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Adaptive regularization, linearization, and discretization and a posteriori error control for the two-phase Stefan problem*

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Abstract

We consider in this paper the time-dependent two-phase Stefan problem and derive a posteriori error estimates and adaptive strategies for its conforming spatial and backward Euler temporal discretizations. Regularization of the enthalpy–temperature function and iterative linearization of the arising systems of nonlinear algebraic equations are considered. Our estimators yield a guaranteed and fully computable upper bound on the dual norm of the residual, as well as on the $L^2(L^2)$ error of the temperature and the $L^2(H^{-1})$ error of the enthalpy. Moreover, they allow to distinguish the space, time, regularization, and linearization error components. An adaptive algorithm is proposed, which ensures computational savings through the online choice of a sufficient regularization parameter, a stopping criterion for the linearization iterations, local space mesh refinement, time step adjustment, and equilibration of the spatial and temporal errors. We also prove the efficiency of our estimate. Our analysis is quite general and is not focused on a specific choice of the space discretization and of the linearization. As an example, we apply it to the vertex-centered finite volume (finite element with mass lumping and quadrature) and Newton methods. Numerical results illustrate the effectiveness of our estimates and the performance of the adaptive algorithm.

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1 Introduction

The two-phase Stefan problem models a phase change process which is governed by the Fourier law, c.f. Friedman [22]. The two phases, typically solid and liquid, are separated by a moving interface, whose motion is governed by the so-called Stefan condition. Let $\Omega \subset \mathbb{R}^d$, $d \in \{2, 3\}$, be an open bounded polygonal or polyhedral domain, not necessarily convex, and let $T > 0$. The mathematical statement of the problem is as follows: given an initial enthalpy u_0 and a source function f , find the enthalpy u such that

$$\partial_t u - \nabla \cdot (\nabla \beta(u)) = f \quad \text{in } \Omega \times (0, T), \quad (1.1a)$$

$$u(\cdot, 0) = u_0 \quad \text{in } \Omega, \quad (1.1b)$$

$$\beta(u) = 0 \quad \text{on } \partial\Omega \times (0, T). \quad (1.1c)$$

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For the sake of simplicity, we assume that u is normalized so that the (specific) enthalpies of the two phases are 0 and 1, respectively, and only consider the homogeneous Dirichlet boundary condition (1.1c). The temperature $\beta(u)$ is expressed as a function of the enthalpy u . In what follows, we assume that $\beta(\cdot)$ is a nondecreasing Lipschitz continuous function which vanishes in the interval $(0, 1)$. The latter condition reflects the latent heat in the phase change.

The numerical analysis of the Stefan problem has been considered in several works. A finite difference method for the multi-dimensional Stefan problem is discussed in Meyer [32]. The author presents a convergent numerical scheme which is the implicit analogue of the method of Kamenomostskaja [27]. In Ciavaldini [12], the numerical approach is based on finite elements of first order. The author describes the different schemes used and the nature of their convergence. Elliott [16] presents a finite element approximation of an elliptic variational inequality deduced from a semi-discretization in time of the weak formulation of the two-phase Stefan problem. Discretization schemes for regularized versions of the Stefan problem based on piecewise linear Lagrange finite elements in space and backward differencing in time are presented by Jerome and Rose [25]. Jäger and Kačur [24] use the enthalpy formulation and a variational technique to analyze the convergence of linearized semi-discrete-in-time and fully discrete schemes for nonlinear degenerate parabolic systems of porous medium type. In Amiez and Gremaud [2], a numerical scheme based on the approximation of the enthalpy formulation by semi-implicit finite differences in time combined with continuous piecewise linear finite elements in space is presented. Nocketto [33] employs the regularization technique to derive a priori error estimates in $L^2(0, T; L^2(\Omega))$ for the enthalpy and temperature errors of respectively one half and first order for an implicit finite element scheme, under suitable conditions on the data and relation of the space and time steps.

More recently, attention has been paid to finite volume methods which can be used on a large variety of meshes. The framework of semigroup theory has been used by Baughman and Walkington [4] for the study of the co-volume method, which is a special instance of the finite volume method. The analysis predicts one half order rates of convergence for approximate solutions of the enthalpy in $L^\infty(0, T; H^{-1}(\Omega))$ and of the temperature in $L^2(0, T; L^2(\Omega))$. In Eymard et al. [21] the authors give a convergence proof in the case that a finite volume scheme on a general mesh is used for the space discretization. Weak-* convergence for u in L^∞ and strong convergence for $\beta(u)$ in L^2 is shown by means of a priori estimates in L^∞ and use of the Kolmogorov theorem on relative compactness of subsets of L^2 . Half order error estimates via regularization have also been obtained in Pop et al. [42], whereas extensions to degenerate hyperbolic–parabolic equations can be found in Andreianov et al. [3].

A technique often used in various numerical approaches, c.f. Nocketto [33], Nocketto and Verdi [37], Picasso [40], Beckett et al. [6], or Pop et al. [42], employs a regularization of the nonsmooth and nonstrictly increasing function β by a smooth and strictly increasing one. This allows, in particular, to use the Newton method for the solution of the arising system of nonlinear algebraic equations, albeit its use without regularization has been advocated in Wheeler [52] or in Baughman and Walkington [4] and studied in Kelley and Rulla [28]. Alternative approaches such as transformation of dependent variables of Čermák and Zlámal [10] have also been proposed.

An inevitable tool in practical simulations seems to be an a posteriori error estimate-driven adaptive mesh refinement. One of the first works on a posteriori error estimates for the steady Stefan problem is that of Picasso [40]. Therein, the author derives estimates based on the residual equation for a two-dimensional regularized Stefan problem and proposes a space adaptive finite element algorithm. A posteriori indicators for unsteady phase change problems were derived by Nocketto et al. in [34, 35], together with an adaptive algorithm which equilibrates space and time discretization errors. Many other adaptive refinement algorithms such as that of Beckett et al. [6] have also been proposed. Rigorous a posteriori error estimates for nonlinear parabolic problems seem much less developed. In nondegenerate cases, Verfürth [46, 47] was able to obtain an estimator which is both reliable and efficient. A pioneering contribution for degenerate parabolic problems has been obtained by Nocketto et al. in [36]. Therein, $L^\infty(0, T; H^{-1}(\Omega))$ estimates for the error in the enthalpy and $L^2(0, T; L^2(\Omega))$ estimates for the error in the temperature are obtained. The approach is based on the relation of these errors to the residual of (1.1a) obtained through the corresponding dual partial differential equation and subsequent use of the Galerkin

orthogonality of the finite element discretization. Recently, rigorous a posteriori error analysis in a space–time dual norm, including some degenerate cases, was given in [14].

The aim of this paper is to derive fully computable a posteriori error estimates and adaptive strategies for the two-phase Stefan problem (1.1) for conforming spatial discretization schemes such as the finite element, co-volume, or vertex-centered finite volume ones with backward Euler time stepping. As in Nochetto et al. [36], our approach is based on the dual norm of the residual. However, we proceed differently in order to have a fully and easily computable estimates not featuring any undetermined constants. This is achieved by introducing $\mathbf{H}(\text{div}; \Omega)$ -conforming and locally conservative flux reconstructions following Prager–Synge [43], Ladevèze [29], Destuynder and Métivet [13], Luce and Wohlmuth [30], Braess and Schöberl [8], Repin [45], and [50, 18, 15, 14, 19], see also the references therein.

In Section 2 we give a weak formulation, introduce a regularized problem with a regularization parameter $\epsilon > 0$, and fix the notation for temporal and spatial meshes. In Section 3, we identify the residual and its dual norm and we derive an a posteriori error estimate on this problem-dependent error measure. We next split this estimate into estimators characterizing the space, time, regularization, linearization, and quadrature errors.

Section 4 subsequently presents a criterion for the choice of the regularization parameter ϵ and a stopping criterion for an iterative linearization such as the Newton method. The former is designed to facilitate the treatment of the degeneracy while not spoiling the accuracy, whereas the latter is designed to avoid performing an excessive number of nonlinear solver iterations. These criteria are inspired mainly from [26, 15, 19]. We then propose an adaptive algorithm which uses these criteria while simultaneously performing the usual local mesh refinement and equilibration of the spatial and temporal errors. This algorithm is inspired from [34, 35, 40, 36, 6] and from the work [18, 23, 14, 19]. We conclude Section 4 by proving that, under these criteria, our estimators are also efficient while representing a lower bound for the dual norm of the residual.

In Section 5, we show how to bound the $L^2(0, T; H^{-1}(\Omega))$ -type error in the enthalpy and $L^2(0, T; L^2(\Omega))$ -type error in the temperature by the above dual norm of the residual. We in particular focus on the use of the Gronwall lemma with as small overestimation as possible and no appearance of the exponential term e^T elsewhere than in the approximation of the initial condition. Guaranteed and fully computable a posteriori error estimates on these natural norms immediately follow.

Section 6 presents the application of all these developments to the vertex-centered finite volume (or, equivalently, finite element with mass lumping and numerical quadrature) discretization in space, backward Euler discretization in time, and Newton linearization. Illustrative numerical results fill up Section 7 and, finally, Appendix A collects the more involved proofs of the various theorems of the paper.

2 Continuous and discrete settings

This section fixes the basic continuous and discrete settings. More precisely, Section 2.1 presents the continuous problem and the regularization, whereas the basic assumptions on the discretization are introduced in Section 2.2.

2.1 Continuous setting

2.1.1 The continuous problem

The starting point for our a posteriori analysis is the weak form of problem (1.1). To give it, we need to introduce the assumptions on the data and set up some notation. We suppose that: (i) the enthalpy–temperature function $\beta : \mathbb{R} \rightarrow \mathbb{R}$ is a Lipschitz continuous function such that

$$\beta(s) = 0 \quad \text{in } (0, 1),$$

β is strictly increasing in \mathbb{R}^- and $\mathbb{R}^+ \setminus (0, 1)$, and there exist $c, C > 0$ such that, for all $s \in \mathbb{R} \setminus (0, 1)$, $\text{sign}(s)\beta(s) \geq c|s| - C$, see Figure 1; the Lipschitz constant of β is denoted by L_β ; (ii) the initial enthalpy u_0 is such that $u_0 \in L^2(\Omega)$; (iii) the source term is such that $f \in L^2(0, T; L^2(\Omega))$.

We will repeatedly use throughout the paper the two following spaces:

$$X := L^2(0, T; H_0^1(\Omega)), \quad Z := H^1(0, T; H^{-1}(\Omega)). \quad (2.1)$$

We will also need the dual space X' of X ,

$$X' = L^2(0, T; H^{-1}(\Omega)),$$

and equip the space X with the norm

$$\|\varphi\|_X := \left\{ \int_0^T \|\nabla \varphi(\cdot, t)\|_{L^2(\Omega)}^2 dt \right\}^{\frac{1}{2}}.$$

We denote by $\langle \cdot, \cdot \rangle$ the duality pairing between $H^{-1}(\Omega)$ and $H_0^1(\Omega)$, while $(\cdot, \cdot)_S$ is the usual scalar product in $L^2(S)$ or $[L^2(S)]^d$, with the subscript omitted when $S = \Omega$.

The weak formulation of problem (1.1) can now be stated. It reads: find

$$u \in Z \quad \text{with} \quad \beta(u) \in X \quad (2.2a)$$

such that

$$u(\cdot, 0) = u_0 \quad \text{in } \Omega \quad (2.2b)$$

and, for a.e. $s \in (0, T)$,

$$\langle \partial_t u(\cdot, s), \varphi \rangle + (\nabla \beta(u(\cdot, s)), \nabla \varphi) = (f(\cdot, s), \varphi) \quad \forall \varphi \in H_0^1(\Omega). \quad (2.2c)$$

Existence and uniqueness of the solution to this problem are known [22, 1, 38, 7].

2.1.2 A regularization

An important feature of the problem (2.2) is that, as a result of the assumptions on β , the normal component of the temperature flux $-\nabla \beta(u)$ may jump across the interface

$$I(t) := \{\mathbf{x} \in \Omega : \beta(u)(\mathbf{x}, t) = 0\}.$$

This fact may hinder both the design and the convergence analysis of a discretization method. Additionally, the lack of smoothness in the dependency of the solution on the problem data can severely affect the convergence of nonlinear iterations. A possible and often employed approach [33, 37, 40, 6] to overcome these difficulties consists in regularizing the problem (2.2) by replacing the function β by a smooth, strictly increasing regularized function $\beta_\epsilon \in C^1(\mathbb{R})$, $\beta'_\epsilon \geq \epsilon$, for a parameter $\epsilon > 0$; see Figure 1 for an example. The regularized problem reads as follows: find

$$u^\epsilon \in Z \quad \text{with} \quad \beta_\epsilon(u^\epsilon) \in X \quad (2.3a)$$

such that

$$u^\epsilon(\cdot, 0) = \beta_\epsilon^{-1}(\beta(u_0)) \quad \text{in } \Omega, \quad (2.3b)$$

and, for a.e. $s \in (0, T)$,

$$\langle \partial_t u^\epsilon(\cdot, s), \varphi \rangle + (\nabla \beta_\epsilon(u^\epsilon(\cdot, s)), \nabla \varphi) = (f(\cdot, s), \varphi) \quad \forall \varphi \in H_0^1(\Omega). \quad (2.3c)$$

2.2 Discrete setting

We describe here the basic discrete setting that will be sufficient for the developments of Sections 3–5. Further details are given in Section 6.

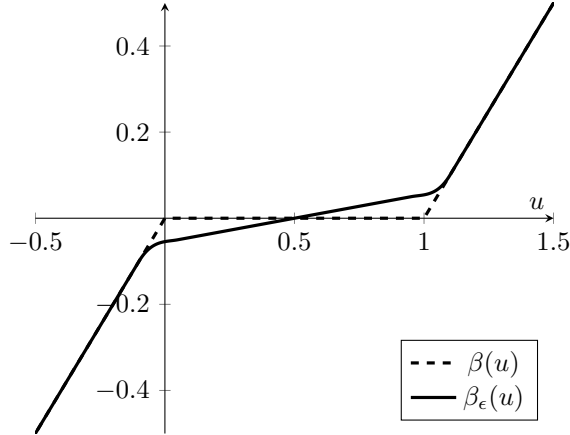


Figure 1: An example of a function β and a regularization β_ϵ

2.2.1 Time mesh

Our focus is on first-order time discretizations based on the backward Euler scheme. Let $\{\tau^n\}_{1 \leq n \leq N}$ denote a sequence of positive real numbers corresponding to the discrete time steps such that $T = \sum_{n=1}^N \tau^n$. We let $t^0 := 0$ and, for $1 \leq n \leq N$, we introduce the discrete times $t^n := \sum_{i=1}^n \tau^i$ and the time intervals $I_n := (t^{n-1}, t^n)$.

2.2.2 Space meshes

Let $\{\mathcal{K}^n\}_{0 \leq n \leq N}$ denote a family of matching simplicial meshes of the space domain Ω . The initial mesh \mathcal{K}^0 is used to approximate the initial condition u_0 , while \mathcal{K}^n is used to march in time from t^{n-1} to t^n . The meshes can be refined or coarsened as time evolves. For the developments of Section 4.3 below, we are led to suppose that \mathcal{K}^n , $1 \leq n \leq N$, is obtained from \mathcal{K}^{n-1} by refining (a limited number of times) some elements and coarsening (a limited number of times) some other ones. We denote by $\mathcal{K}^{n-1,n}$ the coarsest common submesh (overlay) of both \mathcal{K}^n and \mathcal{K}^{n-1} and, once again for the developments of Section 4.3, suppose that the meshes $\{\mathcal{K}^{n-1,n}\}_{1 \leq n \leq N}$ are shape-regular in the sense that there exists a constant $\kappa_{\mathcal{K}} > 0$ such that

$$\min_{K \in \mathcal{K}^{n-1,n}} \frac{\rho_K}{h_K} \geq \kappa_{\mathcal{K}}$$

for all $1 \leq n \leq N$, where ρ_K denotes the diameter of the largest ball inscribed in the element K and h_K the diameter of K . For $0 \leq n \leq N$, we denote by Π_0^n the L^2 -orthogonal projection onto the space of piecewise constant functions on \mathcal{K}^n .

For $0 \leq n \leq N$, let \mathcal{F}^n denote the set of mesh faces. Boundary faces are collected in the set $\mathcal{F}^{b,n} := \{F \in \mathcal{F}^n; F \subset \partial\Omega\}$ and we let $\mathcal{F}^{i,n} := \mathcal{F}^n \setminus \mathcal{F}^{b,n}$. For a given face $F \in \mathcal{F}^{i,n}$ we fix an arbitrary orientation and denote the corresponding unit normal vector by \mathbf{n}_F ; for $F \in \mathcal{F}^{b,n}$, \mathbf{n}_F coincides with the exterior unit normal \mathbf{n}_Ω of Ω . A similar notation for the faces $\mathcal{F}^{n-1,n}$ of the meshes $\mathcal{K}^{n-1,n}$ will also be used.

3 An a posteriori error estimate for the dual norm of the residual

In this section we derive an a posteriori estimate for the error measured by the dual norm of the residual that we first identify. We then give a basic estimate that we subsequently refine to distinguish the space, time, linearization, regularization, and quadrature errors.

3.1 Dual norm of the residual

As in Picasso [40] or Nochetto et al. [36], our key for deriving a posteriori error estimates for the Stefan problem (1.1) will be the residual and its dual norm. Recall that u denotes the weak solution of the Stefan problem given by (2.2) and the definition of the space X (2.1). Let $u_{h\tau} \in Z$ such that $\beta(u_{h\tau}) \in X$ be arbitrary. In practice, $u_{h\tau}$ will be the result of the numerical simulation. We define the residual $\mathcal{R}(u_{h\tau}) \in X'$ such that

$$\langle \mathcal{R}(u_{h\tau}), \varphi \rangle_{X', X} := \int_0^T \{ \langle \partial_t(u - u_{h\tau}), \varphi \rangle + (\nabla \beta(u) - \nabla \beta(u_{h\tau}), \nabla \varphi) \} (s) ds, \quad \varphi \in X. \quad (3.1)$$

Using (2.2c), we can infer the following alternative expression for (3.1):

$$\langle \mathcal{R}(u_{h\tau}), \varphi \rangle_{X', X} = \int_0^T \{ (f, \varphi) - \langle \partial_t u_{h\tau}, \varphi \rangle - (\nabla \beta(u_{h\tau}), \nabla \varphi) \} (s) ds, \quad \varphi \in X.$$

The norm of the residual in the dual space X' is then given by

$$\|\mathcal{R}(u_{h\tau})\|_{X'} := \sup_{\varphi \in X, \|\varphi\|_X=1} \langle \mathcal{R}(u_{h\tau}), \varphi \rangle_{X', X}. \quad (3.2)$$

The key problem-specific measure of the distance between $u_{h\tau}$ and u that we will use in this paper is given by

$$\|\mathcal{R}(u_{h\tau})\|_{X'} + \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}. \quad (3.3)$$

It follows from (2.2) that the measure (3.3) is zero if and only if the function $u_{h\tau}$ coincides with the exact solution u . As we shall see below in Section 5, it in fact controls the energy error between u and $u_{h\tau}$ and $\beta(u)$ and $\beta(u_{h\tau})$. Moreover, this quantity can be easily bounded in terms of error estimators based on $\mathbf{H}(\text{div}; \Omega)$ -conforming flux reconstructions for piecewise affine-in-time $u_{h\tau}$ that we show next.

3.2 General assumptions

In order to proceed with the analysis further, without the necessity to specify at this point any details on how the approximate solution $u_{h\tau}$ was obtained, we are lead to make the following assumption. It requires Z - and X - conformity and $u_{h\tau}$ to be piecewise affine and continuous in time on the time mesh $\{I_n\}_{1 \leq n \leq N}$ of Section 2.2.1:

Assumption 3.1 (Approximate solution). *The function $u_{h\tau}$ is such that*

$$\begin{aligned} u_{h\tau} \in Z, \quad \partial_t u_{h\tau} \in L^2(0, T; L^2(\Omega)), \quad \beta(u_{h\tau}) \in X, \\ u_{h\tau}|_{I_n} \text{ is affine in time on } I_n \quad \forall 1 \leq n \leq N. \end{aligned}$$

Note that, consequently, the function $u_{h\tau}$ is uniquely determined by the $N + 1$ functions $u_h^n := u_{h\tau}(\cdot, t^n)$, $0 \leq n \leq N$, and $\partial_t u_{h\tau}^n := \partial_t u_{h\tau}|_{I_n} \equiv (u_h^n - u_h^{n-1})/\tau^n$, $1 \leq n \leq N$. We will also employ the abridged notation $u_{h\tau}^n$ for $u_{h\tau}|_{I_n}$.

The second assumption that we make is the existence of a piecewise constant-in-time $\mathbf{H}(\text{div}; \Omega)$ -conforming flux reconstruction $\mathbf{t}_{h\tau}$, locally conservative on the meshes \mathcal{K}^n of Section 2.2.2. Let us first denote by \hat{f} the piecewise constant-in-time function given by the time-mean values of the source function f on the intervals I_n , $1 \leq n \leq N$.

Assumption 3.2 (Equilibrated flux reconstruction). *For all $1 \leq n \leq N$, there exists a vector field $\mathbf{t}_h^n \in \mathbf{H}(\text{div}; \Omega)$ such that*

$$(\nabla \cdot \mathbf{t}_h^n, 1)_K = (\hat{f}^n, 1)_K - (\partial_t u_{h\tau}^n, 1)_K \quad \forall K \in \mathcal{K}^n.$$

We denote by $\mathbf{t}_{h\tau}$ the space-time function such that $\mathbf{t}_{h\tau}|_{I_n} := \mathbf{t}_h^n$ for all $1 \leq n \leq N$.

In Section 6 below, we show how to construct an equilibrated flux reconstruction $\mathbf{t}_{h\tau}$ in the context of vertex-centered finite volume (finite element with mass lumping and quadrature) spatial discretization.

3.3 A basic a posteriori error estimate

We now give an a posteriori error estimate in the general setting of Assumptions 3.1 and 3.2. Note that the regularization of Section 2.1.2 is not used at the present stage.

We will estimate the error measure (3.3) by the local residual expressed with the flux $\mathbf{t}_{h\tau}$ and by the difference of $\mathbf{t}_{h\tau}$ and the temperature flux, in the spirit of [43, 29, 13, 30, 8, 45] and [50, 18, 15, 14, 19]. More specifically, for $1 \leq n \leq N$, \mathbf{t}_h^n as in Assumption 3.2, and $K \in \mathcal{K}^n$, we define the *residual estimator* $\eta_{\mathbb{R},K}^n$ and the *flux estimator* $\eta_{\mathbb{F},K}^n$ as follows:

$$\eta_{\mathbb{R},K}^n := C_{\mathbb{P},K} h_K \|\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n\|_{L^2(K)}, \quad (3.4a)$$

$$\eta_{\mathbb{F},K}^n(t) := \|\mathbf{t}_h^n + \nabla \beta(u_{h\tau}(\cdot, t))\|_{L^2(K)} \quad t \in I_n. \quad (3.4b)$$

Here, $C_{\mathbb{P},K}$ is the constant from the Poincaré inequality

$$\|\varphi - \Pi_0^n \varphi\|_{L^2(K)} \leq C_{\mathbb{P},K} h_K \|\nabla \varphi\|_{L^2(K)} \quad \forall \varphi \in H^1(K). \quad (3.5)$$

There holds $C_{\mathbb{P},K} = 1/\pi$ as the simplices K are convex, see [5, 39]. Finally, we define the *initial condition estimator* by

$$\eta_{\mathbb{IC}} := \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}. \quad (3.6)$$

We then have:

Theorem 3.3 (A posteriori estimate for the error measure (3.3)). *Let u be the weak solution given by (2.2) and let $u_{h\tau}$ and $\mathbf{t}_{h\tau}$ fulfill Assumptions 3.1 and 3.2, respectively. Then, there holds*

$$\|\mathcal{R}(u_{h\tau})\|_{X'} + \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)} \leq \eta + \eta_{\mathbb{IC}}, \quad (3.7)$$

where

$$\eta := \left\{ \sum_{n=1}^N \int_{I_n} \sum_{K \in \mathcal{K}^n} (\eta_{\mathbb{R},K}^n + \eta_{\mathbb{F},K}^n(t))^2 dt \right\}^{\frac{1}{2}} + \|f - \hat{f}\|_{X'}. \quad (3.8)$$

Proof. Let $\varphi \in X$ with $\|\varphi\|_X = 1$ be given. Then there holds, adding and subtracting $(\mathbf{t}_{h\tau}, \nabla \varphi)$ and using Green's theorem,

$$\begin{aligned} \langle \mathcal{R}(u_{h\tau}), \varphi \rangle_{X', X} &= \int_0^T \{ (f - \partial_t u_{h\tau} - \nabla \cdot \mathbf{t}_{h\tau}, \varphi) - (\mathbf{t}_{h\tau} + \nabla \beta(u_{h\tau}), \nabla \varphi) \} (s) ds \\ &= \int_0^T \left\{ (f - \hat{f}, \varphi) + (\hat{f} - \partial_t u_{h\tau} - \nabla \cdot \mathbf{t}_{h\tau}, \varphi) - (\mathbf{t}_{h\tau} + \nabla \beta(u_{h\tau}), \nabla \varphi) \right\} (s) ds \\ &=: \mathfrak{T}_1 + \mathfrak{T}_2 + \mathfrak{T}_3. \end{aligned}$$

For the first term we infer $\mathfrak{T}_1 \leq \|f - \hat{f}\|_{X'} \|\varphi\|_X = \|f - \hat{f}\|_{X'}$. The second term can be rewritten as follows:

$$\mathfrak{T}_2 = \sum_{n=1}^N \int_{I_n} (\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n, \varphi)(s) ds.$$

For all $1 \leq n \leq N$ and $t \in I_n$, there holds (the dependence of φ on the time variable is omitted for brevity),

$$\begin{aligned} (\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n, \varphi) &= \sum_{K \in \mathcal{K}^n} (\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n, \varphi)_K \\ &= \sum_{K \in \mathcal{K}^n} (\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n, \varphi - \Pi_0^n \varphi)_K \\ &\leq \sum_{K \in \mathcal{K}^n} \|\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n\|_{L^2(K)} \|\varphi - \Pi_0^n \varphi\|_{L^2(K)} \\ &\leq \sum_{K \in \mathcal{K}^n} C_{\mathbb{P},K} h_K \|\hat{f}^n - \partial_t u_{h\tau}^n - \nabla \cdot \mathbf{t}_h^n\|_{L^2(K)} \|\nabla \varphi\|_{L^2(K)} \\ &= \sum_{K \in \mathcal{K}^n} \eta_{\mathbb{R},K}^n \|\nabla \varphi\|_{L^2(K)}, \end{aligned}$$

where we have used the regularity of the arguments, Assumption 3.2, the Cauchy–Schwarz inequality, and the Poincaré inequality (3.5). For the third term, an application of the Cauchy–Schwarz inequality yields

$$\mathfrak{T}_3 \leq \sum_{n=1}^N \int_{I_n} \sum_{K \in \mathcal{K}^n} \eta_{F,K}^n \|\nabla \varphi\|_{L^2(K)}(s) ds.$$

Collecting the above estimates, using the definition (3.2) of the dual norm of the residual, and using the Cauchy–Schwarz inequality yields (3.7). \square

3.4 An a posteriori error estimate distinguishing the space, time, regularization, linearization, and quadrature errors

Our next goal is to distinguish the different error components. This is an instrumental step to design an adaptive algorithm where the time step, the space mesh, the regularization parameter, and the stopping criterion for the linearization iterations are chosen optimally. We start by localizing in time the error measure introduced in Section 3.1. For $1 \leq n \leq N$, we let

$$X_n := L^2(I_n; H_0^1(\Omega)), \quad Z_n := H^1(I_n; H^{-1}(\Omega)).$$

We localize in time the dual norm of the residual (3.2) by setting

$$\|\mathcal{R}(u_{h\tau})\|_{X'_n} := \sup_{\varphi \in X_n, \|\varphi\|_{X_n}=1} \int_{I_n} \{ \langle \partial_t(u - u_{h\tau}), \varphi \rangle + (\nabla \beta(u) - \nabla \beta(u_{h\tau}), \nabla \varphi) \} (s) ds. \quad (3.9)$$

Note that, consequently,

$$\|\mathcal{R}(u_{h\tau})\|_{X'}^2 = \sum_{n=1}^N \|\mathcal{R}(u_{h\tau})\|_{X'_n}^2$$

for any $u_{h\tau} \in Z$ with $\beta(u_{h\tau}) \in X$.

Suppose now that we are marching in time from time t^{n-1} to time t^n with a given time step τ^n , starting from the approximation u_h^{n-1} . We also suppose that the regularization of Section 2.1.2 has been used for a given value of the parameter ϵ , and that we are on the k -th step of some iterative linearization algorithm. We denote by $u_h^{n,\epsilon,k}$ the approximation of the solution u at time t^n and prescribe the space–time function $u_{h\tau}^{n,\epsilon,k}$ by the value u_h^{n-1} at time t^{n-1} , by the value $u_h^{n,\epsilon,k}$ at time t^n , and by affine behavior in time on I_n , i.e.,

$$u_{h\tau}^{n,\epsilon,k}(\cdot, t) = (1 - \rho(t))u_h^{n-1} + \rho(t)u_h^{n,\epsilon,k}, \quad \rho(t) := \frac{t - t^{n-1}}{\tau^n}. \quad (3.10)$$

We summarize our general requirements in the following:

Assumption 3.4 (Adaptive setting). *For all $1 \leq n \leq N$, a regularization parameter $\epsilon \geq 0$, and a linearization step $k \geq 1$:*

(i) $u_{h\tau}^{n,\epsilon,k}$ is the approximate solution given by (3.10), $u_{h\tau}^{n,\epsilon,k} \in Z_n$ with $\partial_t u_{h\tau}^{n,\epsilon,k} \in L^2(I_n; L^2(\Omega))$ and $\beta(u_{h\tau}^{n,\epsilon,k}) \in X_n$;

(ii) there exists an equilibrated flux $\mathbf{t}_h^{n,\epsilon,k} \in \mathbf{H}(\text{div}; \Omega)$ such that

$$(\nabla \cdot \mathbf{t}_h^{n,\epsilon,k}, \mathbf{1})_K = (\hat{f}^n, \mathbf{1})_K - (\partial_t u_{h\tau}^{n,\epsilon,k}, \mathbf{1})_K \quad \forall K \in \mathcal{K}^n; \quad (3.11)$$

(iii) $\mathbf{I}_h^{n,\epsilon,k} \in [L^2(\Omega)]^d$ is the available approximation of the flux $\nabla \beta_\epsilon(u(\cdot, t^n))$;

(iv) Π^n is an operator used for interpolatory numerical integration.

An example of the approximate solution $u_{h\tau}^{n,\epsilon,k}$, the linearized flux $\mathbf{I}_h^{n,\epsilon,k}$, and the operator Π^n in the context of the implicit vertex-centered finite volume discretization and Newton linearization is provided in Section 6.3 below.

Proceeding as in Theorem 3.3, it is immediately inferred

$$\|\mathcal{R}(u_{h\tau}^{n,\epsilon,k})\|_{X'_n} \leq \left\{ \int_{I_n} \sum_{K \in \mathcal{K}^n} \left(\eta_{\mathbf{R},K}^{n,\epsilon,k} + \eta_{\mathbf{F},K}^{n,\epsilon,k}(t) \right)^2 dt \right\}^{\frac{1}{2}} + \|f - \hat{f}\|_{X'_n}, \quad (3.12)$$

where

$$\begin{aligned} \eta_{\mathbf{R},K}^{n,\epsilon,k} &:= C_{\mathbf{P},K} h_K \|\hat{f}^n - \partial_t u_{h\tau}^{n,\epsilon,k} - \nabla \cdot \mathbf{t}_h^{n,\epsilon,k}\|_{L^2(K)}, \\ \eta_{\mathbf{F},K}^{n,\epsilon,k}(t) &:= \|\mathbf{t}_h^{n,\epsilon,k} + \nabla \beta(u_{h\tau}^{n,\epsilon,k}(\cdot, t))\|_{L^2(K)}, \quad t \in I_n. \end{aligned}$$

For all $K \in \mathcal{K}^n$, we next define the local *spatial*, *temporal*, *quadrature*, *regularization*, and *linearization estimators* as follows:

$$\eta_{\text{sp},K}^{n,\epsilon,k} := \eta_{\mathbf{R},K}^{n,\epsilon,k} + \|\mathbf{I}_h^{n,\epsilon,k} + \mathbf{t}_h^{n,\epsilon,k}\|_{L^2(K)}, \quad (3.13a)$$

$$\eta_{\text{tm},K}^{n,\epsilon,k}(t) := \|\nabla(\Pi^n \beta(u_{h\tau}^{n,\epsilon,k}(\cdot, t))) - \nabla(\Pi^n \beta(u_h^{n,\epsilon,k}))\|_{L^2(K)}, \quad t \in I_n, \quad (3.13b)$$

$$\eta_{\text{qd},K}^{n,\epsilon,k}(t) := \|\nabla(\beta(u_{h\tau}^{n,\epsilon,k}(\cdot, t))) - \nabla(\Pi^n \beta(u_{h\tau}^{n,\epsilon,k}(\cdot, t)))\|_{L^2(K)}, \quad t \in I_n, \quad (3.13c)$$

$$\eta_{\text{reg},K}^{n,\epsilon,k} := \|\nabla(\Pi^n \beta(u_h^{n,\epsilon,k})) - \nabla(\Pi^n \beta_\epsilon(u_h^{n,\epsilon,k}))\|_{L^2(K)}, \quad (3.13d)$$

$$\eta_{\text{lin},K}^{n,\epsilon,k} := \|\nabla(\Pi^n \beta_\epsilon(u_h^{n,\epsilon,k})) - \mathbf{I}_h^{n,\epsilon,k}\|_{L^2(K)}. \quad (3.13e)$$

Global versions of these estimators are given by,

$$(\eta_{\text{sp}}^{n,\epsilon,k})^2 := \tau^n \sum_{K \in \mathcal{K}^n} \left(\eta_{\text{sp},K}^{n,\epsilon,k} \right)^2, \quad (3.14a)$$

$$(\eta_{\text{tm}}^{n,\epsilon,k})^2 := \int_{I_n} \sum_{K \in \mathcal{K}^n} \left(\eta_{\text{tm},K}^{n,\epsilon,k}(t) \right)^2 dt, \quad (3.14b)$$

$$(\eta_{\text{qd}}^{n,\epsilon,k})^2 := \int_{I_n} \sum_{K \in \mathcal{K}^n} \left(\eta_{\text{qd},K}^{n,\epsilon,k}(t) \right)^2 dt, \quad (3.14c)$$

$$(\eta_{\text{reg}}^{n,\epsilon,k})^2 := \tau^n \sum_{K \in \mathcal{K}^n} \left(\eta_{\text{reg},K}^{n,\epsilon,k} \right)^2, \quad (3.14d)$$

$$(\eta_{\text{lin}}^{n,\epsilon,k})^2 := \tau^n \sum_{K \in \mathcal{K}^n} \left(\eta_{\text{lin},K}^{n,\epsilon,k} \right)^2. \quad (3.14e)$$

Using the inequality (3.12) followed by the triangle inequality we obtain the following estimate:

Corollary 3.5 (Distinguishing the space, time, quadrature, regularization, linearization, and data oscillation errors). *Let u be the weak solution given by (2.2), let $1 \leq n \leq N$, $\epsilon \geq 0$, and $k \geq 1$, and let $u_{h\tau}^{n,\epsilon,k}$, $\mathbf{t}_h^{n,\epsilon,k}$, $\mathbf{I}_h^{n,\epsilon,k}$, and Π^n be as described in Assumption 3.4. Then there holds*

$$\|\mathcal{R}(u_{h\tau}^{n,\epsilon,k})\|_{X'_n} \leq \eta_{\text{sp}}^{n,\epsilon,k} + \eta_{\text{tm}}^{n,\epsilon,k} + \eta_{\text{qd}}^{n,\epsilon,k} + \eta_{\text{reg}}^{n,\epsilon,k} + \eta_{\text{lin}}^{n,\epsilon,k} + \|f - \hat{f}\|_{X'_n}.$$

Remark 3.6 (Time oscillation of the source term). *The error due to the time oscillation of the source term $\|f - \hat{f}\|_{X'_n}$ is zero provided that the source function f is piecewise constant in time.*

4 Balancing and stopping criteria, adaptive algorithm, and efficiency

The individual error component estimators of Corollary 3.5 are used in this section to define adaptive criteria to stop the iterative linearizations, to select the value of the regularization parameter ϵ , to locally adapt the quadrature rule, to adjust the time step, and to select the mesh

elements to refine/derefine. These criteria are incorporated in a fully adaptive algorithm detailed in Section 4.2. Finally, in Section 4.3 we show the efficiency of our estimators when the adaptive balancing and stopping criteria are used.

4.1 Balancing and stopping criteria

Following [26, 15, 19], this section introduces stopping criteria for the iterative algorithms based on the estimators of Corollary 3.5. The goal is to stop the iterations as soon as the corresponding error component no longer affects significantly the overall error. We assume in what follows that we are marching in time from time t^{n-1} to time t^n . Let three user-given parameters $\Gamma_{\text{lin}}, \Gamma_{\text{reg}}, \Gamma_{\text{qd}} \in (0, 1)$ be given. The criteria are:

- (i) *Linearization.* The linearization iteration is pursued until step k_n such that

$$\eta_{\text{lin}}^{n,\epsilon,k_n} \leq \Gamma_{\text{lin}} \left(\eta_{\text{sp}}^{n,\epsilon,k_n} + \eta_{\text{tm}}^{n,\epsilon,k_n} + \eta_{\text{qd}}^{n,\epsilon,k_n} + \eta_{\text{reg}}^{n,\epsilon,k_n} \right). \quad (4.1)$$

- (ii) *Regularization.* The regularization parameter ϵ is reduced until the value ϵ_n such that

$$\eta_{\text{reg}}^{n,\epsilon_n,k_n} \leq \Gamma_{\text{reg}} \left(\eta_{\text{sp}}^{n,\epsilon_n,k_n} + \eta_{\text{tm}}^{n,\epsilon_n,k_n} + \eta_{\text{qd}}^{n,\epsilon_n,k_n} \right). \quad (4.2)$$

- (iii) *Quadrature.* The quadrature rule is improved until

$$\eta_{\text{qd}}^{n,\epsilon_n,k_n} \leq \Gamma_{\text{qd}} \left(\eta_{\text{sp}}^{n,\epsilon_n,k_n} + \eta_{\text{tm}}^{n,\epsilon_n,k_n} \right). \quad (4.3)$$

Note that all the linearization, regularization, and quadrature errors may be classified as subsidiary as they can be made as small as desired by increasing the computational effort for fixed mesh and time step; it is thus reasonable to expect that the above criteria will be attained. Local, element by element, versions of the criteria (4.1)–(4.3) can be formulated using the local estimators (3.13) (see [26, 15, 19]), and require that the inequalities hold for all $K \in \mathcal{K}^n$; c.f. (7.4) for an example.

In the spirit of [41, 36, 48] and [18, 23, 14], we also propose the usual space–time adaptivity:

- (iv) *Space–time error balancing.* The space and time error components should be equilibrated by selecting the time step τ^n and adjusting the spatial meshes \mathcal{K}^n in such a way that

$$\gamma_{\text{tm}} \eta_{\text{sp}}^{n,\epsilon_n,k_n} \leq \eta_{\text{tm}}^{n,\epsilon_n,k_n} \leq \Gamma_{\text{tm}} \eta_{\text{sp}}^{n,\epsilon_n,k_n}. \quad (4.4)$$

Above, $\Gamma_{\text{tm}} > \gamma_{\text{tm}} > 0$ are again user-given parameters, typically close to 1.

- (v) *Adaptive mesh refinement.* The error in space should be evenly distributed throughout the domain Ω by local adaptation (refinement, coarsening) of the space mesh \mathcal{K}^n in such a way that, for all $K_1, K_2 \in \mathcal{K}^n$,

$$\eta_{\text{sp},K_1}^{n,\epsilon_n,k_n} \approx \eta_{\text{sp},K_2}^{n,\epsilon_n,k_n}.$$

In contrast to (4.1)–(4.3), the goal is to make $\eta_{\text{sp}}^{n,\epsilon_n,k_n}$ and $\eta_{\text{tm}}^{n,\epsilon_n,k_n}$ of comparable size as these error components are substantial and cannot be made arbitrarily small for a given choice of the mesh and of the time step.

4.2 Adaptive algorithm

In this section we propose an adaptive algorithm that implements the balancing and stopping criteria of Section 4.1. Moreover, for a prescribed $\zeta > 0$, we aim at satisfying the relation

$$\frac{\sum_{n=1}^N \|\mathcal{R}(u_{h\tau})\|_{X'_n}^2}{\sum_{n=1}^N \|\mathbf{1}_h^{n,\epsilon,k}\|_{L^2(I_n;L^2(\Omega))}^2} \leq \zeta^2, \quad (4.5)$$

i.e., to bring the relative error under the user-given precision ζ . To account for limited computing resources, we fix refinement thresholds $\underline{h}, \underline{\tau} > 0$ for both the mesh size and the time step and require, for all $0 \leq n \leq N$,

$$\min_{K \in \mathcal{K}^n} h_K \geq \underline{h}, \quad \tau^n \geq \underline{\tau}. \quad (4.6)$$

Note that, in particular because of (4.6), the attainment of (4.5) is not guaranteed.

Recall that $u_h^{n,\epsilon,k}$ stands for the approximation of the solution u_h^n at discrete time t^n obtained after k linearization iterations using a regularization parameter ϵ . At each linearization iteration k , the new approximation $u_h^{n,\epsilon,k}$ is obtained solving the linear problem written schematically as $u_h^{n,\epsilon,k} = \Psi(u_h^{n,\epsilon,k-1}, \tau^n, \mathcal{K}^n)$. For the sake of simplicity, in what follows we neglect the quadrature and data oscillation estimators. Our adaptive algorithm is the following:

Algorithm 4.1 (Adaptive algorithm).

Fix the fractions of cells to refine, c_{ref} , and to derefine, c_{deref}
Choose an initial mesh \mathcal{K}^0 , regularization parameter ϵ_0 , and a tolerance $\zeta_{\text{IC}} > 0$
 $u_h^0 \leftarrow \Pi^0(\beta_{\epsilon_0}^{-1}(\beta(u_0)))$
repeat {Initial mesh and regularization parameter adaptation}
 Compute η_{IC}
 Refine the cells $K \in \mathcal{K}^0$ such that $\eta_{\text{IC},K} \geq c_{\text{ref}} \max_{L \in \mathcal{K}^0} \{\eta_{\text{IC},L}\}$ in accordance with (4.6) and adjust the regularization parameter ϵ_0
 $u_h^0 \leftarrow \Pi^0(\beta_{\epsilon_0}^{-1}(\beta(u_0)))$
until $\eta_{\text{IC}} \leq \zeta_{\text{IC}} \|\nabla(\beta_{\epsilon_0}(u_h^0))\|_{L^2(\Omega)}$
Choose an initial time step τ^0
 $\epsilon \leftarrow \epsilon_0, t^0 \leftarrow 0, n \leftarrow 0$
while $t^n \leq T$ **do** {Time loop}
 $n \leftarrow n + 1$
 $\mathcal{K}^n \leftarrow \mathcal{K}^{n-1}$
 $\tau^n \leftarrow \tau^{n-1}$
 $u_h^{n,\epsilon,0} \leftarrow u_h^{n-1}$
 repeat {Space refinement}
 repeat {Space and time error balancing}
 repeat {Regularization}
 $k \leftarrow 0$
 repeat {Nonlinear solver}
 $k \leftarrow k + 1$
 $u_h^{n,\epsilon,k} = \Psi(u_h^{n,\epsilon,k-1}, \tau^n, \mathcal{K}^n)$
 Compute $\eta_{\text{sp}}^{n,\epsilon,k}, \eta_{\text{tm}}^{n,\epsilon,k}, \eta_{\text{reg}}^{n,\epsilon,k}, \eta_{\text{lin}}^{n,\epsilon,k}$
 until (4.1) is satisfied
 $k_n \leftarrow k$
 if (4.2) does not hold **then**
 $\epsilon \leftarrow \epsilon/2$
 end if
 until (4.2) is satisfied
 $\epsilon_n \leftarrow \epsilon$
 if $\eta_{\text{tm}}^{n,\epsilon_n,k_n} < \gamma_{\text{tm}} \eta_{\text{sp}}^{n,\epsilon_n,k_n}$ **then**
 $\tau^n \leftarrow 2\tau^n$
 else if $\eta_{\text{tm}}^{n,\epsilon_n,k_n} > \Gamma_{\text{tm}} \eta_{\text{sp}}^{n,\epsilon_n,k_n}$ and $\tau^n \geq 2\underline{\tau}$ **then**
 $\tau^n \leftarrow \tau^n/2$
 end if
 until (4.4) is satisfied or $\tau^n = \underline{\tau}$
 Refine the cells $K \in \mathcal{K}^n$ such that $\eta_{\text{sp},K}^{n,\epsilon_n,k_n} \geq c_{\text{ref}} \max_{L \in \mathcal{K}^n} \{\eta_{\text{sp},L}^{n,\epsilon_n,k_n}\}$ in accordance with (4.6)
until $\eta_{\text{sp}}^{n,\epsilon_n,k_n} + \eta_{\text{tm}}^{n,\epsilon_n,k_n} + \eta_{\text{reg}}^{n,\epsilon_n,k_n} + \eta_{\text{lin}}^{n,\epsilon_n,k_n} \leq \zeta \|\mathbf{1}_h^{n,\epsilon_n,k_n}\|_{L^2(I_n; L^2(\Omega))}$ or $(h_K = \underline{h}, \forall K \in \mathcal{K}^n)$

Derefine the cells $K \in \mathcal{K}^n$ such that $\eta_{\text{sp},K}^{n,\epsilon_n,k_n} \leq c_{\text{deref}} \max_{L \in \mathcal{K}^n} \{\eta_{\text{sp},L}^{n,\epsilon_n,k_n}\}$
 $u_h^n \leftarrow u_h^{n,\epsilon_n,k_n}$
 $t^n \leftarrow t^{n-1} + \tau^n$
 $\epsilon \leftarrow 2\epsilon$
end while

4.3 Efficiency of the a posteriori error estimate

In this section we investigate the global efficiency of the estimators of Corollary 3.5 under the stopping and balancing criteria of Section 4.1. Hence, the quantities at discrete time t^n are those obtained after performing k_n linearization iterations to meet the criterion (4.1), using a regularization parameter ϵ_n and a quadrature rule such that, respectively, (4.2) and (4.3) are satisfied, and a time step ensuring the time and space error balance (4.4). As usual, in order to use the argument of equivalence of norms on finite-dimensional spaces, we need to assume here:

Assumption 4.2 (Polynomial approximations). *For all $1 \leq n \leq N$, the function $u_{h\tau}^{n,\epsilon_n,k_n}$ is affine in time on the time interval I_n and piecewise polynomial of order m in space on the mesh $\mathcal{K}^{n-1,n}$; the functions $\mathbf{I}_h^{n,\epsilon_n,k_n}$ and $\mathbf{t}_h^{n,\epsilon_n,k_n}$ are piecewise polynomial of order m in space on $\mathcal{K}^{n-1,n}$.*

For $1 \leq n \leq N$, we introduce the standard residual-based a posteriori error estimators, c.f. [48]:

$$(\eta_{\text{res},1}^n)^2 := \tau^n \sum_{K \in \mathcal{K}^{n-1,n}} h_K^2 \|\hat{f}^n - \partial_t u_{h\tau}^{n,\epsilon_n,k_n} + \nabla \cdot \mathbf{I}_h^{n,\epsilon_n,k_n}\|_{L^2(K)}^2, \quad (4.7a)$$

$$(\eta_{\text{res},2}^n)^2 := \tau^n \sum_{F \in \mathcal{F}^{1,n-1,n}} h_F \|\mathbf{J}_h^{n,\epsilon_n,k_n} \mathbf{K} \cdot \mathbf{n}_F\|_{L^2(F)}^2. \quad (4.7b)$$

Let C be a generic constant only depending on the shape regularity parameter $\kappa_{\mathcal{K}}$ of the meshes $\mathcal{K}^{n-1,n}$, $1 \leq n \leq N$, the space dimension d , and the polynomial degree m . In order to still proceed generally, without the specification of a particular spatial discretization scheme, we will suppose the following:

Assumption 4.3 (Approximation property). *For all $1 \leq n \leq N$, there holds*

$$\tau^n \sum_{K \in \mathcal{K}^{n-1,n}} \|\mathbf{I}_h^{n,\epsilon_n,k_n} + \mathbf{t}_h^{n,\epsilon_n,k_n}\|_{L^2(K)}^2 \leq C \left((\eta_{\text{res},1}^n)^2 + (\eta_{\text{res},2}^n)^2 \right). \quad (4.8)$$

This property will be verified in Section 6 below for the vertex-centered finite volume spatial discretization and specific constructions of the fluxes $\mathbf{t}_h^{n,\epsilon_n,k_n}$ and $\mathbf{I}_h^{n,\epsilon_n,k_n}$.

Under these assumptions, we have the following result, showing the equivalence of the error $\|\mathcal{R}(u_{h\tau}^{n,\epsilon_n,k_n})\|_{X'_n}$ and the estimators of Corollary 3.5, up to data oscillation:

Theorem 4.4 (Global efficiency). *Let, for all $1 \leq n \leq N$, the stopping criteria (4.1)–(4.3) as well as the second inequality in the balancing criterion (4.4) be satisfied with the parameters Γ_{lin} , Γ_{reg} , Γ_{qd} , and Γ_{tm} small enough. Let Assumptions 4.2 and 4.3 hold true. Then*

$$\eta_{\text{sp}}^{n,\epsilon_n,k_n} + \eta_{\text{tm}}^{n,\epsilon_n,k_n} + \eta_{\text{qd}}^{n,\epsilon_n,k_n} + \eta_{\text{reg}}^{n,\epsilon_n,k_n} + \eta_{\text{lin}}^{n,\epsilon_n,k_n} \leq C \left(\|\mathcal{R}(u_{h\tau}^{n,\epsilon_n,k_n})\|_{X'_n} + \|f - \hat{f}\|_{X'_n} \right).$$

The proof of this result follows the techniques of [49] and the approach of [15]. It is given in Appendix A.1.

5 An a posteriori error estimate for the error in temperature and enthalpy

In the previous sections we have given a posteriori error estimators for the dual norm of the residual. In this section we prove that these same estimators also bound an error in temperature and enthalpy. We rely on a duality argument which reveals simpler than using the dual partial differential equation as in [36].

5.1 Bounding the error of the temperature and enthalpy by the dual norm of the residual

For brevity of notation we let for $t \in (0, T]$,

$$Q_t := L^2(0, t; L^2(\Omega)), \quad X_t := L^2(0, t; H_0^1(\Omega)), \quad X'_t := L^2(0, t; H^{-1}(\Omega)).$$

It is convenient to stress that the result of this section applies to all functions $u_{h\tau} \in Z$ such that $\beta(u_{h\tau}) \in X$. We first state the following bound:

Lemma 5.1 (Simple bounds for the temperature and enthalpy errors). *Let u be the solution of (2.2) and let $u_{h\tau} \in Z$ be such that $\beta(u_{h\tau}) \in X$. Then there holds*

$$\begin{aligned} \frac{L_\beta}{2} \|u - u_{h\tau}\|_{X'}^2 + \frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 \\ \leq \frac{L_\beta}{2} (2e^T - 1) \left(\|\mathcal{R}(u_{h\tau})\|_{X'}^2 + \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 \right), \end{aligned}$$

and

$$\frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 \leq \frac{L_\beta}{2} e^T \left(\|\mathcal{R}(u_{h\tau})\|_{X'}^2 + \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 \right).$$

The results of Lemma 5.1 are classical; we obtain them as a byproduct in the proof of Theorem 5.2 in Section A.2 below. These results are, however, not sufficiently precise. In particular, the use of the Gronwall lemma in its proof implies the appearance of the term e^T on the right-hand sides, which grows exponentially with the final time T . The purpose of the following theorem is to improve considerably this point. Indeed, note that, in Theorem 5.2, the term e^T does not appear elsewhere than in the approximation of the initial condition $\|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2$ which can be made sufficiently small. Theorem 5.2 takes a more complicated form than Lemma 5.1 but the numerical results based on its use, see Section 7, reveal excellent, which is not the case for the framework of Lemma 5.1:

Theorem 5.2 (An improved bound for the temperature and enthalpy errors). *Let u be the solution of (2.2) and let $u_{h\tau} \in Z$ be such that $\beta(u_{h\tau}) \in X$. Then there holds*

$$\begin{aligned} \frac{L_\beta}{2} \|u - u_{h\tau}\|_{X'}^2 + \frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 \\ + 2 \int_0^T \left(\|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \int_0^t \|\beta(u) - \beta(u_{h\tau})\|_{Q_s}^2 e^{t-s} ds \right) dt \\ \leq \frac{L_\beta}{2} \left\{ (2e^T - 1) \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'}^2 \right. \\ \left. + 2 \int_0^T \left(\|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \int_0^t \|\mathcal{R}(u_{h\tau})\|_{X'_s}^2 e^{t-s} ds \right) dt \right\}. \end{aligned}$$

The proof of this result is given in Section A.2.

5.2 The a posteriori error estimate

The upper bound in Theorem 5.2 can be combined with the results of Section 3.3 to obtain an a posteriori estimate for the temperature and enthalpy errors.

Theorem 5.3 (A posteriori estimate for the temperature and enthalpy errors). *Let u be the solution of (2.2) and let $u_{h\tau}$ and $\mathbf{t}_{h\tau}$ fulfill Assumptions 3.1 and 3.2, respectively. Then there*

holds

$$\begin{aligned}
& \frac{L_\beta}{2} \|u - u_{h\tau}\|_{X'}^2 + \frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 \\
& + 2 \int_0^T \left(\|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \int_0^t \|\beta(u) - \beta(u_{h\tau})\|_{Q_s}^2 e^{t-s} ds \right) dt \\
& \leq \frac{L_\beta}{2} \left\{ (2e^T - 1) \eta_{\text{IC}}^2 + \eta^2 \right. \\
& \quad \left. + 2 \left(\sum_{n=1}^N \tau^n \sum_{l=1}^n (\eta^l)^2 + \sum_{n=1}^N \sum_{l=1}^n J_{nl} \left\{ \sum_{i=1}^l (\eta^i)^2 \right\} \right) \right\},
\end{aligned} \tag{5.1}$$

with η_{IC} defined by (3.6), η defined by (3.8), η^n , $1 \leq n \leq N$, defined by

$$\eta^n := \left\{ \int_{I_n} \sum_{K \in \mathcal{K}^n} (\eta_{\text{R},K}^n + \eta_{\text{F},K}^n(t))^2 dt \right\}^{\frac{1}{2}} + \|f - \hat{f}\|_{X'_n}, \tag{5.2}$$

and setting, for $1 \leq n, l \leq N$,

$$J_{nl} := \int_{I_n} \int_{I_l} e^{t-s} ds dt.$$

Proof. To prove the result, we rely on Theorem 5.2. Applying Theorem 3.3, it follows that $\|\mathcal{R}(u_{h\tau})\|_{X'} \leq \eta$, so we are left to estimate the following right-hand side contributions in terms of the a posteriori error estimators:

$$\mathfrak{T}_1 := \int_0^T \|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 dt, \quad \mathfrak{T}_2 := \int_0^T \left(\int_0^t \|\mathcal{R}(u_{h\tau})\|_{X'_s}^2 e^{t-s} ds \right) dt.$$

As in Theorem 3.3, it is readily inferred that $\|\mathcal{R}(u_{h\tau})\|_{X'_l} \leq \eta^l$ for all $1 \leq l \leq N$, so that

$$\|\mathcal{R}(u_{h\tau})\|_{X'_{t_n}}^2 = \sum_{l=1}^n \|\mathcal{R}(u_{h\tau})\|_{X'_l}^2 \leq \sum_{l=1}^n (\eta^l)^2.$$

Using the fact that $\|\mathcal{R}(u_{h\tau})\|_{X'_l}$ is a nondecreasing function of the time t together with the above inequality yields for the first term

$$\mathfrak{T}_1 \leq \sum_{n=1}^N \int_{I_n} \|\mathcal{R}(u_{h\tau})\|_{X'_{t_n}}^2 dt \leq \sum_{n=1}^N \int_{I_n} \sum_{l=1}^n (\eta^l)^2 dt = \sum_{n=1}^N \tau^n \sum_{l=1}^n (\eta^l)^2.$$

Proceeding in a similar way, for the second term \mathfrak{T}_2 we obtain

$$\begin{aligned}
\mathfrak{T}_2 & \leq \sum_{n=1}^N \int_{I_n} \sum_{l=1}^n \int_{I_l} \|\mathcal{R}(u_{h\tau})\|_{X'_{t_l}}^2 e^{t-s} ds dt \\
& \leq \sum_{n=1}^N \int_{I_n} \sum_{l=1}^n \left\{ \int_{I_l} \sum_{i=1}^l (\eta^i)^2 e^{t-s} ds \right\} dt \\
& = \sum_{n=1}^N \sum_{l=1}^n \left\{ \int_{I_n} \int_{I_l} e^{t-s} ds dt \right\} \times \left\{ \sum_{i=1}^l (\eta^i)^2 \right\} = \sum_{n=1}^N \sum_{l=1}^n J_{nl} \left\{ \sum_{i=1}^l (\eta^i)^2 \right\},
\end{aligned}$$

whence the conclusion follows. \square

Remark 5.4 (Simplified versions of the a posteriori estimate). *In the spirit of Lemma 5.1, the following simplified versions of the a posteriori estimate of Theorem 5.3 hold:*

$$\begin{aligned}
\frac{L_\beta}{2} \|u - u_{h\tau}\|_{X'}^2 + \frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 & \leq \frac{L_\beta}{2} (2e^T - 1) (\eta^2 + \eta_{\text{IC}}^2), \\
\frac{L_\beta}{2} \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 + \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 & \leq \frac{L_\beta}{2} e^T (\eta^2 + \eta_{\text{IC}}^2).
\end{aligned}$$

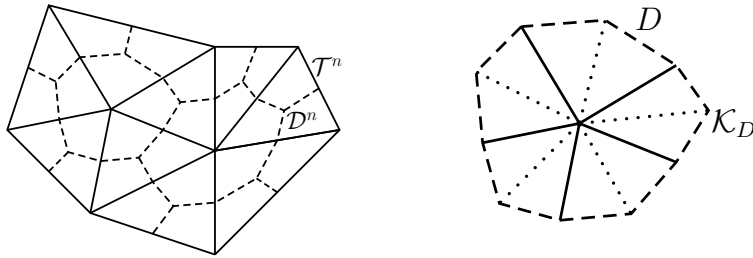


Figure 2: Simplicial mesh \mathcal{T}^n and the associated vertex-centered dual mesh \mathcal{D}^n (left) and the fine simplicial mesh \mathcal{K}_D of $D \in \mathcal{D}^n$ (right)

Remark 5.5 (An a posteriori error estimate distinguishing the different error components). *While relying on Corollary 3.5 instead of Theorem 3.3, equivalents of Theorem 5.3 and of the bounds of Remark 5.4 distinguishing the different error components can immediately be obtained.*

6 Application to a vertex-centered finite volume discretization

In this section, we consider the vertex-centered finite volume spatial and backward Euler temporal discretization of the Stefan problem (1.1). The regularization of Section 2.1.2 is considered and the Newton linearization is used. We show how to construct the equilibrated flux $\mathbf{t}_h^{n,\epsilon,k}$, the linearized flux $\mathbf{l}_h^{n,\epsilon,k}$, and the interpolation operator Π^n of Assumption 3.4 (in generalization of Assumptions 3.1 and 3.2) and verify Assumptions 4.2 and 4.3. Thus, all the results of Sections 3–5 will apply.

6.1 Dual and tertial space meshes

The vertex-centered finite volume method is defined using a sequence of dual meshes $\{\mathcal{D}^n\}_{0 \leq n \leq N}$ of the space domain Ω . For a given family of matching simplicial primal meshes $\{\mathcal{T}^n\}_{0 \leq n \leq N}$, we construct $\{\mathcal{D}^n\}_{0 \leq n \leq N}$ as follows: for any $1 \leq n \leq N$ and with every vertex \mathbf{a} of the mesh \mathcal{T}^n , we associate one dual volume D , constructed by connecting the barycenters of the simplices sharing \mathbf{a} through edge (and face for $d = 3$) barycenters, see Figure 2, left, for $d = 2$. We split every set \mathcal{D}^n into interior dual volumes $\mathcal{D}^{n,i}$ and boundary dual volumes $\mathcal{D}^{n,b}$. The simplicial mesh \mathcal{K}^n appearing in Sections 2–5 is constructed by dividing each $D \in \mathcal{D}^n$ into a mesh \mathcal{K}_D as indicated in Figure 2, right, if $d = 2$ and similarly for $d = 3$.

6.2 The vertex-centered finite volume scheme

Let, for $1 \leq n \leq N$,

$$V_h^n := \{\varphi_h \in C^0(\bar{\Omega}); \varphi_h|_K \in \mathbb{P}_1(K) \quad \forall K \in \mathcal{T}^n\}$$

and let

$$\Pi^n : C^0(\bar{\Omega}) \rightarrow V_h^n \quad \text{be the Lagrange interpolation operator,} \quad (6.1)$$

c.f. Ciarlet [11], which to a function $\varphi \in C^0(\bar{\Omega})$ associates a function $\varphi_h \in V_h^n$ by setting $\varphi_h(\mathbf{a}) := \varphi(\mathbf{a})$ for any vertex \mathbf{a} of the mesh \mathcal{T}^n .

Let $u_h^0 \in V_h^0$ be a suitable approximation of the regularized initial enthalpy $\beta_\epsilon^{-1}(\beta(u_0))$, see Algorithm 4.1. Let next $1 \leq n \leq N$, $u_h^{n-1} \in V_h^{n-1}$, and a mesh \mathcal{T}^n (and consequently \mathcal{D}^n) be given. The vertex-centered finite volume scheme for the regularized Stefan problem (2.3) reads: find $u_h^{n,\epsilon} \in V_h^n$ such that $\beta_\epsilon(u_h^{n,\epsilon})(\mathbf{a}) = 0$ for all vertices \mathbf{a} of \mathcal{T}^n on $\partial\Omega$ and such that

$$\frac{1}{\tau^n} (u_h^{n,\epsilon} - u_h^{n-1}, 1)_D - (\nabla \Pi^n(\beta_\epsilon(u_h^{n,\epsilon})) \cdot \mathbf{n}_D, 1)_{\partial D} = (\hat{f}^n, 1)_D \quad \forall D \in \mathcal{D}^{n,i}. \quad (6.2)$$

Then the continuous and piecewise affine-in-time function $u_{h\tau}$ appearing in the previous sections is given by $u_{h\tau}|_{I_n} := u_{h\tau}^{n,\epsilon}$,

$$u_{h\tau}^{n,\epsilon}(\cdot, t) = (1 - \rho(t))u_h^{n-1} + \rho(t)u_h^{n,\epsilon}, \quad \rho(t) := \frac{t - t^{n-1}}{\tau^n} \quad t \in I_n. \quad (6.3)$$

Remark 6.1 (Regularization). *It is also possible to consider the vertex-centered finite volume discretization without any regularization, i.e., use β in place of β_ϵ in (6.2), with $u_h^0 \in V_h^0$ an approximation of the initial enthalpy u_0 .*

Remark 6.2 (Links to the discretizations of [36, 4, 21]). *Let for simplicity the meshes \mathcal{T}^n (and consequently \mathcal{D}^n) do not move in time and let \hat{f}^n be piecewise constant on \mathcal{D}^n . Consider the case without regularization. Then the second and third terms of the scheme (6.2) coincide with that of [36, equation (4.4)], because of the links of the vertex-centered finite volumes and finite elements with mass lumping/quadrature for the source term. Similarly, in two space dimensions and when all the angles of \mathcal{T}^n are smaller than or equal to 90° , replacing the triangle barycenters by the triangle circumcenters in the construction of \mathcal{D}^n , the second and third terms of the scheme (6.2) coincide with that in the co-volume method of [4]. More generally, whenever \mathcal{T}^n is Delaunay and the mesh \mathcal{D}^n is its Voronoï dual, the same link holds true with the cell-centered finite volume scheme of [21], c.f., e.g., [36, 4, 21], [11, 20], or [51, Section 3]. Hence the only slight difference between (6.2) and these schemes is in the treatment of the time evolution term which is not mass-lumped herein.*

Remark 6.3 (Assumption 3.1). *By the definition of $u_{h\tau}$ by (6.2)–(6.3) and by the fact that $u_{h\tau}$ lies in a finite-dimensional space, $u_{h\tau} \in Z$ and $\beta(u_{h\tau}) \in X$, so that Assumption 3.1 is satisfied. A uniform bound could also be obtained by a priori stability analysis such as those in [36, 4, 21], but is not necessary in our setting.*

6.3 Newton linearization

Let $1 \leq n \leq N$ and the mesh \mathcal{T}^n (and \mathcal{D}^n) be fixed. Let the vector G^{n-1} be given by its components associated with the dual volumes $D \in \mathcal{D}^{n,i}$, $G_D^{n-1} := (u_h^{n-1}, 1)_D$, and similarly for the vector F^n , $F_D^n := (\hat{f}^n, 1)_D$. Let $u_h^{b,n,\epsilon} \in V_h^n$ take the values $\beta_\epsilon^{-1}(0)$ (0.5 for the example of Figure 1) at the boundary vertices of \mathcal{T}^n and the value zero at the other vertices of \mathcal{T}^n . The last vector that we need is $H^{n,\epsilon}$, $H_D^{n,\epsilon} := (u_h^{b,n,\epsilon}, 1)_D$. Let, for a given dual volume $E \in \mathcal{D}^{n,i}$, ϕ_E stand for the hat basis function of the space V_h^n associated with E : this is a function that takes the value 1 in the vertex associated with E and the value 0 at all other vertices of \mathcal{T}^n . We also define two matrices, with the components on the line associated with the dual volume $D \in \mathcal{D}^{n,i}$ and on the column associated with the dual volume $E \in \mathcal{D}^{n,i}$ given by $\mathbb{M}_{D,E}^n := (\phi_E, 1)_D$, $\mathbb{K}_{D,E}^n := (\nabla \phi_E \cdot \mathbf{n}_D, 1)_{\partial D}$. All the vectors are of size $\mathbb{R}^{|\mathcal{D}^{n,i}|}$ and the matrices of size $\mathbb{R}^{|\mathcal{D}^{n,i}| \times |\mathcal{D}^{n,i}|}$, with $|\mathcal{D}^{n,i}|$ the number of dual volumes in $\mathcal{D}^{n,i}$ (equal to the number of interior vertices of \mathcal{T}^n). The equation (6.2) can be written in matrix form as follows: find the vector $U^{n,\epsilon}$ such that

$$\mathbb{M}^n U^{n,\epsilon} - \tau^n \mathbb{K}^n \beta_\epsilon(U^{n,\epsilon}) = \tau^n F^n + G^{n-1} - H^{n,\epsilon}, \quad (6.4)$$

where $(\beta_\epsilon(U^{n,\epsilon}))_D := \beta_\epsilon(U_D^{n,\epsilon})$. We have $u_h^{n,\epsilon} = \sum_{E \in \mathcal{D}^{n,i}} U_E^{n,\epsilon} \phi_E + u_h^{b,n,\epsilon}$.

The algebraic system (6.4) is nonlinear. Its solution is approximated using the Newton linearization. Let $U^{n,\epsilon,0}$ be fixed; typically, $U^{n,\epsilon,0} := U^{n-1}$. Then, for $k \geq 1$, we approximate

$$\beta_\epsilon(U^{n,\epsilon,k}) \approx \beta_\epsilon(U^{n,\epsilon,k-1}) + \beta'_\epsilon(U^{n,\epsilon,k-1})(U^{n,\epsilon,k} - U^{n,\epsilon,k-1}). \quad (6.5)$$

The Newton linearization (6.5) is well defined since the regularized enthalpy–temperature function β_ϵ is continuously differentiable. At every Newton iteration k , we are thus lead to solve the following system of linear algebraic equations: find the vector $U^{n,\epsilon,k}$ such that

$$\begin{aligned} (\mathbb{M}^n - \tau^n \mathbb{K}^n \beta'_\epsilon(U^{n,\epsilon,k-1})) U^{n,\epsilon,k} &= \tau^n F^n + G^{n-1} - H^{n,\epsilon} \\ &\quad - \tau^n \mathbb{K}^n (\beta'_\epsilon(U^{n,\epsilon,k-1}) U^{n,\epsilon,k-1} - \beta_\epsilon(U^{n,\epsilon,k-1})). \end{aligned} \quad (6.6)$$

At each linearization step k , we set

$$u_h^{n,\epsilon,k} := \sum_{E \in \mathcal{D}^{n,i}} U_E^{n,\epsilon,k} \phi_E + u_h^{b,n,\epsilon}, \quad (6.7)$$

which is the function appearing in Section 3.4. The corresponding linearized flux of Assumption 3.4 is given by

$$\mathbf{l}_h^{n,\epsilon,k} := \nabla \left(\sum_{E \in \mathcal{D}^{n,i}} \left\{ \beta_\epsilon(U_E^{n,\epsilon,k-1}) + \beta'_\epsilon(U_E^{n,\epsilon,k-1})(U_E^{n,\epsilon,k} - U_E^{n,\epsilon,k-1}) \right\} \phi_E \right). \quad (6.8)$$

We perform the Newton iterations until we meet the convergence criterion discussed in Section 4.1.

6.4 Flux reconstruction

Let a time step $1 \leq n \leq N$, a regularization parameter $\epsilon > 0$, and a Newton linearization step k be fixed. We now show how to construct the flux $\mathbf{t}_h^{n,\epsilon,k}$ of Assumption 3.4. For this purpose, we will solve a local Neumann problem by mixed finite elements on every dual volume, following [30, 17, 51]. For a given $D \in \mathcal{D}^n$, we introduce the spaces

$$\begin{aligned} \mathbf{RTN}(\mathcal{K}_D) &:= \{\mathbf{v}_h \in \mathbf{H}(\operatorname{div}; D); \mathbf{v}_h|_K \in [\mathbb{P}_0(K)]^d + \mathbf{x} \mathbb{P}_0(K) \quad \forall K \in \mathcal{K}_D\}, \\ \mathbf{RTN}_N(\mathcal{K}_D) &:= \{\mathbf{v}_h \in \mathbf{RTN}(\mathcal{K}_D); \mathbf{v}_h \cdot \mathbf{n}_F = -\mathbf{l}_h^{n,\epsilon,k} \cdot \mathbf{n}_F \quad \forall F \in \partial \mathcal{K}_D^i\}, \\ \mathbf{RTN}_{N,0}(\mathcal{K}_D) &:= \{\mathbf{v}_h \in \mathbf{RTN}(\mathcal{K}_D); \mathbf{v}_h \cdot \mathbf{n}_F = 0 \quad \forall F \in \partial \mathcal{K}_D^i\}, \end{aligned}$$

where $\partial \mathcal{K}_D^i$ stands for all the faces of the submesh \mathcal{K}_D which are on the boundary of the dual volume D but not on the boundary of Ω . We will also need the space $\mathbb{P}_0^*(\mathcal{K}_D)$ which consists of piecewise constants functions on \mathcal{K}_D ; when $D \in \mathcal{D}^{n,i}$, we additionally impose a zero mean value over D . The local problem consists in finding $\mathbf{t}_h^{n,\epsilon,k} \in \mathbf{RTN}_N(\mathcal{K}_D)$ and $q_h \in \mathbb{P}_0^*(\mathcal{K}_D)$, the mixed finite element approximations of local Neumann problems on $D \in \mathcal{D}^{n,i}$ and local Neumann/Dirichlet problems on $D \in \mathcal{D}^{n,b}$:

$$(\mathbf{t}_h^{n,\epsilon,k} + \mathbf{l}_h^{n,\epsilon,k}, \mathbf{v}_h)_D - (q_h, \nabla \cdot \mathbf{v}_h)_D = 0 \quad \forall \mathbf{v}_h \in \mathbf{RTN}_{N,0}(\mathcal{K}_D), \quad (6.9a)$$

$$(\nabla \cdot \mathbf{t}_h^{n,\epsilon,k}, \phi_h)_D - (\hat{f}^n - \partial_t u_{h\tau}^{n,\epsilon,k}, \phi_h)_D = 0 \quad \forall \phi_h \in \mathbb{P}_0^*(\mathcal{K}_D). \quad (6.9b)$$

Note that the problem (6.9) is well-posed and one can take all $\phi_h \in \mathbb{P}_0(\mathcal{K}_D)$ as the test functions in (6.9b). Indeed, it follows from (6.6) and (6.8) that (compare to (6.2)) that

$$\frac{1}{\tau^n} (u_h^{n,\epsilon,k} - u_h^{n-1}, 1)_D - (\mathbf{l}_h^{n,\epsilon,k} \cdot \mathbf{n}_D, 1)_{\partial D} = (\hat{f}^n, 1)_D \quad \forall D \in \mathcal{D}^{n,i}. \quad (6.10)$$

From (6.10), we see that the Neumann boundary condition encoded in $\mathbf{RTN}_N(\mathcal{K}_D)$ is in equilibrium with the boundary datum $\hat{f}^n - \partial_t u_{h\tau}^{n,\epsilon,k}$ of (6.9). We have the following key result:

Lemma 6.4 (Assumptions 3.4, 4.2, and 4.3). *Let $1 \leq n \leq N$, $\epsilon > 0$, and $k \geq 1$ be fixed. Let $u_h^{n,\epsilon,k}$ be given by (6.6)–(6.7), $\mathbf{l}_h^{n,\epsilon,k}$ by (6.8), $\mathbf{t}_h^{n,\epsilon,k}$ by (6.9), and Π^n by (6.1). Then Assumptions 3.4, 4.2, and 4.3 hold true.*

Proof. The equilibrium property (3.11) follows immediately from (6.9b), so that Assumption 3.4 is easily satisfied. Whereas Assumption 4.2 is trivial, Assumption 4.3 is obtained by proceeding exactly as in [51, proof of Theorem 5.5] or [15, proof of Lemma 5.3]. \square

7 Numerical experiments

We illustrate in this section our theoretical results on a series of numerical experiments for the vertex-centered finite volume discretization approach of Section 6.

7.1 Setting

We consider the two-dimensional test case proposed by Nochetto et al. [34, 35] on the space–time domain $\Omega \times (0, T)$ with $\Omega = (0, 5)^2$ and $T = \pi/1.25$. The function $\beta(\cdot)$ is given by $\beta(u) = u^\ominus + (u - 1)^\oplus$. The exact temperature has the following expression:

$$(\beta(u))(x, y, t) = \begin{cases} 0.75(r^2 - 1), & \text{if } r < 1, \\ \left(1.5 - \rho'(t)\frac{y-\rho(t)}{r}\right)(r - 1), & \text{if } r \geq 1, \end{cases} \quad (7.1)$$

where $r^2 := x^2 + (y - \rho(t))^2$ and $\rho(t) := 0.5 + \sin(1.25t)$. The exact interface $I(t)$ is a circle with center $(0, \rho(t))$ and radius 1. The motion of the interface is governed by the Stefan law which prescribes that the normal velocity \mathbf{v} satisfies

$$(\nabla\beta(u)^+ - \nabla\beta(u)^-) \cdot \mathbf{n} = \mathbf{v} \quad \text{on } I(t),$$

where $\nabla\beta(u)^+$ and $\nabla\beta(u)^-$ denote the values of the temperature gradient on each side of the interface, while \mathbf{n} is the unit normal to the interface with suitable orientation. The enthalpy u on $\Omega \setminus I(t)$ can be obtained from the expression (7.1) of $\beta(u)$. The homogeneous Neumann condition $\nabla\beta(u) \cdot \mathbf{n} = 0$ is enforced at $x = 0$, whereas Dirichlet boundary conditions on the temperature are prescribed at $y = 0$, $y = 5$, and $x = 5$ using (7.1). The initial enthalpy u_0 and the source term f are likewise imposed using (7.1). The vertex-centered finite volume discretization of Section 6 is considered. No adaptation of the quadrature rule is performed; this is a reasonable simplification since the enthalpy–temperature function $\beta(\cdot)$ is piecewise affine.

7.2 Computing approximately the negative norms

In practice we cannot compute the negative norms as the initial data indicator η_{IC} , the data oscillation $\|f - \hat{f}\|_{X'}$, and the dual norm of the residual $\|\mathcal{R}(u_{h\tau})\|_{X'}$, even if the exact solution u is known. For numerical experiments below, the dual norms are approximated by solving auxiliary problems. More specifically, for a function $v \in X'$ to compute the negative norm $\|v\|_{X'}$ we consider for a.e. $t \in (0, T)$ the problem: find $\psi(\cdot, t) \in H_0^1(\Omega)$ such that

$$(\nabla\psi(\cdot, t), \nabla\varphi) = \langle v(\cdot, t), \varphi \rangle, \quad \forall \varphi \in H_0^1(\Omega). \quad (7.2)$$

Then

$$\begin{aligned} \|v\|_{X'}^2 &= \int_0^T \left\{ \sup_{\varphi \in H_0^1(\Omega), \|\nabla\varphi\|_{L^2(\Omega)}=1} \langle v(\cdot, t), \varphi \rangle \right\}^2 dt \\ &= \int_0^T \left\{ \sup_{\varphi \in H_0^1(\Omega), \|\nabla\varphi\|_{L^2(\Omega)}=1} (\nabla\psi(\cdot, t), \nabla\varphi) \right\}^2 dt \\ &= \int_0^T \|\nabla\psi\|_{L^2(\Omega)}^2(t) dt = \|\nabla\psi\|_{L^2(0,T;L^2(\Omega))}^2. \end{aligned}$$

We obtain an approximation of the function ψ by solving the problem (7.2) numerically by the vertex-centered finite volume scheme on a refined spatial mesh and on discrete times which refine the given temporal mesh. We suppose that the ensuing discretization error is small and can be ignored. The computation of η_{IC} is easier as it only involves the initial time $t = 0$.

7.3 Stopping criteria

We start by assessing the performance of the balancing and stopping criteria introduced in Section 4.1. Figure 3 depicts the evolution of the spatial (3.14a), temporal (3.14b), regularization (3.14d), and linearization (3.14e) error estimators as a function of the number of Newton iterations for a fixed mesh \mathcal{K} with $h_{\mathcal{K}} = 0.25$, time step $\tau = 0.1$, and regularization parameter

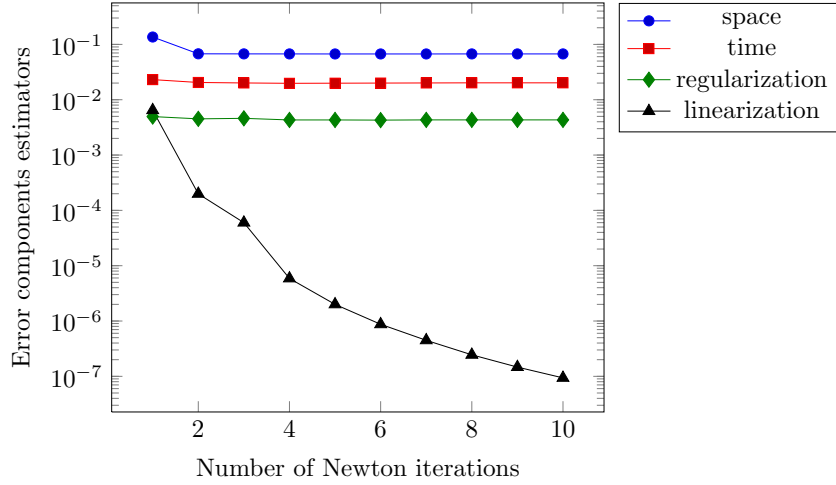


Figure 3: Evolution of the spatial, temporal, regularization, and linearization error estimators (3.14) as a function of Newton iterations for a fixed mesh, time step, and regularization parameter

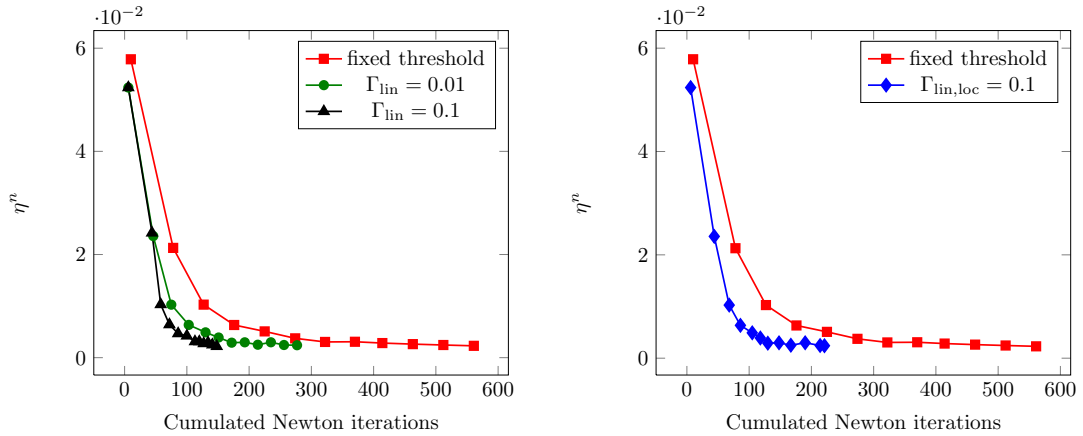


Figure 4: Error estimator η^n (5.2) as a function of the cumulated Newton iterations at each time step (time steps are identified by markers). Global stopping criterion (4.1) (left), local stopping criterion (7.4) (right)

$\epsilon = 0.05$. As expected, the linearization error steadily decreases, while the other components stagnate starting from the second iteration. The stopping criterion (4.1) with $\Gamma_{\text{lin}} = 10^{-2}$ allows to profit from this behavior by stopping the Newton algorithm after the second iteration, while a classical criterion based on a fixed threshold,

$$\eta_{\text{lin}}^{n,\epsilon,k} \leq \zeta_{\text{lin}}, \quad \zeta_{\text{lin}} = 10^{-7}, \quad (7.3)$$

would require 10 iterations to converge.

The overall gain for an entire simulation in terms of linearization iterations can be appreciated considering the results in Figure 4, left. We use the adaptive Algorithm 4.1 with different choices for the linearization stopping criterion: the classical criterion (7.3) then the stopping criterion (4.1) with $\Gamma_{\text{lin}} = 0.01$ and $\Gamma_{\text{lin}} = 0.1$. The others parameters used in the Algorithm are: $\Gamma_{\text{reg}} = 0.1$, $\zeta = 1$, $\zeta_{\text{IC}} = 1$, $h_{\mathcal{K}^0} = 0.1$, $\tau^0 = 0.1$, $\epsilon^0 = 0.25$, $\underline{h} = 10^{-2}$, $\underline{\tau} = 10^{-2}$, $c_{\text{ref}} = 0.7$, $c_{\text{deref}} = 0.2$, $\gamma_{\text{tm}} = 0.7$, and $\Gamma_{\text{tm}} = 1.3$. For the sake of completeness we also add a comparison with the local

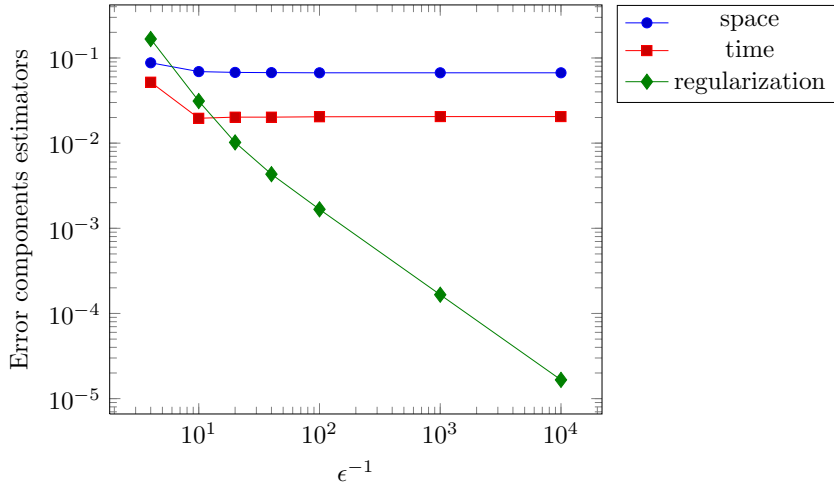


Figure 5: Evolution of the spatial, temporal, and regularization error estimators (3.14) as a function of ϵ^{-1} for a fixed mesh and time step

version of the stopping criterion (4.1), namely

$$\eta_{\text{lin},K}^{n,\epsilon,k_n} \leq \Gamma_{\text{lin,loc}} \left(\eta_{\text{sp},K}^{n,\epsilon,k_n} + \eta_{\text{tm},K}^{n,\epsilon,k_n} + \eta_{\text{qd},K}^{n,\epsilon,k_n} + \eta_{\text{reg},K}^{n,\epsilon,k_n} \right) \quad \forall K \in \mathcal{K}^n. \quad (7.4)$$

Even with this more stringent criterion, Figure 4, right, shows that a considerable gain in terms of number of linearization iterations can be achieved, whereas the precision on each time step (expressed by our error estimator η^n (5.2)) is basically unchanged.

Figure 5 shows similar results concerning the criterion (4.2) for the choice of the regularization parameter ϵ for a fixed mesh \mathcal{K} with $h_{\mathcal{K}} = 0.25$ and time step $\tau = 0.1$. For each value of ϵ , the Newton iterations are stopped according to (4.1) with $\Gamma_{\text{lin}} = 0.1$. The regularization error estimator decreases as expected when ϵ decreases, while the space and time error estimators stagnate starting from the third iteration. The criterion (4.2) with $\Gamma_{\text{reg}} = 0.1$ leads to stopping the iterations after the fourth step.

7.4 Balancing criteria

The next series of numerical experiments aims at assessing the space–time balancing criterion (4.4) by showing its impact on the estimated error (3.8) as a function of the total number of space–time unknowns $\sum_{n=1}^N |\mathcal{D}^{n,i}|$. In Figure 6, bottom left, we started by an initial mesh \mathcal{K}^0 with $h_{\mathcal{K}^0} = 0.4$ and a time step $\tau^0 = 0.1$. Then the time step is adapted in order to satisfy (4.4), with $\gamma_{\text{tm}} = 0.7$, $\Gamma_{\text{tm}} = 1.3$. As a result, the spatial (3.14a) and temporal (3.14b) error estimators stay equilibrated during the whole simulation. Figure 6, top, on the other hand, shows two possible disequilibrated patterns corresponding to space and time over-refinement. In the top left we started by an initial mesh \mathcal{K}^0 with $h_{\mathcal{K}^0} = 0.2$ and a time step $\tau^0 = 0.2$, we fixed also $\gamma_{\text{tm}} = 2$ and $\Gamma_{\text{tm}} = 3$, while in the top right we started by an initial mesh \mathcal{K}^0 with $h_{\mathcal{K}^0} = 0.5$ and time step $\tau^0 = 0.05$ and we fixed $\gamma_{\text{tm}} = \frac{1}{3}$ and $\Gamma_{\text{tm}} = \frac{1}{2}$. Finally, Figure 6, bottom right shows the effect of this violating of the balancing criterion (4.4) on the total error. These results make it apparent that the performance of an adaptive code may be considerably reduced when time and space errors are not balanced, and advocate the use of (4.4).

Next, we compare in Figure 7 the actual and predicted error distribution using the adaptive Algorithm 4.1 with $\Gamma_{\text{lin}} = \Gamma_{\text{reg}} = 0.1$, $\zeta = 1$, $\zeta_{\text{IC}} = 1$, $h_{\mathcal{K}^0} = 0.25$, $\tau^0 = 0.05$, $\epsilon^0 = 0.25$, $\underline{h} = 10^{-2}$, $\underline{\tau} = 10^{-2}$, $c_{\text{ref}} = 0.7$, $c_{\text{deref}} = 0.2$, $\gamma_{\text{tm}} = 0.7$, and $\Gamma_{\text{tm}} = 1.3$. We present the results at time $t = 0.1$. We see that the actual and predicted error distributions match very nicely. The corresponding exact and discrete enthalpies are depicted in Figure 8.

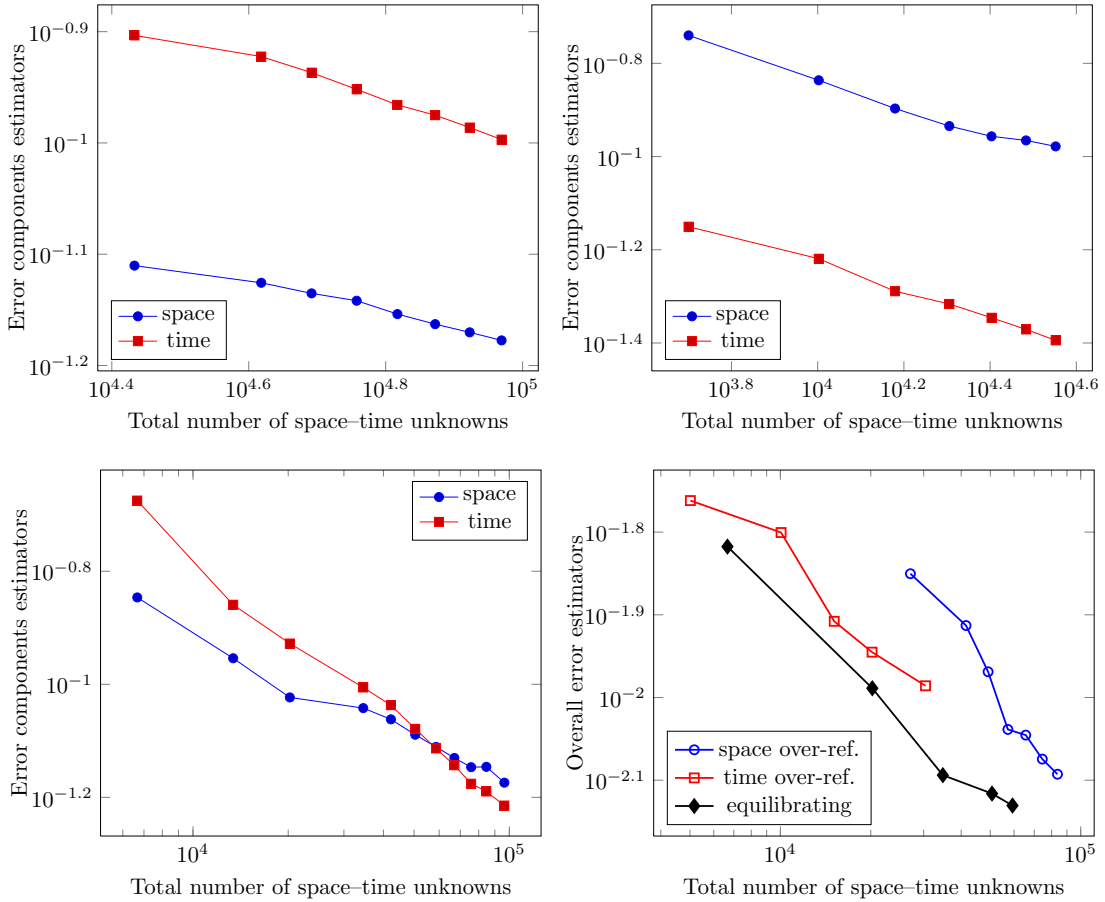


Figure 6: Effect of the time step adaptation strategy on the global error estimator (3.8). Violations of the balancing criterion (4.4) by space over-refinement (*top left*) and time over-refinement (*top right*). Time step refinement honoring (4.4) (*bottom left*). Overall comparison (*bottom right*)

7.5 Overall performance

In this section we assess the overall performance of the adaptive algorithm of Section 4.2 in terms of precision vs. the number of unknowns.

In Figure 9, left, we depict the error and estimates as a function of the total number of space-time unknowns in the fully adaptive case and in the uniform case. In the adaptive case, we use Algorithm 4.1 with the parameters detailed in Section 7.4. In the uniform case, the temporal and spatial meshes as well as the regularization parameter are fixed during the simulation, and linearization is stopped when (7.3) is satisfied. The error is measured in the dual norm (3.3) and estimated by Theorem 3.3 in the top part of Figure 9, whereas the energy-like norm (5.1) and the estimate of Theorem 5.3 are used in the bottom part of Figure 9. In both cases the adaptive strategy yields much better results than the uniform one, as expected. The right part of Figure 9 displays the corresponding effectivity indices, given by the ratio of the estimates over the error. These are remarkably close to the optimal value of one for the dual norm (3.3), even for the present time-dependent, degenerate problem with a moving free boundary. We regard the effectivity indices corresponding to Theorem 5.3 as likewise excellent; they are in particular several orders of magnitude smaller than the effectivity indices corresponding to the setting of Remark 5.4 that we have also assessed (not presented).

A quantitative evaluation of the performance in terms of precision vs. the number of unknowns

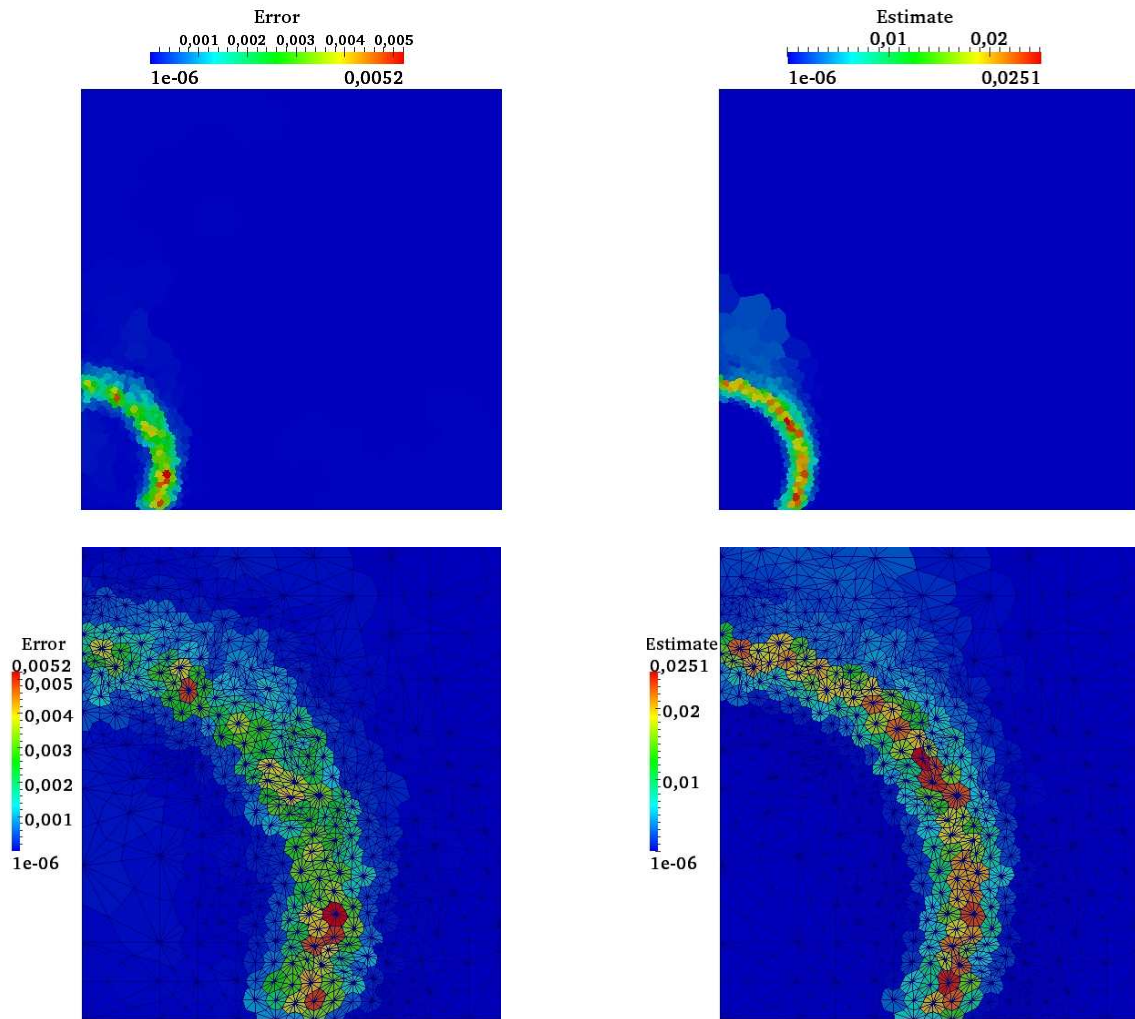


Figure 7: Actual (*left*) and estimated (*right*) error distribution for $\Gamma_{\text{lin}} = \Gamma_{\text{reg}} = 0.1$, adaptive Algorithm 4.1, entire domain (*top*), interface zoom (*bottom*)

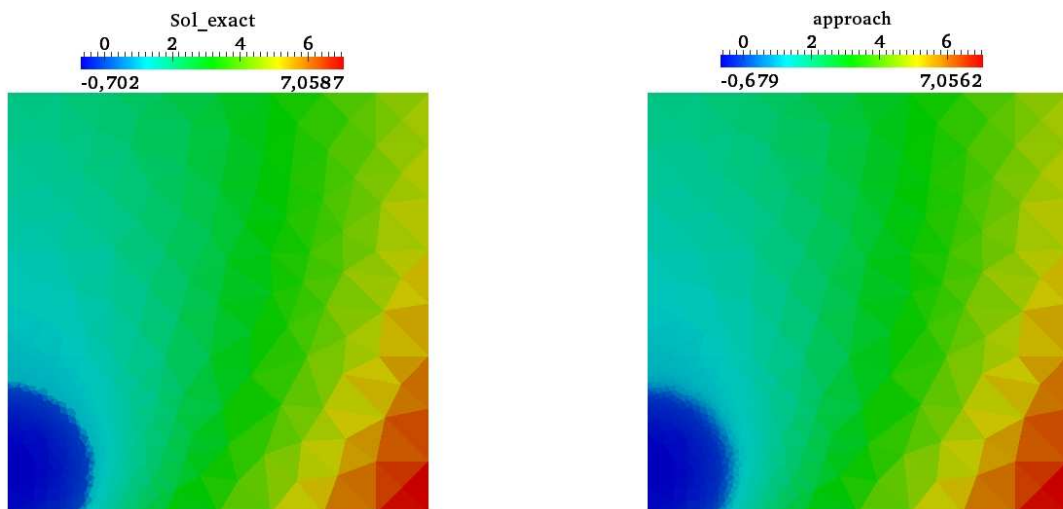


Figure 8: Exact (*left*) and approximate (*right*) enthalpy corresponding to the results of Figure 7

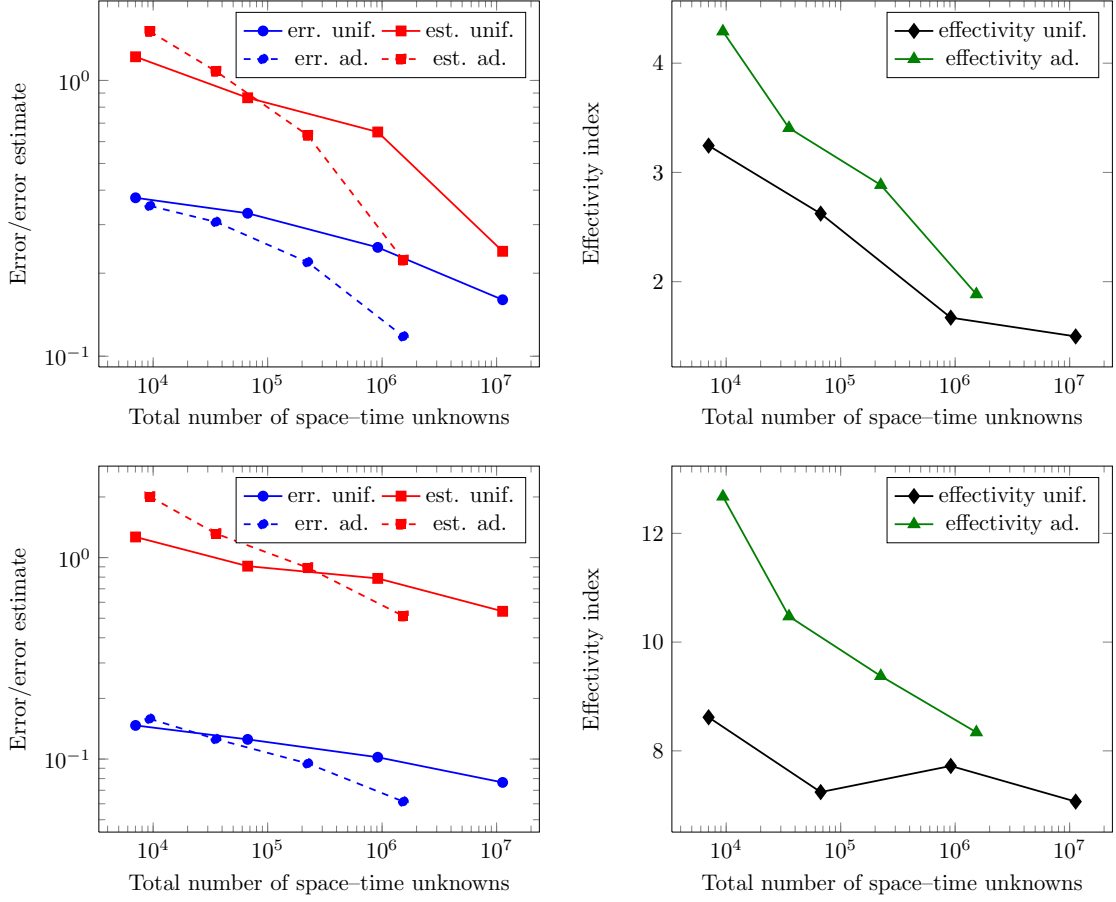


Figure 9: Comparison between adaptive and uniform refinement. Dual norm (3.3) (*top*), energy-like norm (5.1) (*bottom*). Error and estimators (*left*), effectivity indices (*right*)

Table 1: Comparison of the experimental orders of convergence (e.o.c.) in the uniform and fully adaptive cases. The total number of space-time unknowns is denoted by N_{st} . The actual error $\|\mathcal{R}(u_{h\tau})\|_{X'}$ and the estimated error η are defined by (3.2) and (3.8) respectively.

(a) Uniform case						
N_{st}	$\ \beta(u) - \beta(u_{h\tau})\ _{Q_T}$	e.o.c.	$\ \mathcal{R}(u_{h\tau})\ _{X'}$	e.o.c.	η	e.o.c.
7020	7.13e-02	–	3.75e-01	–	1.22e-00	–
66906	6.02e-02	0.224	3.30e-01	0.172	8.65e-01	0.455
915840	5.07e-02	0.197	2.48e-01	0.364	6.50e-01	0.392
1.12963e+07	2.19e-02	0.221	1.60e-01	0.115	2.40e-01	0.261
(b) Adaptive case						
N_{st}	$\ \beta(u) - \beta(u_{h\tau})\ _{Q_T}$	e.o.c.	$\ \mathcal{R}(u_{h\tau})\ _{X'}$	e.o.c.	η	e.o.c.
9360	6.55e-02	–	3.51e-01	–	1.51e-00	–
35370	5.28e-02	0.486	3.07e-01	0.303	1.08e-00	0.751
224082	4.06e-02	0.427	2.19e-01	0.546	6.32e-01	0.868
1.53329e+06	1.10e-02	0.392	1.18e-01	0.186	2.23e-01	0.312

can be obtained by computing the *experimental order of convergence* (e.o.c.), defined as follows:

$$\text{e.o.c} := \frac{\log(e_{N_{\text{st}}}) - \log(e_{M_{\text{st}}})}{-\frac{1}{3}(\log N_{\text{st}} - \log M_{\text{st}})},$$

where e denotes the chosen error measure while N_{st} and M_{st} are the total number of space–time unknowns corresponding to two subsequent levels of refinement. The results for the uniform and adaptive cases are collected in Tables 1a and 1b, respectively. We evaluate the dual norm of the residual (3.2), the $L^2(0, T; L^2(\Omega))$ error in the temperature, and the estimator η of (3.8). We observe roughly twice faster convergence in the adaptive case in comparison with the uniform one.

A Proofs

In this appendix, we collect the more involved proofs of some theorems of the paper.

A.1 Proof of Theorem 4.4

In this section, we will use the notation $a \lesssim b$ for the inequality $a \leq Cb$ with a generic constant C only depending on the shape regularity parameter $\kappa_{\mathcal{K}}$ of the meshes $\mathcal{K}^{n-1, n}$, $1 \leq n \leq N$, the space dimension d , and the polynomial degree m . Fix $1 \leq n \leq N$. We start by observing that, owing to the stopping criteria (4.1)–(4.3) and to the second inequality in the balancing criterion (4.4),

$$\eta_{\text{sp}}^{n, \epsilon_n, k_n} + \eta_{\text{tm}}^{n, \epsilon_n, k_n} + \eta_{\text{qd}}^{n, \epsilon_n, k_n} + \eta_{\text{reg}}^{n, \epsilon_n, k_n} + \eta_{\text{lin}}^{n, \epsilon_n, k_n} \lesssim \eta_{\text{sp}}^{n, \epsilon_n, k_n}. \quad (\text{A.1})$$

Recall that we have supposed in Section 2.2.2 that the mesh \mathcal{K}^n , $1 \leq n \leq N$, is obtained from \mathcal{K}^{n-1} by limited refinement/coarsening and that the common refinements $\mathcal{K}^{n-1, n}$ are uniformly shape regular. Thus, for $K \in \mathcal{K}^n$, using the triangle inequality, Assumption 4.2, and the inverse inequality, c.f. [44, Proposition 6.3.2], the first term of (3.13a) can be bounded by

$$\begin{aligned} \eta_{\text{R}, K}^{n, \epsilon_n, k_n} &= C_{\text{P}, K} h_K \|\hat{f}^n - \partial_t u_{h\tau}^{n, \epsilon_n, k_n} - \nabla \cdot \mathbf{t}_h^{n, \epsilon_n, k_n}\|_{L^2(K)} \\ &\leq C_{\text{P}, K} h_K \|\hat{f}^n - \partial_t u_{h\tau}^{n, \epsilon_n, k_n} + \nabla \cdot \mathbf{1}_h^{n, \epsilon_n, k_n}\|_{L^2(K)} + C_{\text{P}, K} h_K \|\nabla \cdot (\mathbf{1}_h^{n, \epsilon_n, k_n} + \mathbf{t}_h^{n, \epsilon_n, k_n})\|_{L^2(K)} \\ &\lesssim \left\{ \sum_{K' \in \mathcal{K}^{n-1, n}, K' \subset K} h_{K'}^2 \|\hat{f}^n - \partial_t u_{h\tau}^{n, \epsilon_n, k_n} + \nabla \cdot \mathbf{1}_h^{n, \epsilon_n, k_n}\|_{L^2(K')}^2 \right\}^{\frac{1}{2}} \\ &\quad + \left\{ \sum_{K' \in \mathcal{K}^{n-1, n}, K' \subset K} \|\mathbf{1}_h^{n, \epsilon_n, k_n} + \mathbf{t}_h^{n, \epsilon_n, k_n}\|_{L^2(K')}^2 \right\}^{\frac{1}{2}}. \end{aligned}$$

Consequently, employing Assumption 4.3,

$$\eta_{\text{sp}}^{n, \epsilon_n, k_n} \lesssim \eta_{\text{res}, 1}^n + \eta_{\text{res}, 2}^n. \quad (\text{A.2})$$

Proving the efficiency of the estimators introduced in Section 3.4 thus amounts to proving the efficiency of the residual estimators $\eta_{\text{res}, 1}^n$ and $\eta_{\text{res}, 2}^n$.

Henceforth, to simplify, we will use the shorthand notation

$$u_{h\tau}^n = u_{h\tau}^{n, \epsilon_n, k_n}, \quad \mathbf{1}_h^n = \mathbf{1}_h^{n, \epsilon_n, k_n}$$

and denote

$$(\eta_{\text{LRQT}}^n)^2 := \int_{I_n} \sum_{K \in \mathcal{K}^{n-1, n}} \|\nabla \beta(u_{h\tau}^n(\cdot, t)) - \mathbf{1}_h^n\|_{L^2(K)}^2 dt. \quad (\text{A.3})$$

We have:

Lemma A.1 (Estimate of $\eta_{\text{res},1}^n$). *Under the assumptions of Theorem 4.4, there holds*

$$\eta_{\text{res},1}^n \lesssim \|\mathcal{R}(u_{h\tau}^n)\|_{X'_n} + \eta_{\text{LRQT}}^n + \|f - \hat{f}\|_{X'_n}. \quad (\text{A.4})$$

Proof. For all $K \in \mathcal{K}^{n-1,n}$, we let $v_K := (\hat{f}^n - \partial_t u_{h\tau}^n + \nabla \cdot \mathbf{I}_h^n)|_K$. By Assumption 4.2, v_K is polynomial in K . We denote by ψ_K the usual bubble function on K , i.e., the product of the $(d+1)$ hat basis functions (barycentric coordinates) $\psi_{\mathbf{a}}$ associated with the vertices \mathbf{a} of the element K , set $\lambda_K := h_K^2 \psi_K v_K$ for all $K \in \mathcal{K}^{n-1,n}$, and let $\lambda := \sum_{K \in \mathcal{K}^{n-1,n}} \lambda_K$. Clearly, $\lambda \in H_0^1(\Omega)$ and $\lambda|_K \in H_0^1(K)$ for all $K \in \mathcal{K}^{n-1,n}$. Using the equivalence of norms on finite-dimensional spaces, integrating by parts in space, the weak form (2.2c), and (3.9) together with the Cauchy–Schwarz inequality, we infer, c.f. [49],

$$\begin{aligned} (\eta_{\text{res},1}^n)^2 &\lesssim \int_{I_n} \sum_{K \in \mathcal{K}^{n-1,n}} h_K^2 (v_K, \psi_K v_K)_K \, ds \\ &= \int_{I_n} \{ \langle \partial_t (u - u_{h\tau}^n), \lambda \rangle + (\nabla \beta(u) - \nabla \beta(u_{h\tau}^n), \nabla \lambda) \\ &\quad + (\nabla \beta(u_{h\tau}^n) - \mathbf{I}_h^n, \nabla \lambda) + (\hat{f}^n - f, \lambda) \} \, ds \\ &\leq \left(\|\mathcal{R}(u_{h\tau}^n)\|_{X'_n} + \eta_{\text{LRQT}}^n + \|f - \hat{f}\|_{X'_n} \right) \|\lambda\|_{X_n}. \end{aligned} \quad (\text{A.5})$$

By the shape regularity of the mesh $\mathcal{K}^{n-1,n}$ and the inverse inequality, c.f. [44, Proposition 6.3.2], we have, for any $K \in \mathcal{K}^{n-1,n}$,

$$\|\nabla \lambda\|_{L^2(K)} = h_K^2 \|\nabla(\psi_K v_K)\|_{L^2(K)} \lesssim h_K \|\psi_K v_K\|_{L^2(K)} \leq h_K \|v_K\|_{L^2(K)}.$$

An immediate consequence is that $\|\lambda\|_{X_n} \lesssim \eta_{\text{res},1}^n$ and (A.4) follows. \square

Lemma A.2 (Estimate of $\eta_{\text{res},2}^n$). *Under the assumptions of Theorem 4.4, there holds*

$$\eta_{\text{res},2}^n \lesssim \|\mathcal{R}(u_{h\tau}^n)\|_{X'_n} + \eta_{\text{LRQT}}^n + \|f - \hat{f}\|_{X'_n}. \quad (\text{A.6})$$

Proof. Let $F \in \mathcal{F}^{i,n-1,n}$. We denote by \mathcal{K}_F the simplices $K \in \mathcal{K}^{n-1,n}$ that share the face F . Let $v_F := \mathbf{J}_h^n \mathbf{K} \cdot \mathbf{n}_F$ and keep the same notation for the constant extension of v_F into \mathcal{K}_F along the vectors face barycenter–opposite vertex. Owing to Assumption 4.2, v_F is a polynomial on \mathcal{K}_F . Let ψ_F be the usual face bubble function supported on \mathcal{K}_F , i.e., the product of the d hat basis functions (barycentric coordinates) $\psi_{\mathbf{a}}$ associated with the vertices \mathbf{a} of the face F . For all $F \in \mathcal{F}^{i,n-1,n}$, set $\lambda_F := h_F \psi_F v_F$ and let $\lambda := \sum_{F \in \mathcal{F}^{i,n-1,n}} \lambda_F$. Note that $\lambda \in H_0^1(\Omega)$ and $\lambda|_{\mathcal{K}_F} \in H_0^1(\mathcal{K}_F)$ for all $F \in \mathcal{F}^{i,n-1,n}$. Using the equivalence of norms in finite-dimensional spaces, integrating by parts in space, using the weak form (2.2c), and (3.9) together with the Cauchy–Schwarz inequality, it is inferred, c.f. [49],

$$\begin{aligned} (\eta_{\text{res},2}^n)^2 &\lesssim \int_{I_n} \sum_{F \in \mathcal{F}^{i,n-1,n}} h_F (v_F, \psi_F v_F)_F \, ds \\ &= \int_{I_n} \sum_{F \in \mathcal{F}^{i,n-1,n}} \sum_{K \in \mathcal{K}_F} \{ (\nabla \cdot \mathbf{I}_h^n, \lambda_F)_K + (\mathbf{I}_h^n, \nabla \lambda_F)_K \} \, ds \\ &= \int_{I_n} \{ (\nabla \cdot \mathbf{I}_h^n, \lambda) + (\mathbf{I}_h^n, \nabla \lambda) \} \, ds \\ &= \int_{I_n} \{ \langle \partial_t (u_{h\tau}^n - u), \lambda \rangle + (\nabla \beta(u_{h\tau}^n) - \nabla \beta(u), \nabla \lambda) + (\hat{f}^n - \partial_t u_{h\tau}^n + \nabla \cdot \mathbf{I}_h^n, \lambda) \\ &\quad + (\mathbf{I}_h^n - \nabla \beta(u_{h\tau}^n), \nabla \lambda) + (f - \hat{f}^n, \lambda) \} \, ds \\ &\lesssim \left(\|\mathcal{R}(u_{h\tau}^n)\|_{X'_n} + \eta_{\text{LRQT}}^n + \|f - \hat{f}\|_{X'_n} \right) \|\lambda\|_{X_n} \\ &\quad + \eta_{\text{res},1}^n \left\{ \tau^n \sum_{K \in \mathcal{K}^{n-1,n}} h_K^{-2} \|\lambda\|_{L^2(K)}^2 \right\}^{\frac{1}{2}}. \end{aligned} \quad (\text{A.7})$$

Using the fact that, for all $F \in \mathcal{F}^{i,n-1,n}$ and $K \in \mathcal{K}_F^{n-1,n}$, $\|\psi_F v_F\|_{L^2(K)} \lesssim h_F^{\frac{1}{2}} \|v_F\|_{L^2(F)}$, it is inferred that $\tau^n \sum_{K \in \mathcal{K}^{n-1,n}} h_K^{-2} \|\lambda\|_{L^2(K)}^2 \lesssim (\eta_{\text{res},2}^n)^2$, whence by the inverse inequality, $\|\lambda\|_{X_n} \lesssim \eta_{\text{res},2}^n$. Using this fact in (A.7) in conjunction with (A.4), (A.6) follows. \square

Proof of Theorem 4.4. It follows from Lemmas A.1 and A.2 and from (A.2) that

$$\eta_{\text{sp}}^{n,\epsilon_n,k_n} \lesssim \|\mathcal{R}(u_{h\tau}^{n,\epsilon_n,k_n})\|_{X'_n} + \eta_{\text{LRQT}}^n + \|f - \hat{f}\|_{X'_n}. \quad (\text{A.8})$$

In order to bound the term η_{LRQT}^n , we proceed as follows. The triangle inequality and the definitions (3.13) and (3.14) give

$$\eta_{\text{LRQT}}^n \leq \eta_{\text{lin}}^{n,\epsilon_n,k_n} + \eta_{\text{reg}}^{n,\epsilon_n,k_n} + \eta_{\text{qd}}^{n,\epsilon_n,k_n} + \eta_{\text{tm}}^{n,\epsilon_n,k_n}.$$

Thus, proceeding as for the bound (A.1),

$$\eta_{\text{LRQT}}^n \leq C \eta_{\text{sp}}^{n,\epsilon_n,k_n},$$

where the constant C only depends on the parameters Γ_{lin} , Γ_{reg} , and Γ_{qd} in the stopping criteria (4.1)–(4.3) and Γ_{tm} in the balancing criterion (4.4). Thus, choosing these parameters small enough, the term η_{LRQT}^n can be made small enough to be discarded from the right-hand side of (A.8), c.f. [15, Theorem 4.4] and the assertion of Theorem 4.4 follows from (A.1). \square

A.2 Proof of Theorem 5.2

We start by proving the following intermediate result.

Lemma A.3 (Duality bound). *Let u be the solution to (2.2) and let $u_{h\tau} \in Z$ be such that $\beta(u_{h\tau}) \in X$. Then, there holds, for a.e. $t \in (0, T)$,*

$$\frac{2}{L_\beta} \|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \|(u - u_{h\tau})(\cdot, t)\|_{H^{-1}(\Omega)}^2 \leq \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \|u - u_{h\tau}\|_{X'_t}^2. \quad (\text{A.9})$$

Proof. For a.e. $t \in (0, T)$, we denote by $W(\cdot, t) \in H_0^1(\Omega)$ the solution to

$$(\nabla W(\cdot, t), \nabla \psi) = ((u - u_{h\tau})(\cdot, t), \psi) \quad \forall \psi \in H_0^1(\Omega). \quad (\text{A.10})$$

The existence and uniqueness of $W(\cdot, t)$ follow from the Lax–Milgram lemma. Moreover, since $u, u_{h\tau} \in Z$, there holds $W \in X$. Using (A.10), it is inferred that

$$\begin{aligned} \|\nabla W(\cdot, t)\|_{L^2(\Omega)} &= \sup_{\psi \in H_0^1(\Omega), \|\nabla \psi\|_{L^2(\Omega)}=1} (\nabla W(\cdot, t), \nabla \psi) \\ &= \sup_{\psi \in H_0^1(\Omega), \|\nabla \psi\|_{L^2(\Omega)}=1} ((u - u_{h\tau})(\cdot, t), \psi) = \|(u - u_{h\tau})(\cdot, t)\|_{H^{-1}(\Omega)}. \end{aligned} \quad (\text{A.11})$$

This duality technique is rather standard; see [9] and the references therein. Its origins can be traced back at least to the elliptic projection of Wheeler [53]. In some aspects, it is close to the elliptic reconstruction of Makridakis and Nochetto [31]; however, in [31] it is used to restore optimal order of the a posteriori estimate in $L^\infty(0, T; L^2(\Omega))$, whereas here we employ it to obtain a bound on an energy-like norm.

Taking $\varphi = W 1_{(0,t)}$ with $1_{(0,t)}$ the characteristic function of the interval $(0, t)$ in definition (3.1) and using (A.11) and the Young inequality, it is inferred

$$\langle \mathcal{R}(u_{h\tau}), W \rangle_{X'_t, X_t} \leq \|\mathcal{R}(u_{h\tau})\|_{X'_t} \|u - u_{h\tau}\|_{X_t} \leq \frac{1}{2} \|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \frac{1}{2} \|u - u_{h\tau}\|_{X_t}^2. \quad (\text{A.12})$$

Moreover,

$$\langle \mathcal{R}(u_{h\tau}), W \rangle_{X'_t, X_t} = \int_0^t \langle \partial_t(u - u_{h\tau}), W \rangle(s) ds + \int_0^t (\nabla\beta(u) - \nabla\beta(u_{h\tau}), \nabla W)(s) ds =: \mathfrak{R}_1 + \mathfrak{R}_2. \quad (\text{A.13})$$

Recalling (A.10), and since $u - u_{h\tau} \in H^1(0, T; H^{-1}(\Omega))$, there holds $\partial_t W \in X$ and, for a.e. $s \in (0, T)$, $\partial_t W(\cdot, s)$ satisfies in a weak sense

$$\begin{aligned} -\nabla \cdot (\nabla \partial_t W(\cdot, s)) &= \partial_t(u - u_{h\tau})(\cdot, s) && \text{in } \Omega, \\ \partial_t W(\cdot, s) &= 0 && \text{on } \partial\Omega. \end{aligned}$$

Thus, it follows from the definition (A.10) of W and from the norm characterization (A.11) that

$$\begin{aligned} \mathfrak{R}_1 &= \int_0^t (\partial_t \nabla W, \nabla W)(s) ds = \frac{1}{2} \left(\|\nabla W(\cdot, t)\|_{L^2(\Omega)}^2 - \|\nabla W(\cdot, 0)\|_{L^2(\Omega)}^2 \right) \\ &= \frac{1}{2} \left(\|(u - u_{h\tau})(\cdot, t)\|_{H^{-1}(\Omega)}^2 - \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 \right). \end{aligned} \quad (\text{A.14})$$

Invoking again the definition (A.10) and using the fact that β is nondecreasing and L_β -Lipschitz continuous, there holds

$$\begin{aligned} \mathfrak{R}_2 &= \int_0^t (u - u_{h\tau}, \beta(u) - \beta(u_{h\tau}))(s) ds \geq \frac{1}{L_\beta} \int_0^t (\beta(u) - \beta(u_{h\tau}), \beta(u) - \beta(u_{h\tau}))(s) ds \\ &= \frac{1}{L_\beta} \|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2. \end{aligned} \quad (\text{A.15})$$

The conclusion follows using inequalities (A.12), (A.14), and (A.15) in equation (A.13). \square

Corollary A.4 (Application of the Gronwall lemma). *Under the assumptions of Lemma A.3, there holds*

$$\begin{aligned} \|u - u_{h\tau}\|_{X'}^2 &\leq (e^T - 1) \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \int_0^T \left(\|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \int_0^t \|\mathcal{R}(u_{h\tau})\|_{X'_s}^2 e^{t-s} ds \right) dt \\ &\quad - \frac{2}{L_\beta} \int_0^T \left(\|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \int_0^t \|\beta(u) - \beta(u_{h\tau})\|_{Q_s}^2 e^{t-s} ds \right) dt. \end{aligned}$$

Proof. Using (A.9) followed by the Gronwall lemma

$$\xi(t) \leq \alpha(t) + \int_0^t \xi(s) ds \implies \xi(t) \leq \alpha(t) + \int_0^t \alpha(s) e^{t-s} ds,$$

with $\xi(t) := \|(u - u_{h\tau})(\cdot, t)\|_{H^{-1}(\Omega)}^2$ and $\alpha(t) := \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 - \frac{2}{L_\beta} \|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2$, it is inferred, for a.e. $t \in (0, T)$,

$$\begin{aligned} \|(u - u_{h\tau})(\cdot, t)\|_{H^{-1}(\Omega)}^2 &\leq e^t \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \int_0^t \|\mathcal{R}(u_{h\tau})\|_{X'_s}^2 e^{t-s} ds \\ &\quad - \frac{2}{L_\beta} \left(\|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \int_0^t \|\beta(u) - \beta(u_{h\tau})\|_{Q_s}^2 e^{t-s} ds \right). \end{aligned}$$

The assertion follows by integrating over the interval $(0, T)$. \square

We are now ready to prove Theorem 5.2:

Proof of Theorem 5.2. Using (A.9) with $t = T$ and adding $\|u - u_{h\tau}\|_{X'}^2$ to both sides we infer

$$\begin{aligned} \mathfrak{L} &:= \frac{2}{L_\beta} \|\beta(u) - \beta(u_{h\tau})\|_{Q_T}^2 + \|u - u_{h\tau}\|_{X'}^2 + \|(u - u_{h\tau})(\cdot, T)\|_{H^{-1}(\Omega)}^2 \\ &\leq \|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'}^2 + 2\|u - u_{h\tau}\|_{X'}^2. \end{aligned}$$

Using Corollary A.4 to estimate the last term in the right-hand side we obtain

$$\begin{aligned} \mathfrak{L} &\leq (2e^T - 1)\|u_0 - u_{h\tau}(\cdot, 0)\|_{H^{-1}(\Omega)}^2 + \|\mathcal{R}(u_{h\tau})\|_{X'}^2 \\ &\quad + 2 \int_0^T \left(\|\mathcal{R}(u_{h\tau})\|_{X'_t}^2 + \int_0^t \|\mathcal{R}(u_{h\tau})\|_{X'_s}^2 e^{t-s} ds \right) dt \\ &\quad - \frac{4}{L_\beta} \int_0^T \left(\|\beta(u) - \beta(u_{h\tau})\|_{Q_t}^2 + \int_0^t \|\beta(u) - \beta(u_{h\tau})\|_{Q_s}^2 e^{t-s} ds \right) dt. \end{aligned}$$

The conclusion follows multiplying both sides by $L_\beta/2$ and rearranging the terms. \square

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