemented in FORTRAN95. The mesh refinement has been done using the toolbox ALBERTA [20]. We used the package ARPACK [15] to compute eigenpairs and the sparse direct linear solver ME27 from the HSL [21, 13] to carry out the shift-invert solves required by ARPACK. Additional numerical experiments on photonic crystal problems and on 3D problems are given in [10] and [11].

7.1. Example: Laplace operator. In the first set of simulations, we have solved the Laplace eigenvalue problem (i.e., \( A = I \) and \( B = 1 \) in (2.2)) on a unit square with Dirichlet boundary conditions. The exact eigenvalues are known explicitly.

We compare different runs of Algorithm 1 using different values for \( \theta \) and \( \tilde{\theta} \) in Table 7.1. Since the problem is smooth, from Theorem 3.1 it follows that using uniform refinement the rate of convergence for eigenvalues should be \( O(H_n^{\text{max}})^2 \), or, equivalently, the rate of convergence in the number of degrees of freedom (DOFs) \( N \) should be \( O(N^{-1}) \). We measure the rate of convergence by conjecturing that \( |\lambda - \lambda_n| = C N^{-\beta} \) and estimating \( \beta \) for each pair of computations from the formula \( \beta = -\log(|\lambda - \lambda_n|/|\lambda - \lambda_{n-1}|)/\log(\text{DOFs}_n/\text{DOFs}_{n-1}) \). Similarly, Table 7.2 contains the same kind of information relative to the fourth smallest eigenvalue of the problem. Our results show a convergence rate close to \( O(N^{-1}) \) for \( \theta, \tilde{\theta} \) sufficiently large. However, the rate of convergence is sensitive to the values of \( \theta \) and \( \tilde{\theta} \).
In the theory presented in [24], it is shown that the error for eigenvalues for smooth problems is bounded in terms of the square of the considered eigenvalue, i.e.,

\[ |\lambda - \lambda_n| \leq C \lambda^2 (H_{\text{max}})^2. \]

Also, we know that the first and the fourth eigenvalues are 19.7392089 and 78.9568352, and so, \( \lambda_4 = 4\lambda_1 \). Comparing errors in Table 7.2 with those in Table 7.1, we see that the errors are roughly multiplied by a factor of 16, as predicted by (7.1).

Often h-adaptivity uses only a marking strategy based on an estimation of the error, as in Marking Strategy 1 and avoids refining based on oscillations as in Marking Strategy 2. (Convergence of an adaptive scheme for eigenvalue problems which does not use marking strategy 2 is recently proved in [5].) To investigate the effects of refinement based on oscillations, in Table 7.3 we have computed the smallest eigenvalue for the Laplace problem keeping \( \theta \) fixed and varying \( \tilde{\theta} \). Reducing \( \tilde{\theta} \) towards 0 has the effect of turning off the refinement arising from Marking Strategy 2. The results in Table 7.3 seem to suggest that the rate of convergence slightly increases as \( \tilde{\theta} \) increases.

We investigate this further in Table 7.4, where we take iterations 5, 6, and 7 from Table 7.3, and we present the quantity \( C^* := N \times |\lambda - \lambda_n| \), where \( N \) denotes the number of DOFs. Then \( C^* \) gives an indication of the size of the unknown constant in the optimal error estimate \( |\lambda - \lambda_n| = O(N^{-1}) \). The results suggest that \( C^* \) stays fairly constant independent of \( \theta \).

In Table 7.5, we have set \( \tilde{\theta} = 0 \). Although the convergence result given in this paper does not hold any more, the method is still clearly convergent. Comparing Table 7.1, Table 7.3, and Table 7.5, we see that with the second marking strategy the
Convergent Adaptive Method for Eigenvalue Problems

Comparison of the reduction of the error and DOFs of the adaptive method for the smallest eigenvalue for the Laplace problem on the unit square for a fixed value of $\theta$ and varying $\hat{\theta}$.

<table>
<thead>
<tr>
<th>Iteration</th>
<th>$\theta = 0.8, \hat{\theta} = 0.1$</th>
<th>$\theta = 0.8, \hat{\theta} = 0.3$</th>
<th>$\theta = 0.8, \hat{\theta} = 0.5$</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$</td>
<td>\lambda - \lambda_n</td>
<td>$</td>
</tr>
<tr>
<td>1</td>
<td>0.1340</td>
<td>400</td>
<td>0.1340</td>
</tr>
<tr>
<td>2</td>
<td>0.0608</td>
<td>2000</td>
<td>0.0608</td>
</tr>
<tr>
<td>3</td>
<td>0.0370</td>
<td>1260</td>
<td>0.0370</td>
</tr>
<tr>
<td>4</td>
<td>0.0133</td>
<td>7493</td>
<td>0.0133</td>
</tr>
<tr>
<td>5</td>
<td>0.0056</td>
<td>18853</td>
<td>0.0056</td>
</tr>
<tr>
<td>6</td>
<td>0.0021</td>
<td>52427</td>
<td>0.0021</td>
</tr>
<tr>
<td>7</td>
<td>0.0008</td>
<td>140049</td>
<td>0.0008</td>
</tr>
</tbody>
</table>

Table 7.4

Values of $C^*$ computed from Table 7.3.

<table>
<thead>
<tr>
<th>Iteration</th>
<th>$\theta = 0.8, \theta = 0.1$</th>
<th>$\theta = 0.8, \theta = 0.3$</th>
<th>$\theta = 0.8, \theta = 0.5$</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$</td>
<td>\lambda - \lambda_n</td>
<td>$</td>
</tr>
<tr>
<td>5</td>
<td>1.06 \times 10^2</td>
<td>1.05 \times 10^2</td>
<td>1.05 \times 10^2</td>
</tr>
<tr>
<td>6</td>
<td>1.10 \times 10^2</td>
<td>1.11 \times 10^2</td>
<td>1.11 \times 10^2</td>
</tr>
<tr>
<td>7</td>
<td>1.12 \times 10^2</td>
<td>1.12 \times 10^2</td>
<td>1.13 \times 10^2</td>
</tr>
</tbody>
</table>

Comparison of the reduction of the error and DOFs of the adaptive method for the smallest eigenvalue for the Laplace problem on the unit square using marking strategy 1 only.

<table>
<thead>
<tr>
<th>Iteration</th>
<th>$\theta = 0.2$</th>
<th>$\theta = 0.5$</th>
<th>$\theta = 0.8$</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$</td>
<td>\lambda - \lambda_n</td>
<td>$</td>
</tr>
<tr>
<td>1</td>
<td>0.1340</td>
<td>400</td>
<td>0.1340</td>
</tr>
<tr>
<td>2</td>
<td>0.1299</td>
<td>503</td>
<td>0.1299</td>
</tr>
<tr>
<td>3</td>
<td>0.1247</td>
<td>556</td>
<td>0.1247</td>
</tr>
<tr>
<td>4</td>
<td>0.1238</td>
<td>637</td>
<td>0.1238</td>
</tr>
<tr>
<td>5</td>
<td>0.1189</td>
<td>712</td>
<td>0.1189</td>
</tr>
<tr>
<td>7</td>
<td>0.1113</td>
<td>795</td>
<td>0.1113</td>
</tr>
</tbody>
</table>

Comparison between the number of marked elements by strategy 1 (i.e., $\bar{\mathcal{T}}_n$) and the number of marked elements by strategy 2 only (i.e., $\bar{\mathcal{T}}_n \setminus \mathcal{T}_n$) for different values of $\theta$ and $\hat{\theta}$ for the smallest eigenvalue of the Laplace problem on the unit square.

<table>
<thead>
<tr>
<th>Iteration</th>
<th>$\theta = \hat{\theta} = 0.2$</th>
<th>$\theta = \hat{\theta} = 0.5$</th>
<th>$\theta = \hat{\theta} = 0.8$</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$# \mathcal{T}_n$</td>
<td>$# (\mathcal{T}_n \setminus \mathcal{T}_n)$</td>
<td>$# \mathcal{T}_n$</td>
</tr>
<tr>
<td>1</td>
<td>12</td>
<td>15</td>
<td>85</td>
</tr>
<tr>
<td>2</td>
<td>13</td>
<td>15</td>
<td>102</td>
</tr>
<tr>
<td>3</td>
<td>14</td>
<td>15</td>
<td>100</td>
</tr>
<tr>
<td>4</td>
<td>14</td>
<td>14</td>
<td>115</td>
</tr>
<tr>
<td>5</td>
<td>15</td>
<td>13</td>
<td>310</td>
</tr>
<tr>
<td>6</td>
<td>15</td>
<td>12</td>
<td>552</td>
</tr>
</tbody>
</table>

The number of degrees of freedom grows faster than without it. To illustrate this effect better, Table 7.6 tabulates the number of elements $\bar{\mathcal{T}}_n$ (marked by Marking Strategy 1) with the extra number of elements $\# (\mathcal{T}_n \setminus \bar{\mathcal{T}}_n)$ (marked by Marking Strategy 2 alone). Note that the new DOFs created by mesh refinement come only from the refinement of
The marked elements, but also from the closures used to keep the meshes conforming. It is clear that the number of elements marked as a result of the oscillations continues to rise as refinement proceeds, although much more slowly than the number marked by the residual-based criterion (Marking Strategy 1).

In Figure 7.1 we compare the performance of the adaptive algorithm with uniform bisection refinement (see Figure 4.1) for the first and fourth eigenvalues of the Laplace operator. We note that in this case both methods converge with a similar rate, as is expected since in this case the regularity of eigenfunctions is $H^2$. To complete this section, we give in Table 7.7 an example of the performance of the adaptive method for computing nonsimple eigenvalues. In this case, we considered the second smallest eigenvalue of the Laplace operator on the unit square which has multiplicity 2. We see that, although the theory given above does not strictly hold, the method performs similarly to the case of simple eigenvalues.

### 7.2. Example: Elliptic operator with discontinuous coefficients.

In this example, we investigate how our method copes with discontinuous coefficients. In order to do that, we modified the smooth problem from Example 7.1. We inserted a square subdomain of side 0.5 in the center of the unit square domain. In the bilinear form (2.2), we also chose the function $A$ to be the scalar piecewise constant function, which assumes the value 100 inside the inner subdomain and the value 1 outside it. As before, $B$ in (2.2) is chosen as $B = 1$. The jump in the value of $A$ generally
Comparison of the reduction of the error and DOFs of the adaptive method for the smallest eigenvalue for the 2D problem with discontinuous coefficient.

<table>
<thead>
<tr>
<th>Iteration</th>
<th>$\theta = \tilde{\theta} = 0.2$</th>
<th>$\theta = \tilde{\theta} = 0.5$</th>
<th>$\theta = \tilde{\theta} = 0.8$</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>$</td>
<td>\lambda - \lambda_0</td>
<td>$, DOFs</td>
</tr>
<tr>
<td>1</td>
<td>1.1071, 81</td>
<td>1.1071, 81</td>
<td>1.1071, 81</td>
</tr>
<tr>
<td>2</td>
<td>0.8200, 102</td>
<td>0.3410, 199</td>
<td>0.5834, 387</td>
</tr>
<tr>
<td>3</td>
<td>0.0145, 129</td>
<td>0.0416, 314</td>
<td>0.2244, 799</td>
</tr>
<tr>
<td>4</td>
<td>0.8393, 147</td>
<td>0.0498, 491</td>
<td>0.5891, 703</td>
</tr>
<tr>
<td>5</td>
<td>0.8968, 167</td>
<td>0.8843, 673</td>
<td>1.1564, 4764</td>
</tr>
<tr>
<td>6</td>
<td>0.8076, 194</td>
<td>0.6996, 976</td>
<td>0.9665, 2235</td>
</tr>
<tr>
<td>7</td>
<td>0.8008, 217</td>
<td>0.0747, 1476</td>
<td>0.8722, 29148</td>
</tr>
<tr>
<td>8</td>
<td>0.7502, 237</td>
<td>0.7401, 2080</td>
<td>1.0237, 65387</td>
</tr>
</tbody>
</table>

The adaptive method yields the optimal order of $O(N^{-1})$ (which holds for uniform meshes and smooth problems) for large enough $\theta$ and $\tilde{\theta}$. (See Table 7.8.) Here we compute the “exact” $\lambda$ using a mesh with about half a million of DOFs.

In Figure 7.2, we depict the mesh coming from the fourth iteration of Algorithm 1 with $\theta = \tilde{\theta} = 0.8$ for the smallest eigenvalue of this problem. This mesh is the result of multiple refinements using both marking strategies 1 and 2 each time. As can be seen, the corners of the subdomain are much more refined than the rest of the mesh. This is clearly the effect of the first marking strategy, since the edge residuals have detected the singularity in the gradient of the eigenfunction at these points. In Figure 7.2, we also depict the corresponding eigenfunction.

In Figure 7.3, analogously to Figure 7.1, we compare the convergence of the adaptive method with uniform refinement for this example. Now, because of the lack of regularity, the superiority of the adaptive method is clearly visible.
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REFERENCES


