# A Posteriori Estimation of Dimension Reduction Errors for Elliptic Problems on Thin Domains 

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#### Abstract

A new a-posteriori error estimator is presented for the verification of the dimensionally reduced models stemming from the elliptic problems on thin domains. The original problem is considered in a general setting, without any specific assumptions on the domain geometry, coefficients and the right-hand sides. For the energy norm of the error of the zero-order dimension reduction method, the proposed estimator is shown to always provide a guaranteed upper bound. In the case when the original domain has constant thickness (but, possibly, non-plane upper and lower faces), the estimator demonstrates the optimal convergence rate as the thickness tends to zero. It is also flexible enough to successfully cope with the case of infinitely growing right-hand side of the equation when the domain thickness tends to zero. The numerical tests indicate the efficiency of the estimator and its ability to accurately represent the local error distribution needed for an adaptive improvement of the reduced model.


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## 1 Introduction

The method of dimension reduction is a popular approach frequently used by engineers for the approximate solution of the problems posed in thin domains. The term "thin" means that the size of the original physical domain along one coordinate direction is much smaller than along the others; this allows to make some simplifying assumptions on the behaviour of the exact solution and to replace the original high-dimensional problem by a lower-dimensional one. For instance, such a situation arises if a 3D problem is analysed with the help of a 2 D model. It is, however, clear that the solution of the new, "reduced" problem will, in general, differ from the solution to the original high-dimensional problem. Thus, the dimension reduction method unavoidably produces an error that can be referred to as the dimension reduction or the modelling error. The essential part of the model verification is, hence, a reliable a-posteriori control of the dimension reduction error.

Despite the practical importance of the topic, only a few a-posteriori estimators for the dimension reduction error have been introduced so far. In [12] and [3] (see also [2]) the residual-type estimators were proposed and proved reliable and efficient under the assumptions that the right-hand side of

[^0]the given equation is zero and the original domain is a plate with plane parallel faces. In [4] and [10] the implicit estimators based on the solution of local three-dimensional Neumann problems were developed for the hierarchical modelling of complex elastic plates. In [1] the estimator of Babuška and Schwab (see [2], [3]) was extended to take into account the discretization error stemming from the approximate solution of the reduced problem. In this respect, we have to notice that the present work is focused on the estimation of the modelling error, i.e. we assume, exactly as in [2], [3], that the error of discretizing the reduced problem is negligible. The work on the simultaneous a-posteriori estimation of both the modelling and the discretization error will be reported in a forthcoming paper.

In this work we propose a reliable and efficient a posteriori estimator for the dimension reduction error in the energy norm, having no specific assumptions on the right-hand side of the given equation and considering a general geometry of the given domain. In contrast to the above mentioned papers, which deal with the hierarchical modelling of the problems in thin domains, we only consider the so-called zero-order method of dimension reduction that is, however, very popular owing to its simplicity and purely two-dimensional formulation. At the same time, this method forms a basis for the hierarchical modelling of three-dimensional plates (see, e.g., [11], [3], [10]). We advocate the functional-type a posteriori error estimation approach (see [5], [6], [7], [8]) that essentially differs from the approaches taken in the aforementioned articles; however, surprisingly enough, it is possible to show that Babuška and Schwab's estimator for the zero-order reduced problem can be obtained as a particular case of our estimator when the right-hand side of the equation is zero and the original domain is a plate with plane parallel faces. It must be also noticed that the treatment of the case with non-zero right-hand side may require a special care, as we are about to see in one of the numerical examples; the presented estimator exhibits sufficient flexibility to remain efficient in this case.

The paper is set out as follows. Section 2 contains the geometric definitions and the problem statement. In Section 3 we derive the reduced problem. Section 4 is devoted to the derivation of the a-posteriori error estimate, while in Section 5 we consider two particular cases and analyse the behaviour of the estimator. The numerical examples are considered in Section 6, and we draw the conclusions in Section 7.

## 2 Problem setting

We consider three-dimensional Lipschitz domains which can be given in the form

$$
\Omega:=\left\{x \in \mathbb{R}^{3} \mid\left(x_{1}, x_{2}\right) \in \widehat{\Omega}, d_{\ominus}\left(x_{1}, x_{2}\right)<x_{3}<d_{\oplus}\left(x_{1}, x_{2}\right)\right\},
$$

where $\widehat{\Omega} \subset \mathbb{R}^{2}$ is the orthogonal projection of $\Omega$ on the ( $x_{1}, x_{2}$ )-plane ( $\widehat{\Omega}$ has the Lipschitz boundary $\widehat{\Gamma})$ and $d_{\ominus}$ and $d_{\oplus}$ are Lipschitz continuous functions defined on $\overline{\widehat{\Omega}}$. The lower and upper faces of $\Omega$ are denoted by

$$
\Gamma_{\ominus}:=\left\{x \in \mathbb{R}^{3} \mid\left(x_{1}, x_{2}\right) \in \widehat{\Omega}, x_{3}=d_{\ominus}\left(x_{1}, x_{2}\right)\right\}
$$

and

$$
\Gamma_{\oplus}:=\left\{x \in \mathbb{R}^{3} \mid\left(x_{1}, x_{2}\right) \in \widehat{\Omega}, x_{3}=d_{\oplus}\left(x_{1}, x_{2}\right)\right\},
$$

the lateral boundary by

$$
\Gamma_{0}:=\left\{x \in \mathbb{R}^{3} \mid\left(x_{1}, x_{2}\right) \in \widehat{\Gamma}, d_{\ominus}\left(x_{1}, x_{2}\right)<x_{3}<d_{\oplus}\left(x_{1}, x_{2}\right)\right\}
$$

(see Figure 1).
Remark 2.1 We consider $d_{\ominus}$ and $d_{\oplus}$ as explicit functions of ( $x_{1}, x_{2}$ )-coordinates only for the sake of simplicity. The generalization of the theory to the case of an arbitrary Lipschitz domain $\Omega$ presents no difficulty from the conceptional point of view.


Figure 1: Sketch of the domain geometry.

The assumption that the given domain $\Omega$ is "thin" can now be written as

$$
\begin{equation*}
\operatorname{diam} \widehat{\Omega} \gg \max _{\left(x_{1}, x_{2}\right) \in \bar{\Omega}} d\left(x_{1}, x_{2}\right), \tag{2.1}
\end{equation*}
$$

where $d=d_{\oplus}-d_{\ominus}$ is the domain thickness, $d\left(x_{1}, x_{2}\right) \geq d_{*}>0 \quad \forall\left(x_{1}, x_{2}\right) \in \overline{\widehat{\Omega}}$. Although the assumption is of purely qualitative nature, it will motivate the derivation of the corresponding twodimensional reduced model in the next section. We also have to notice that Figure 1 depicts a simplified case; in the geometrical definitions we do not assume the domain thickness $d\left(x_{1}, x_{2}\right)$ to be a constant.

In the domain $\Omega$ we consider a model elliptic problem

$$
\begin{align*}
-\operatorname{Div}(\mathbf{A} \nabla u) & =f \quad \text { in } \Omega,  \tag{2.2}\\
u & =0 \quad \text { on } \Gamma_{0},  \tag{2.3}\\
\mathbf{A} \nabla u \cdot \boldsymbol{\nu}_{\ominus} & =F_{\ominus} \quad \text { on } \Gamma_{\ominus},  \tag{2.4}\\
\mathbf{A} \nabla u \cdot \boldsymbol{\nu}_{\oplus} & =F_{\oplus} \quad \text { on } \Gamma_{\oplus}, \tag{2.5}
\end{align*}
$$

where $f \in L_{2}(\Omega), F_{\ominus} \in L_{2}\left(\Gamma_{\ominus}\right), F_{\oplus} \in L_{2}\left(\Gamma_{\oplus}\right), \boldsymbol{\nu}_{\ominus}$ and $\boldsymbol{\nu}_{\oplus}$ are outward normal vectors at $\Gamma_{\ominus}$ and $\Gamma_{\oplus}$ respectively. The matrix $\mathbf{A}=\left(a_{i j}(x)\right)_{i, j=\overline{1,3}}$ with the components from $L_{\infty}(\Omega)$ is symmetric and uniformly positive definite, i.e. there exist constants $0<c<C<\infty$ such that

$$
\begin{equation*}
c|\xi|^{2} \leq \mathbf{A}(x) \xi \cdot \xi \leq C|\xi|^{2} \quad \forall \xi \in \mathbb{R}^{3}, \text { a. e. in } \Omega . \tag{2.6}
\end{equation*}
$$

If the space of admissible functions is denoted by

$$
\begin{equation*}
V_{0}:=\left\{v \in H^{1}(\Omega) \mid v=0 \text { on } \Gamma_{0}\right\}, \tag{2.7}
\end{equation*}
$$

the weak form of the problem (2.2)-(2.5) reads
Problem (P): Find $u \in V_{0}$ such that

$$
\begin{equation*}
\int_{\Omega} \mathbf{A} \nabla u \cdot \nabla w d x=\int_{\Omega} f w d x+\int_{\Gamma_{\ominus}} F_{\ominus} w d s+\int_{\Gamma_{\oplus}} F_{\oplus} w d s \quad \forall w \in V_{0} . \tag{2.8}
\end{equation*}
$$

From now on we will frequently use the notation $\widehat{x}=\left(x_{1}, x_{2}\right), \widehat{x} \in \widehat{\Omega}$, and all functions depending only on ( $x_{1}, x_{2}$ ) will be marked by ${ }^{\wedge}$; in addition, we will distinguish between the 3 - and 2-dimensional divergence operator:

$$
\operatorname{Div} \boldsymbol{\tau}=\frac{\partial \boldsymbol{\tau}_{1}}{\partial x_{1}}+\frac{\partial \boldsymbol{\tau}_{2}}{\partial x_{2}}+\frac{\partial \boldsymbol{\tau}_{3}}{\partial x_{3}}, \quad \operatorname{div} \widehat{\boldsymbol{\tau}}=\frac{\partial \widehat{\boldsymbol{\tau}}_{1}}{\partial x_{1}}+\frac{\partial \widehat{\boldsymbol{\tau}}_{2}}{\partial x_{2}} .
$$

We also denote $\widehat{F}_{\ominus}(\widehat{x}):=F_{\ominus}\left(\widehat{x}, d_{\ominus}(\widehat{x})\right), \widehat{F}_{\oplus}(\widehat{x}):=F_{\oplus}\left(\widehat{x}, d_{\oplus}(\widehat{x})\right)$ for any $\widehat{x} \in \widehat{\Omega}$. Finally, we define the energy norm

$$
\begin{equation*}
\left\|\|v\| \mid:=\left(\int_{\Omega} \mathbf{A}(x) \nabla v \cdot \nabla v d x\right)^{1 / 2} \quad \forall v \in V_{0}\right. \tag{2.9}
\end{equation*}
$$

## 3 The reduced problem

In view of (2.1), it is reasonable to consider the hypothesis that

$$
\begin{equation*}
\text { the exact solution } u \text { is almost constant with respect to } x_{3} \text {-coordinate. } \tag{3.1}
\end{equation*}
$$

This gives rise to the so-called zero-order reduced model for the original problem (2.8). The model is very popular due to its simplicity and purely two-dimensional formulation. The discussion on the hierarchy of reduced models of different orders can be found in, e.g., [11], [3].

With (3.1) in mind, one can expect that the exact solution $u$ may be well-approximated by the functions from the subspace

$$
\begin{equation*}
\widehat{V}_{0}:=\left\{v \in V_{0} \mid \exists \widehat{v} \in H_{0}^{1}(\widehat{\Omega}) \text { such that } v(x)=\widehat{v}(\widehat{x}) \text { for a.e. } x=\left(\widehat{x}, x_{3}\right) \in \Omega\right\} \tag{3.2}
\end{equation*}
$$

Thus, any function from $\widehat{V}_{0}$ can be identified with the corresponding function $\widehat{v} \in H_{0}^{1}(\widehat{\Omega})$ (and vice versa, for any $\widehat{v} \in H_{0}^{1}(\widehat{\Omega})$ one can reconstruct $v \in \widehat{V}_{0} \subset V_{0}$ by the constant extension as in the definition of $\widehat{V}_{0}$ ). Then, the energy-norm projection of $u$ onto the subspace $\widehat{V}_{0}$ yields the reduced problem (the zero-order reduced model)
Problem $(\widehat{P})$ : Find $\widehat{u} \in \widehat{V}_{0}$ such that

$$
\begin{equation*}
\int_{\Omega} \mathbf{A} \nabla \widehat{u} \cdot \nabla \widehat{w} d x=\int_{\Omega} f \widehat{w} d x+\int_{\Gamma_{\ominus}} F_{\ominus} \widehat{w} d s+\int_{\Gamma_{\oplus}} F_{\oplus} \widehat{w} d s \quad \forall \widehat{w} \in \widehat{V}_{0} . \tag{3.3}
\end{equation*}
$$

Now we can define the dimension reduction error (the modelling error) as the difference $e:=u-\widehat{u}$ between the solutions to the original problem (2.8) and to the reduced problem (3.3).
Remark 3.1 It may be noticed that assumption (2.1) (and, consequently, (3.1)) serves only as an intuitive motivation for the introduction of the approximation subspace $\widehat{V}_{0}$ and the reduced problem (3.3). Since the assumption cannot be quantified, the real error of "replacing" $u$ by $\widehat{u}$ may be large; a robust a-posteriori error estimator should, however, measure this error sufficiently accurately even in the cases when assumption (2.1) is virtually unsatisfied.
Remark 3.2 The asymptotic behaviour of the modelling error $e$ was analysed in [11] (see also $[2]$ ) for the case of a plate with plane parallel faces $\Gamma_{\ominus}$ and $\Gamma_{\oplus}$ (i.e. when $d_{\ominus}=-\frac{d_{0}}{2}, d_{\oplus}=\frac{d_{0}}{2}$, $d_{0}=$ const $>0$ is the plate thickness) and $f=0$. It was proved that

$$
\mid\|e\| \| \leq C d_{0}^{1 / 2}\left(\left\|\widehat{F}_{\ominus}\right\|_{L_{2}(\widehat{\Omega})}+\left\|\widehat{F}_{\oplus}\right\|_{L_{2}(\widehat{\Omega})}\right) \text { as } d_{0} \rightarrow 0
$$

Remark 3.3 We have to note that the third component of the vector $\nabla \widehat{u}$ is zero (since $\widehat{u}$ does not depend on $x_{3}$ ) and, thus, the vector will sometimes be considered as a two-component vector, if no confusion is possible.

In order to see that the reduced problem (3.3) is, in fact, a two-dimensional problem, we define the operation ( $\sim$ ) of averaging in $x_{3}$-direction

$$
\forall g \in L_{1}(\Omega): \quad \widetilde{g}(\widehat{x}):=\frac{1}{d(\widehat{x})} \int_{d_{\ominus}(\widehat{x})}^{d_{\oplus}(\hat{x})} g\left(\widehat{x}, x_{3}\right) d x_{3} \text { for a.e. } \widehat{x} \in \widehat{\Omega},
$$

and, having noticed that

$$
\int_{\Gamma_{\ominus}} F_{\ominus} \widehat{w} d s=\int_{\widehat{\Omega}} \widehat{F}_{\ominus}(\widehat{x}) \widehat{w}(\widehat{x}) \sqrt{1+\left|\nabla d_{\ominus}(\widehat{x})\right|^{2}} d \widehat{x} \quad \text { (analogously for } \int_{\Gamma_{\oplus}} F_{\oplus} \widehat{w} d s \text { ), }
$$

we can rewrite problem (3.3) as follows:
Find $\widehat{u} \in \widehat{V}_{0}$ such that

$$
\begin{equation*}
\int_{\widehat{\Omega}} d(\widehat{x}) \widetilde{\mathbf{A}}_{p}(\widehat{x}) \nabla \widehat{u} \cdot \nabla \widehat{w} d \widehat{x}=\int_{\widehat{\Omega}} d(\widehat{x}) \widehat{f}(\widehat{x}) \widehat{w} d \widehat{x} \quad \forall \widehat{w} \in \widehat{V}_{0} . \tag{3.4}
\end{equation*}
$$

Here $\widetilde{\mathbf{A}}_{p}(\widehat{x})=\left(\widetilde{a}_{i j}(\widehat{x})\right)_{i, j=\overline{1,2}}$ is the averaged "plane" part $\mathbf{A}_{p}(x)\left(\mathbf{A}_{p}(x)=\left(a_{i j}(x)\right)_{i, j=\overline{1,2}}\right)$ of the matrix $\mathbf{A}$ and

$$
\widehat{f}(\widehat{x})=\widetilde{f}(\widehat{x})+\frac{\widehat{F}_{\ominus}(\widehat{x}) \sqrt{1+\left|\nabla d_{\ominus}(\widehat{x})\right|^{2}}+\widehat{F}_{\oplus}(\widehat{x}) \sqrt{1+\left|\nabla d_{\oplus}(\widehat{x})\right|^{2}}}{d(\widehat{x})} .
$$

It is clear that problem (3.4) is a two-dimensional elliptic problem with the homogeneous Dirichlet boundary condition:

$$
\begin{align*}
-\operatorname{div}\left(d(\widehat{x}) \widetilde{\mathbf{A}}_{p}(\widehat{x}) \nabla \widehat{u}\right) & =d(\widehat{x}) \widehat{f}(\widehat{x}) \quad \text { in } \widehat{\Omega}  \tag{3.5}\\
\widehat{u} & =0 \quad \text { on } \widehat{\Gamma} . \tag{3.6}
\end{align*}
$$

## 4 A posteriori estimation of the modelling error

In order to control the dimension reduction error, we apply the functional-type a posteriori error estimate derived in [8] (see also [5] and [7]) to the original three-dimensional problem (2.8). The estimate reads as follows:
For all $\gamma>0, \delta>0$ and $y^{*} \in H_{*}(\Omega$, Div $)$ there holds

$$
\begin{equation*}
\|\|u-v\|\|^{2} \leq(1+\gamma) M_{1}^{2}+\left(1+\frac{1}{\gamma}\right)(1+\delta) C_{\Omega}^{2} M_{2}^{2}+\left(1+\frac{1}{\gamma}\right)\left(1+\frac{1}{\delta}\right) C_{\Gamma}^{2}\left(1+C_{\Omega}^{2}\right) M_{3}^{2}, \tag{4.1}
\end{equation*}
$$

where $v$ is any function from the energy space $V_{0}, C_{\Omega}$ is the constant from Friedrichs' inequality,

$$
\begin{equation*}
C_{\Omega}^{-2}=\inf _{w \in V_{0} \backslash\{0\}} \frac{\|w\|^{2}}{\|w\|_{L_{2}(\Omega)}^{2}}, \tag{4.2}
\end{equation*}
$$

$C_{\Gamma}$ is the constant from the trace inequality,

$$
\begin{equation*}
C_{\Gamma}^{2}=\sup _{w \in V_{0} \backslash\{0\}} \frac{\|w\|_{L_{2}\left(\Gamma_{\oplus}\right)}^{2}+\|w\|_{L_{2}\left(\Gamma_{\ominus}\right)}^{2}}{\|w \mid\|^{2}+\|w\|_{L_{2}(\Omega)}^{2}}, \tag{4.3}
\end{equation*}
$$

the space $H_{*}(\Omega$, Div $)$ is defined as

$$
H_{*}(\Omega, \operatorname{Div}):=\left\{y^{*} \in L_{2}\left(\Omega, \mathbb{R}^{3}\right) \mid \operatorname{Div} y^{*} \in L_{2}(\Omega), y^{*} \cdot \boldsymbol{\nu}_{\ominus} \in L_{2}\left(\Gamma_{\ominus}\right), y^{*} \cdot \boldsymbol{\nu}_{\oplus} \in L_{2}\left(\Gamma_{\oplus}\right)\right\}
$$

and the functionals $M_{1}^{2}, M_{2}^{2}, M_{3}^{2}$ are defined by

$$
\begin{aligned}
& M_{1}^{2}:=\int_{\Omega}\left(\nabla v-\mathbf{A}^{-1} y^{*}\right) \cdot\left(\mathbf{A} \nabla v-y^{*}\right) d x \\
& M_{2}^{2}:=\left\|\operatorname{Div} y^{*}+f\right\|_{L_{2}(\Omega)}^{2} \\
& M_{3}^{2}:=\left\|F_{\ominus}-y^{*} \cdot \boldsymbol{\nu}_{\ominus}\right\|_{L_{2}\left(\Gamma_{\ominus}\right)}^{2}+\left\|F_{\oplus}-y^{*} \cdot \boldsymbol{\nu}_{\oplus}\right\|_{L_{2}\left(\Gamma_{\oplus}\right)}^{2} .
\end{aligned}
$$

Since estimate (4.1) holds true for any "approximate solution" $v$ from $V_{0}$ and the solution to the reduced problem $\widehat{u} \in \widehat{V}_{0} \subset V_{0}$, we can simply plug $\widehat{u}$ into estimate (4.1) to obtain an upper bound of the modelling error. We also emphasize that the estimate is valid for any positive numbers $\gamma$ and $\delta$ and for any vector-function $y^{*}$ from the space $H_{*}(\Omega, \operatorname{Div})$. While the best possible option would be to take as $y^{*}$ the exact flux $\mathbf{A} \nabla u$ (then $M_{2}$ and $M_{3}$ would vanish and $M_{1}$ would give us the energy norm of the exact error), we have to restrict ourselves to choosing some computable quantity, i.e. not containing the unknown exact solution $u$. We approximate the flux by

$$
\begin{equation*}
y^{*}=\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\boldsymbol{\tau}^{*}, \tag{4.4}
\end{equation*}
$$

with $\boldsymbol{\tau}^{*}=\{0,0, \psi(x)\}^{T}, \psi$ is the auxiliary function from $L_{2}(\Omega)$ such that $\frac{\partial \psi}{\partial x_{3}} \in L_{2}(\Omega), \psi \in L_{2}\left(\Gamma_{\ominus}\right)$ and $\psi \in L_{2}\left(\Gamma_{\oplus}\right)$. The concrete form of the function $\psi$ will be given later; its meaning becomes clear in the case of the Poisson equation (i.e. if $\mathbf{A}$ is the identity matrix), where $\psi$ should, obviously, approximate the derivative $\frac{\partial u}{\partial x_{3}}$ of the exact solution in $x_{3}$-direction. Using (3.5), it is easy to verify that $y^{*}$ from (4.4) belongs to $H_{*}(\Omega$, Div $)$.
Remark 4.1 The estimate (4.1) possesses the property of asymptotic exactness (see [8]), but, if we choose $y^{*}$ as in (4.4), this property might be lost, since the only remaining "degree of freedom" is the function $\psi$ and the approximate plane flux $\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}$ may not represent the first two components of the exact flux $\mathbf{A} \nabla u$ sufficiently well. On the other hand, if we did not fix the first two components of $y^{*}$, the process of estimation would require the minimization of the right-hand side of (4.1) with respect to those components, which is, in principle, equivalent to solving a three-dimensional problem. However, our goal is to avoid any truly three-dimensional calculations in the evaluation of the error estimator (this process should not be more expensive than the solution of the reduced problem). Fortunately, in most of the situations, $\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}$ is a good approximation to the "plane" part of the exact flux, and the modelling-error estimate with $y^{*}$ as in (4.4) exhibits both efficiency and flexibility, as the numerical tests of Section 6 show.

In order to rewrite estimate (4.1) in a more convenient form, we introduce the notation:

$$
\begin{align*}
& \mathbf{B}:=\mathbf{A}^{-1} \quad\left(\mathbf{B}(x)=\left(b_{i j}(x)\right)_{i, j=\overline{1,3}}, \quad \mathbf{B}=\mathbf{B}^{T}\right),  \tag{4.5}\\
& \mathbf{B}_{p}:=\left(b_{i j}\right)_{i, j=\overline{1,2}},  \tag{4.6}\\
& \mathbf{b}_{3}:=\left\{b_{31}, b_{32}\right\}^{T} . \tag{4.7}
\end{align*}
$$

The term $M_{1}^{2}$ with $v=\widehat{u}$ reads

$$
\begin{equation*}
M_{1}^{2}=\int_{\Omega}\left(\nabla \widehat{u}-\mathbf{B} y^{*}\right) \cdot\left(\mathbf{A} \nabla \widehat{u}-y^{*}\right) d x=\int_{\Omega}\left(\mathbf{A} \nabla \widehat{u} \cdot \nabla \widehat{u}-2 y^{*} \cdot \nabla \widehat{u}+\mathbf{B} y^{*} \cdot y^{*}\right) d x \tag{4.8}
\end{equation*}
$$

For the first term in (4.8), one immediately obtains

$$
\begin{equation*}
\int_{\Omega} \mathbf{A} \nabla \widehat{u} \cdot \nabla \widehat{u} d x=\int_{\widehat{\Omega}} d(\widehat{x}) \widetilde{\mathbf{A}}_{p}(\widehat{x}) \nabla \widehat{u} \cdot \nabla \widehat{u} d \widehat{x} \tag{4.9}
\end{equation*}
$$

The second term in (4.8) can be further rewritten if one notices that

$$
y^{*} \cdot \nabla \widehat{u}=\left(\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\tau^{*}\right) \cdot \nabla \widehat{u}=\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \nabla \widehat{u} .
$$

Thus,

$$
\begin{equation*}
\int_{\Omega} y^{*} \cdot \nabla \widehat{u} d x=\int_{\widehat{\Omega}} d(\widehat{x}) \widetilde{\mathbf{A}}_{p}(\widehat{x}) \nabla \widehat{u} \cdot \nabla \widehat{u} d \widehat{x} . \tag{4.10}
\end{equation*}
$$

For the third term in (4.8) we have

$$
\begin{aligned}
& \mathbf{B} y^{*} \cdot y^{*}=\left(\mathbf{B} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\mathbf{B} \tau^{*}\right) \cdot\left(\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\tau^{*}\right)=\mathbf{B} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\mathbf{B} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \tau^{*} \\
&+\mathbf{B} \tau^{*} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\mathbf{B} \tau^{*} \cdot \tau^{*}=\mathbf{B}_{p} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+2\left(\mathbf{b}_{3} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}\right) \psi+b_{33} \psi^{2}
\end{aligned}
$$

that yields

$$
\begin{equation*}
\int_{\Omega} \mathbf{B} y^{*} \cdot y^{*} d x=\int_{\widehat{\Omega}} d(\widehat{x}) \widetilde{\mathbf{B}}_{p} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} d \widehat{x}+\int_{\Omega}\left(b_{33} \psi(x)^{2}+2\left(\mathbf{b}_{3} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}\right) \psi(x)\right) d x \tag{4.11}
\end{equation*}
$$

where $\widetilde{\mathbf{B}}_{p}$ is the averaged "plane" part $\mathbf{B}_{p}(x)$ of the matrix $\mathbf{B}(x)$.
Substituting (4.9), (4.10) and (4.11) into (4.8) one obtains

$$
\begin{equation*}
M_{1}^{2}=\int_{\widehat{\Omega}} d(\widehat{x})\left(\widetilde{\mathbf{B}}_{p} \widetilde{\mathbf{A}}_{p}-\mathbf{I}\right) \nabla \widehat{u} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} d \widehat{x}+\int_{\Omega}\left(b_{33} \psi(x)^{2}+2\left(\mathbf{b}_{3} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}\right) \psi(x)\right) d x, \tag{4.12}
\end{equation*}
$$

where $\mathbf{I}$ is the identity $2 \times 2$-matrix. It is interesting to note that the first integral in (4.12) represents the error in averaging the coefficient matrix $\mathbf{A}(x)$; this becomes fully transparent in the case of a block-diagonal matrix $\mathbf{A}$, i.e. when $a_{31}=a_{32}=0$ (then $\mathbf{B}_{p}=\mathbf{A}_{p}^{-1}$ and without the averaging the integral would be identically zero).

The functional $M_{2}^{2}$ of (4.1) can be also rearranged if one takes $y^{*}$ as in (4.4). First, note that

$$
\operatorname{Div} y^{*}=\operatorname{div} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\frac{\partial \psi}{\partial x_{3}} .
$$

From (3.5) one can deduce

$$
\operatorname{div} \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}=-\widehat{f}-\frac{\nabla d}{d} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u} .
$$

Hence,

$$
\begin{equation*}
M_{2}^{2}=\left\|f-\tilde{f}-\frac{\widehat{F}_{\ominus} \sqrt{1+\left|\nabla d_{\ominus}\right|^{2}}+\widehat{F}_{\oplus} \sqrt{1+\left|\nabla d_{\oplus}\right|^{2}}}{d}-\frac{\nabla d}{d} \cdot \widetilde{\mathbf{A}}_{p} \nabla \widehat{u}+\frac{\partial \psi}{\partial x_{3}}\right\|_{L_{2}(\Omega)}^{2} . \tag{4.13}
\end{equation*}
$$

The term $M_{3}^{2}$ with $y^{*}$ from (4.4) reads

$$
\begin{equation*}
M_{3}^{2}=\left\|F_{\ominus}-\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \boldsymbol{\nu}_{\ominus}-\psi \nu_{\ominus 3}\right\|_{L_{2}\left(\Gamma_{\ominus}\right)}^{2}+\left\|F_{\oplus}-\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \boldsymbol{\nu}_{\oplus}-\psi \nu_{\oplus 3}\right\|_{L_{2}\left(\Gamma_{\oplus}\right)}^{2}, \tag{4.14}
\end{equation*}
$$

where $\widetilde{\mathbf{A}}_{p} \nabla \widehat{u}$ is considered as a vector in $\mathbb{R}^{3}$ with the third component equal to zero, and

$$
\nu_{\ominus 3}=\frac{-1}{\sqrt{1+\left|\nabla d_{\ominus}\right|^{2}}}, \quad \nu_{\oplus 3}=\frac{1}{\sqrt{1+\left|\nabla d_{\oplus}\right|^{2}}}
$$

are the third components of the normal vectors $\boldsymbol{\nu}_{\ominus}$ and $\boldsymbol{\nu}_{\oplus}$.
Now we can write the general a-posteriori estimate for the dimension reduction error: For all $\gamma>0$ and $\delta>0$ there holds

$$
\begin{equation*}
\|\|u-\widehat{u}\|\|^{2} \leq(1+\gamma) M_{1}^{2}+\left(1+\frac{1}{\gamma}\right)(1+\delta) C_{\Omega}^{2} M_{2}^{2}+\left(1+\frac{1}{\gamma}\right)\left(1+\frac{1}{\delta}\right) C_{\Gamma}^{2}\left(1+C_{\Omega}^{2}\right) M_{3}^{2}, \tag{4.15}
\end{equation*}
$$

where the constants $C_{\Omega}$ and $C_{\Gamma}$ are as above (see (4.2) and (4.3)), the functionals $M_{1}^{2}, M_{2}^{2}$ and $M_{3}^{2}$ are given by (4.12), (4.13) and (4.14).

In estimate (4.15) we still have the freedom of choosing the auxiliary function $\psi$. The simplest choice is to take such $\psi$ that the term $M_{3}$ (i.e. the residual on the Neumann boundary condition) would be identically zero. To do so, we first rewrite the $L_{2}$-norms on $\Gamma_{\oplus, \ominus}$ in (4.14) as the integrals over $\widehat{\Omega}$ :

$$
\left\|F_{\ominus}-\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \boldsymbol{\nu}_{\ominus}-\psi \nu_{\ominus 3}\right\|_{L_{2}\left(\Gamma_{\ominus}\right)}^{2}=\int_{\widehat{\Omega}}\left(\widehat{F}_{\ominus}(\widehat{x})-\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \boldsymbol{\nu}_{\ominus}-\psi\left(\widehat{x}, d_{\ominus}(\widehat{x})\right) \nu_{\ominus 3}\right)^{2} \sqrt{1+\left|\nabla d_{\ominus}\right|^{2}} d \widehat{x}
$$

(analogously for the norm in $L_{2}\left(\Gamma_{\oplus}\right)$ ). Then, we denote

$$
\widehat{G}_{\oplus, \ominus}:=\widehat{F}_{\oplus, \ominus}-\widetilde{\mathbf{A}}_{p} \nabla \widehat{u} \cdot \boldsymbol{\nu}_{\oplus, \ominus}
$$

and set

$$
\begin{equation*}
\psi_{1}(x)=\widehat{\alpha}(\widehat{x}) x_{3}+\widehat{\beta}(\widehat{x}), \tag{4.16}
\end{equation*}
$$

where the functions $\widehat{\alpha}$ and $\widehat{\beta}\left(\widehat{\alpha}, \widehat{\beta} \in L_{2}(\widehat{\Omega})\right)$ are chosen so that

$$
\begin{equation*}
\psi_{1} \nu_{\oplus 3}=\widehat{G}_{\oplus} \text { at } x_{3}=d_{\oplus}, \quad \psi_{1} \nu_{\ominus 3}=\widehat{G}_{\ominus} \text { at } x_{3}=d_{\ominus} . \tag{4.17}
\end{equation*}
$$

As $\nu_{\ominus 3}, \nu_{\oplus 3}$ belong to $L_{\infty}(\widehat{\Omega})$ and cannot be zero in $\widehat{\Omega}$, the functions $\widehat{\alpha}$ and $\widehat{\beta}$ are uniquely defined by the conditions (4.17):

$$
\begin{align*}
& \widehat{\alpha}=\frac{1}{d}\left(\frac{\widehat{G}_{\oplus}}{\nu_{\oplus 3}}-\frac{\widehat{G}_{\ominus}}{\nu_{\ominus 3}}\right)  \tag{4.18}\\
& \widehat{\beta}=\frac{1}{d}\left(\frac{\widehat{G}_{\ominus}}{\nu_{\ominus 3}} d_{\oplus}-\frac{\widehat{G}_{\oplus}}{\nu_{\oplus 3}} d_{\ominus}\right) . \tag{4.19}
\end{align*}
$$

It is obvious that the function $\psi_{1}$ as well as its derivative in $x_{3}$-direction belong to $L_{2}(\Omega), \psi_{1}$ belongs to $L_{2}\left(\Gamma_{\oplus}\right)$ and $L_{2}\left(\Gamma_{\ominus}\right)$ (since $\left.\left.\psi_{1}\right|_{x_{3}=d_{\oplus, \ominus}(\widehat{x})} \in L_{2}(\widehat{\Omega})\right)$, and, moreover, with such a function $\psi$ the term $M_{3}$ becomes zero.
Remark 4.2 One can also consider a quadratic (with respect to $x_{3}$ ) function

$$
\psi_{2}(x)=\psi_{1}(x)+\widehat{\eta}(\widehat{x})\left(x_{3}-d_{\oplus}(\widehat{x})\right)\left(x_{3}-d_{\ominus}(\widehat{x})\right)
$$

with $\widehat{\eta}$ being an arbitrary function from $L_{2}(\widehat{\Omega})$. The substitution of $\psi_{2}$ into (4.14) instead of $\psi$ will evidently imply $M_{3}=0$. In the second numerical example of Section 6 we will use $\psi_{2}$ because of the freedom in the choice of the function $\widehat{\eta}$. It is clear that one can quite analogously construct the functions $\left\{\psi_{m}\right\}, m=3,4, \ldots$, which would make the $M_{3}$-term vanish and could, possibly, allow to approximate the third component of the exact flux $\mathbf{A} \nabla u$ with a higher accuracy.

Having chosen the function $\psi$ such that $M_{3}=0$, one can obtain from (4.15) the following estimate for the squared energy norm of the modelling error:

$$
\begin{equation*}
\|u-\widehat{u}\| \|^{2} \leq(1+\gamma) M_{1}^{2}+\left(1+\frac{1}{\gamma}\right) C_{\Omega}^{2} M_{2}^{2} \tag{4.20}
\end{equation*}
$$

where $\gamma$ is any positive number, $C_{\Omega}$ is the Friedrichs constant, $M_{1}^{2}$ and $M_{2}^{2}$ are given by (4.12) and (4.13). Minimizing the right-hand side of (4.20) with respect to the scalar parameter $\gamma>0$, we immediately arrive at the estimate for the energy norm of the modelling error

$$
\begin{equation*}
\|\|u-\widehat{u}\|\| \leq M:=M_{1}+C_{\Omega} M_{2} \tag{4.21}
\end{equation*}
$$

with $M_{1}$ and $M_{2}$ defined by (4.12) and (4.13).
The rest of the paper will be devoted to the analysis of the properties of estimates (4.20), (4.21).

## 5 Particular cases

The error majorant $M$ in (4.21) has been derived for quite general geometry of $\Omega$ and coefficient matrix $\mathbf{A}(x)$; to make the estimate more transparent, we consider two particular cases.

### 5.1 Plate of constant thickness

We assume that

$$
\begin{equation*}
d_{\oplus}=d_{\ominus}+d_{0} \quad\left(d_{0}=\text { const }>0\right) \tag{5.1}
\end{equation*}
$$

and, in addition, that

$$
\begin{align*}
& \mathbf{A}=\mathbf{A}(\widehat{x}) \quad \text { (this immediately implies } \mathbf{B}=\mathbf{B}(\widehat{x})),  \tag{5.2}\\
& \left.a_{31}=a_{32}=0 \quad \text { (this yields } \mathbf{B}_{p}=\mathbf{A}_{p}^{-1}, b_{33}=a_{33}^{-1}, b_{31}=b_{32}=0\right) \tag{5.3}
\end{align*}
$$

With these assumptions and the choice $\psi=\psi_{1}$ (see (4.16)) the terms $M_{1}$ and $M_{2}$ in estimate (4.21) become simpler:

$$
\begin{equation*}
M_{1}=\left(\int_{\Omega} a_{33}^{-1} \psi_{1}^{2} d x\right)^{1 / 2}, \quad M_{2}=\|f-\widetilde{f}\|_{L_{2}(\Omega)} \tag{5.4}
\end{equation*}
$$

One may notice that the integral in the first term $M_{1}$ of the error majorant $M$ can be rewritten as

$$
\int_{\Omega} a_{33}^{-1} \psi_{1}^{2} d x=d_{0} \cdot \int_{\widehat{\Omega}} a_{33}^{-1}\left(\widehat{\alpha}^{2} \frac{d_{\oplus}^{2}+d_{\oplus} d_{\ominus}+d_{\ominus}^{2}}{3}+\widehat{\alpha} \widehat{\beta}\left(d_{\oplus}+d_{\ominus}\right)+\widehat{\beta}^{2}\right) d \widehat{x}
$$

which means that the term $M_{1}$ is of order $\mathcal{O}\left(d_{0}^{1 / 2}\right)$ when the plate thickness $d_{0}$ tends to zero. If $f \in L_{\infty}(\Omega)$, the second term $M_{2}$ is obviously of the same order $\mathcal{O}\left(d_{0}^{1 / 2}\right)$, i.e. the whole estimator $M$ converges to zero with the rate $\mathcal{O}\left(d_{0}^{1 / 2}\right)$ as $d_{0} \rightarrow 0$. This is the optimal convergence rate for the modelling error $e$ in the energy norm, as was shown in [11] for the simpler case of a plate with plane parallel faces and $f=0$ (see Remark 3.2). It is worth noting that, if $f \in C^{1}(\Omega)$, the second term in $M$ is of higher order $\mathcal{O}\left(d_{0}^{3 / 2}\right)$ as compared to the first term.

### 5.2 Plate with plane parallel faces

If, in addition to (5.2), (5.3), we strengthen the assumption (5.1) replacing it by

$$
\begin{equation*}
d_{\oplus}=\frac{d_{0}}{2}, d_{\ominus}=-\frac{d_{0}}{2} \quad\left(d_{0}=\text { const }>0\right), \tag{5.5}
\end{equation*}
$$

the function $\psi_{1}$ takes the simple form

$$
\psi_{1}(x)=\frac{\widehat{F}_{\oplus}(\widehat{x})+\widehat{F}_{\ominus}(\widehat{x})}{d_{0}} x_{3}+\frac{\widehat{F}_{\oplus}(\widehat{x})-\widehat{F}_{\ominus}(\widehat{x})}{2}
$$

and the error estimate (4.21) reduces to

$$
\begin{equation*}
\|\|u-\widehat{u}\|\| \leq \sqrt{\frac{d_{0}}{3}}\left(\int_{\widehat{\Omega}} a_{33}^{-1}\left(\widehat{F}_{\oplus}^{2}+\widehat{F}_{\ominus}^{2}-\widehat{F}_{\oplus} \widehat{F}_{\ominus}\right) d \widehat{x}\right)^{1 / 2}+C_{\Omega}\|f-\widetilde{f}\|_{L_{2}(\Omega)} \tag{5.6}
\end{equation*}
$$

If we set here $f=0, a_{33}=1$ and $\widehat{F}_{\oplus}=\widehat{F}_{\ominus}=\widehat{F}$, we obtain

$$
\begin{equation*}
\left\lvert\,\|u-\widehat{u}\|\left\|\leq \sqrt{\frac{d_{0}}{3}}\right\| \widehat{F}\right. \|_{L_{2}(\widehat{\Omega})} \tag{5.7}
\end{equation*}
$$

that is exactly the estimator of Babuška and Schwab (see [2]) for the zero-order reduced model. Thus, the latter estimator can be obtained as a particular case of the error majorant (4.21) if one makes the assumptions (5.2), (5.3), (5.5) and sets $f=0$. This fact is especially interesting, since we advocate the estimation approach (see the details in [8]) completely different from the one utilized in [2].
Remark 5.1 The error estimate (4.21) contains the Friedrichs constant $C_{\Omega}$ that must be, in general, evaluated numerically. The constant depends solely on the geometry of the domain $\Omega$ and can be computed as $1 / \sqrt{\lambda}$ where $\lambda$ is the minimal eigenvalue of the elliptic operator $-\operatorname{Div}(\mathbf{A} \nabla \cdot)$ equipped with the homogeneous Dirichlet condition on $\Gamma_{0}$ and homogeneous Neumann conditions on $\Gamma_{\oplus, \ominus}$ (see (4.2)). It is clear that, in the case of a plate with plane parallel faces, $C_{\Omega}$ can be easily estimated from above, if one computes the Friedrichs constant in a larger domain obtained by embedding the cross-section $\widehat{\Omega}$ of $\Omega$ into some rectangle; the faces of this larger domain are, then, obtained by the extension of plane faces of $\Omega$. Yet simpler but rougher upper estimate for $C_{\Omega}$ in the case of a plate with plane parallel faces is given by $(\operatorname{diam} \widehat{\Omega}) / c$, where $c$ is the lower bound of the minimal eigenvalue of the matrix $\mathbf{A}(x)$ in $\Omega$ (see (2.6)). It is worth noticing that the constant $C_{\Omega}$ multiplies in the majorant the term $M_{2}$ which is often of higher order as compared to the first term $M_{1}$ (it is so, for example, in the considered above particular cases, when the function $f$ is smooth). Then, the possible error of overestimation of $C_{\Omega}$ is harmless for the majorant accuracy.

## 6 Numerical examples

### 6.1 Numerical test 1

In order to analyse the performance of the proposed error estimator, we consider a two-dimensional test problem in the "sine-shape" domain (see Figure 2 (left)) whose upper and lower faces are given by

$$
d_{\oplus, \ominus}(x)=\sin (k \pi x) \pm \frac{d_{0}}{2}, \quad k=1,2, \ldots
$$

where $d_{0}>0$ is the domain thickness. In this example, $\widehat{\Omega}=(0,1)$ and $\Omega=\left\{(x, y) \in \mathbb{R}^{2} \mid x \in\right.$ $\left.\widehat{\Omega}, d_{\ominus}(x)<y<d_{\oplus}(x)\right\}$. The considered problem is

$$
\begin{aligned}
-\Delta u & =f & & \text { in } \Omega, \\
u & =0 & & \text { at } x=0 \text { and } x=1, \\
\nabla u \cdot \boldsymbol{\nu}_{\oplus, \ominus} & =F_{\oplus, \ominus} & & \text { at } y=d_{\oplus, \ominus},
\end{aligned}
$$

and the right-hand sides of the equation and of the boundary condition are computed using the exact solution

$$
u(x, y)=\sin (\pi x) \cdot y^{m} \quad(m=1,2, \ldots)
$$

The reduced problem (3.3) is, in this case, a one-dimensional Dirichlet problem that, of course, can be solved very accurately (in the present work, we address the estimation of the modelling error only, assuming that the discretization error stemming from the solution of the reduced problem is negligible). The Friedrichs constant $C_{\Omega}$ was evaluated by computing the minimal eigenvalue of the Laplace operator with the corresponding homogeneous Dirichlet/Neumann boundary conditions (see Remark 5.1). We found that, for each $k=1,2, \ldots, C_{\Omega}$ is an increasing function of the thickness $d_{0}$ as $d_{0} \rightarrow 0$. There always exists, however, a clear upper bound for $C_{\Omega}$; in particular, the estimates $C_{\Omega} \leq \sqrt{2}$ for $k=2$ and $C_{\Omega} \leq 3$ for $k=4$ hold true.

Figure 2 (right) shows the convergence rates of the exact modelling-error in the energy norm $\left(\left\|||e| \|)\right.\right.$ and of the error majorant $(M)$ as the domain thickness $d_{0}$ tends to zero (the analysis here corresponds to the case $k=2$, when the domain $\Omega$ has the shape depicted in Figure 2 (left)). It is


Figure 2: (left) The domain geometry; (right) Convergence rate of the exact modelling-error and of the error majorant, $k=2, m=4$ (solid lines) and $m=5$ (dash-dot lines), the majorant is indicated by "o".


Figure 3: Local error distribution provided by the exact modelling-error (solid line) and by the $M_{1}$-term of the majorant (dash-dot line), $k=4, m=4$ : (left) $d_{0}=0.1, \quad$ (right) $d_{0}=0.05$.

| $d_{0}^{-1}$ | $m=4$ |  |  | $m=5$ |  |  |
| :--- | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $\\|\\|e\\|\\|$ | $M$ | $\frac{M}{\\|e\\| \\|}$ | $\\|\|\mid e\\| \\|$ | $M$ | $\frac{M}{\\|e\\| \\|}$ |
| $10^{0}$ | 3.2108 | 9.5598 | 2.9774 | 5.0842 | 16.8434 | 3.3129 |
| $10^{1}$ | 0.5058 | 0.5690 | 1.1250 | 0.6399 | 1.3481 | 2.1066 |
| $10^{2}$ | 0.1581 | 0.1598 | 1.0106 | 0.1991 | 0.3937 | 1.9770 |
| $10^{3}$ | 0.0500 | 0.0501 | 1.0010 | 0.0630 | 0.1237 | 1.9650 |
| $10^{4}$ | 0.0158 | 0.0158 | 1.0000 | 0.0199 | 0.0391 | 1.9638 |

Table 1: Convergence of the exact modelling-error in the energy norm ( $\|\|e\|\|$ ) and of the error majorant ( $M$ ) as $d_{0} \rightarrow 0(k=2)$; the results are rounded up to $10^{-4}$.
clear that both the exact error and the majorant vanish with the theoretically predicted, optimal rate $\mathcal{O}\left(d_{0}^{1 / 2}\right)$. However, the behaviour of the majorant is different for even and odd values of degree $m$ determining the polynomial growth of the exact solution $u$ in $y$-direction. The typical picture corresponding to an even value of the parameter $m$ is well represented by the case $m=4$ in Figure 2 (right); in this case, the majorant $M$ demonstrates the asymptotic exactness, and, moreover, the effectivity index $\frac{M}{\|e\| \|}$ behaves like $1+\mathcal{O}\left(d_{0}\right)$ (see Table 1 ). In the case of an odd value of $m$ (represented by $m=5$ in Figure 2 (right)), the majorant loses the property of asymptotic exactness, although the effectivity index remains stable and behaves like, approximately, $1.963+\mathcal{O}\left(d_{0}\right)$ (see Table 1). This problem was addressed in Remark 4.1 and is caused by the fact that the approximate flux computed in the reduced model does not bring sufficient information on the corresponding components of the exact flux. We may note, however, that the effectivity index is still quite acceptable in this case. Finally, it is worth noticing that the presented error estimator provides a reliable upper bound for the exact error at any positive values of the domain thickness $d_{0}$, i.e. also in the cases when the domain is not "thin" at all.

The local error distribution provided by the exact error and by the first, $M_{1}$-term of the majorant $M$ (see (5.4)) are depicted in Figure 3 (here we consider the case $k=4$, when the functions $d_{\oplus, \ominus}$ defining the shape of the domain have 4 extrema). The figure shows that already for rather large value of the domain thickness $d_{0}=0.1$ the majorant delivers sufficiently accurate information on the location of the regions of the biggest modelling error, while for $d_{0}=0.05$ the exact and the estimated error distributions are practically coincident.

### 6.2 Numerical test 2

The previous test shows that in the standard situations the proposed error estimator performs well. The example of this section is supposed to demonstrate the performance of the estimator in a relatively difficult case when the right-hand side of the equation grows infinitely as the domain thickness tends to zero.

In this test, we consider a very simple geometry (see Figure 4 (left)), namely

$$
d_{\oplus, \ominus}= \pm \frac{d_{0}}{2}
$$

where $d_{0}>0$ is the given thickness of the domain, $\widehat{\Omega}=(0,1)$ and $\Omega=\left\{(x, y) \in \mathbb{R}^{2} \mid x \in \widehat{\Omega},-\frac{d_{0}}{2}<\right.$ $\left.y<\frac{d_{0}}{2}\right\}$. The considered problem reads

$$
\begin{aligned}
-\Delta u & =f & & \text { in } \Omega \\
u & =0 & & \text { at } x=0 \text { and } x=1, \\
\frac{\partial u}{\partial y} & = \pm F_{\oplus, \ominus} & & \text { at } y= \pm \frac{d_{0}}{2},
\end{aligned}
$$




Figure 4: (left) The domain geometry; (right) Convergence rate of the exact modelling-error and of the error majorant, $m=2$, the majorant is indicated by " $\circ$ ".
and the right-hand sides of the equation and of the boundary condition are computed using the exact solution

$$
u(x, y)=\sin (\pi x) \cdot \frac{y^{m}}{d_{0}^{m-1}} \quad(m=1,2, \ldots)
$$

The scaling factor $d_{0}^{m-1}$ makes this test essentially different from the previous one: while the Neumann boundary data $F_{\oplus, \ominus}$ remain of order $\mathcal{O}(1)$ as $d_{0} \rightarrow 0$, the right-hand side of the equation $f$ exhibits the behaviour $f \sim \mathcal{O}\left(d_{0}\right)+\mathcal{O}\left(\frac{1}{d_{0}}\right)$, i.e. unboundedly grows when $d_{0}$ tends to zero. The unbounded growth of $f$ may yield serious problems for an a-posteriori error estimator, as we are about to see. We also note that the constant $C_{\Omega}$ can be computed exactly in this example: $C_{\Omega}=\frac{1}{\pi}$ for all values of the thickness $d_{0}$.

First, we take $m=2$ and observe the convergence of the exact modelling-error in the energy norm and of the error majorant as $d_{0}$ tends to zero, see Figure 4 (right). As in the preceding example, the error majorant provides a reliable upper bound for the exact error at any values of the thickness $d_{0}$, both the exact error and the majorant demonstrate the optimal convergence rate $\mathcal{O}\left(d_{0}^{1 / 2}\right)$ and, moreover, the error majorant shows the asymptotic exactness in this case (the effectivity index $\frac{M}{\Pi \mid e \| ा}=1+\mathcal{O}\left(d_{0}\right)$, see Table 2, the column under " $m=2, M\left(\psi_{1}\right)$ "). However, if we set $m=3$, the second term of the majorant $M$ (i.e. $\|f-\widetilde{f}\|_{L_{2}(\Omega)}$, see (5.6)) becomes dominating and the whole estimator grows unboundedly, as can be seen in Figure 5 (left). The estimator becomes, of course, useless as it dramatically overestimates the exact error for small values of $d_{0}$. It is rather clear that the problem originates from the bad choice of the auxiliary function $\psi$ that is supposed to approximate $\frac{\partial u}{\partial y}$; for $m=3$ the derivative is quadratic and cannot be adequately represented by the linear function $\psi_{1}$.

The situation may be cured by invoking the quadratic function $\psi=\psi_{2}$ (see Remark 4.2), $\psi_{2}(x, y)=\psi_{1}(x, y)+\widehat{\eta}(x)\left(y^{2}-\frac{d_{0}^{2}}{4}\right)$, where $\widehat{\eta}$ is an arbitrary function from $L_{2}(\widehat{\Omega})$. The possibility to choose a suitable $\widehat{\eta}$ enables us to suppress the unbounded growth of $f$ in the $M_{2}$-term of the majorant and makes the majorant flexible enough to efficiently reproduce the behaviour of the exact error.

If we plug $\psi_{2}$ into the estimate (4.20), we obtain

$$
\begin{equation*}
\|\|u-\widehat{u}\|\|^{2} \leq M^{2}(\widehat{\eta}, \gamma) \quad \forall \widehat{\eta} \in L_{2}(\widehat{\Omega}), \forall \gamma>0 \tag{6.1}
\end{equation*}
$$



Figure 5: The case $m=3$ : (left) Divergence of the majorant $M\left(\psi_{1}\right)$ as $d_{0} \rightarrow 0$; (right) Convergence of the improved majorant $M\left(\psi_{2}\right)$.
where

$$
\begin{aligned}
& M^{2}(\widehat{\eta}, \gamma):=(1+\gamma)\left\|\psi_{2}\right\|_{L_{2}(\Omega)}^{2}+\left(1+\frac{1}{\gamma}\right) C_{\Omega}^{2}\left\|f-\tilde{f}+\frac{\partial \psi_{2}}{\partial y}\right\|_{L_{2}(\Omega)}^{2}= \\
& (1+\gamma) \int_{\Omega}\left(\psi_{1}(x, y)+\widehat{\eta}(x)\left(y^{2}-\frac{d_{0}^{2}}{4}\right)\right)^{2} d x d y+\left(1+\frac{1}{\gamma}\right) C_{\Omega}^{2} \int_{\Omega}(f(x, y)-\widetilde{f}(x)+\widehat{\eta}(x) \cdot 2 y)^{2} d x d y .
\end{aligned}
$$

Since estimate (6.1) is valid for any $\gamma>0$ and $\widehat{\eta}$ from $L_{2}(\widehat{\Omega})$, one can minimize the functional $M^{2}(\widehat{\eta}, \gamma)$ with respect to these parameters. In particular, one can set $\gamma=\gamma_{*}<1$ (the concrete value of $\gamma_{*}$ does not matter, as the numerical experiments show; we used the value $\gamma_{*}=0.5$ ) and find $\widehat{\eta}_{\text {min }}$ as the minimizer of $M^{2}\left(\widehat{\eta}, \gamma_{*}\right)$ over the space $S$ of piecewise-constant functions defined on some finite subdivision of $\widehat{\Omega}$ (obviously, $S \subset L_{2}(\widehat{\Omega})$ ). The minimization problem is just an $L_{2}$-projection onto the space of functions defined on $\widehat{\Omega}$ and amounts to the solution of a linear system with the diagonal matrix.

The properties of the improved majorant $M\left(\psi_{2}\right)=M_{1}\left(\psi_{2}\right)+C_{\Omega} M_{2}\left(\psi_{2}\right)$, where

$$
\begin{aligned}
& M_{1}\left(\psi_{2}\right):=\left\|\psi_{1}+\widehat{\eta}_{\min }\left(y^{2}-\frac{d_{0}^{2}}{4}\right)\right\|_{L_{2}(\Omega)} \\
& M_{2}\left(\psi_{2}\right):=\left\|f-\tilde{f}+\widehat{\eta}_{\min } \cdot 2 y\right\|_{L_{2}(\Omega)}
\end{aligned}
$$

can be observed in Figure 5 (right). We see that the improved majorant vanishes with the optimal rate $\mathcal{O}\left(d_{0}^{1 / 2}\right)$ as $d_{0} \rightarrow 0$, remains a reliable upper bound for the exact error at any values of the thickness $d_{0}$ and even demonstrates the asymptotic exactness with the effectivity index behaving like $1+\mathcal{O}\left(d_{0}\right)$ (see Table 2 ).

We may note that in the case of larger values of $m(m>3)$ the higher-degree function $\psi_{m-1}$ might be needed (see Remark 4.2); the function will contain several free parameters which are the functions from $L_{2}(\widehat{\Omega})$, and, hence, the minimization should be performed with respect to all of them. However, as this always remains a least-squares minimization problem, the total complexity for the moderate values of $m$ will not be greater than the complexity of solving the reduced problem. In general, if the right-hand side $f$ exhibits an unbounded growth for $d_{0} \rightarrow 0$ and no a-priori information on the behaviour of the exact solution is available, one has to choose the function $\psi$ in an adaptive way, i.e. starting with $\psi_{1}$ increase the polynomial degree of the function until the difference between two successive majorants $M\left(\psi_{n-1}\right)$ and $M\left(\psi_{n}\right)$ becomes small enough.

| $d_{0}^{-1}$ | $m=2, M\left(\psi_{1}\right)$ |  |  | $m=3, M\left(\psi_{2}\right)$ |  |  |
| :--- | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $\mid\\|e\\| \\|$ | $M$ | $\frac{M}{\\|e\\| \\|}$ | $\\|\|\|e\| \\|$ | $M$ | $\frac{M}{\\|e\\| \\|}$ |
| $10^{0}$ | 0.4405 | 0.9284 | 2.1074 | 0.2594 | 0.4187 | 1.6142 |
| $10^{1}$ | 0.1291 | 0.1461 | 1.1265 | 0.0751 | 0.0793 | 1.0562 |
| $10^{2}$ | 0.0408 | 0.0414 | 1.0127 | 0.0237 | 0.0239 | 1.0056 |
| $10^{3}$ | 0.0129 | 0.0130 | 1.0013 | 0.0075 | 0.0076 | 1.0006 |
| $10^{4}$ | 0.0041 | 0.0041 | 1.0001 | 0.0024 | 0.0024 | 1.0001 |

Table 2: Convergence of the exact modelling-error in the energy norm ( $\|\|e\|\|$ ) and of the error majorant ( $M$ ) as $d_{0} \rightarrow 0$; the results are rounded up to $10^{-4}$.

## 7 Conclusions

For the zero-order dimension reduction method, the new a-posteriori error estimator has been derived in a general geometrical setting of the problem and without any specific assumptions on the given data. In particular, the estimator reduces to the Babuška and Schwab estimator when the physical domain $\Omega$ is a plate with plane parallel faces and the equation has zero right-hand side. It has been demonstrated, both theoretically and numerically, that also in a more complicated case of a plate having constant thickness but non-plane faces and for a general right-hand side $f \in L_{\infty}(\Omega)$ the proposed estimator vanishes with the optimal rate $\mathcal{O}\left(d_{0}^{1 / 2}\right)$ as the plate thickness $d_{0}$ tends to zero. Since the estimator always provides an upper bound for the exact modelling-error, the latter convergence result can be considered as the generalization of the result on the convergence of the dimension reduction error proved in [11] (see also [2]) for the case of a plate with plane parallel faces and zero right-hand side $f$.

The presented estimator cannot, however, be considered as just a generalization of the explicit residual-type error estimator to the case of a more complicated geometry, coefficients and right-hand side. As numerical test 2 shows, in the problem with the right-hand side $f$ infinitely growing as the plate thickness tends to zero, some additional "degree of freedom" should be introduced into the estimator to suppress the unbounded growth of $f$. Thus, it seems that any error estimator that cannot be adjusted to the particular problem will fail in such a case. The proposed estimator has been shown sufficiently flexible to allow the modification necessary for capturing the behaviour of the exact error. The recovered efficiency of the estimator manifests itself in the asymptotics of the effectivity index $\frac{M}{\|e e\|}=1+\mathcal{O}\left(d_{0}\right)$ when $d_{0}$ tends to zero. We have to note that such an asymptotics may not always be observed, if the domain $\Omega$ has non-plane faces; however, even in the latter case, the effectivity index of the estimator remains stable (i.e. does not grow with the decreasing domainthickness) and stays at the acceptable level.

The computational cost of evaluating the presented error majorant is typically smaller than or, in the worst case, of the same order as the cost of solving the reduced, lower-dimensional problem. Finally, the numerical results show that the proposed estimator is capable of an accurate indication of the local error distribution and, hence, may be utilized not only for the verification of the dimensionally reduced model but also for its adaptive improvement.

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