

# Explicit geometric construction of *sparse inverse mass matrices* for arbitrary tetrahedral grids

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## Abstract

The geometric reinterpretation of the Finite Element Method (FEM) shows that Raviart–Thomas and Nédélec mass matrices map from degrees of freedoms (DoFs) attached to geometric elements of a tetrahedral grid to DoFs attached to the barycentric dual grid. The algebraic inverses of the mass matrices map DoFs attached to the barycentric dual grid back to DoFs attached to the corresponding primal tetrahedral grid, but they are of limited practical use since they are dense.

In this paper we present a new geometric construction of *sparse inverse mass matrices* for arbitrary tetrahedral grids and possibly inhomogeneous and anisotropic materials, debunking the conventional wisdom that the barycentric dual grid prohibits a sparse representation for inverse mass matrices. In particular, we provide a unified framework for the construction of both edge and face mass matrices and their sparse inverses. Such a unifying principle relies on novel geometric reconstruction formulas, from which, according to a well-established design strategy, local mass matrices are constructed as the sum of a consistent and a stabilization part. A major difference with the approaches proposed so far is that the consistent part is defined geometrically and explicitly, that is, without the necessity of computing the inverses of local matrices. This provides a sensible speedup and an easier implementation. We use these new sparse inverse mass matrices to discretize a three-dimensional Poisson problem, providing the comparison between the results obtained by various formulations on a benchmark problem with analytical solution.

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## 1. Introduction

Many physics-compatible discretization methods depend on a primal–dual mesh structure and the related staggered positioning of the physical variables [1], often referred to as *degrees of freedom* (DoFs). In particular, most physics-compatible discretizations for the Maxwell’s equations introduced so far, like the Yee-scheme [2], the Cell Method (CM) [1,3], the Finite Integration Technique [4], the Discrete Geometric Approach (DGA) [5,6], the Compatible Discrete Operators (CDO) [7] are based on a pair of interlocked primal and dual grids. In other physics-compatible methods like the Mimetic Finite Difference method [8] and the Finite Element Method (FEM) based on

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Whitney edge and face basis functions (i.e. the Nédélec curl-conforming basis functions and the Raviart–Thomas div-conforming basis functions) [9,10], a dual grid is not explicitly used but an interpretation using it is readily available [11,10], at least for the lowest order version.

A fundamental principle of compatible numerical schemes is the distinction between *topological equations* and *constitutive equations* [1,12]. Topological equations express conservation laws of physical theories. As the name suggests, topological equations are valid under arbitrary homeomorphic transformations of the domain in which they are defined. Constitutive equations, also called *material equations*, describe the behavior of material, substance or medium under interest. The *discrete Hodge star operator* [13,10] acts as discrete counterpart of the material parameters used to formulate constitutive equations and its form depends on the analytic form of the material parameters and on the geometric properties of the grid which discretizes the computational domain. In particular, the algebraic realization of the Hodge operator is given by the so-called *mass matrices* [11,13,10].

What is usually achieved is the construction of discrete Hodge operators which map DoFs attached to primal faces to DoFs attached to dual edges or DoFs attached to primal edges to DoFs attached to dual faces [13,10,5,6]. This is exactly what the FEM mass matrices computed with Raviart–Thomas and Nédélec basis functions perform, [11,10], respectively.

However, we can nonetheless define *inverse discrete Hodge operators* that map DoFs attached to dual geometric elements to DoFs attached to primal geometric elements. Explicitly computing the algebraic inverse of the mass matrix is not considered a viable solution given that such a matrix would be dense, so of questionable usefulness in practice. Therefore, we are in particular interested in a *sparse realization* of inverse discrete Hodge operators. Inverse Hodge operators play a vital role in many applications. Most notably, inverse Hodge operators enable consistent and explicit schemes to solve time-domain wave propagation problems [2,14–16]. Other applications enabled by the inverse mass matrices comprise the explicit construction of the *codifferential operator*, the *Laplace–de Rham operator* [17] and compute the discrete Hodge decomposition of discrete fields [17,18].

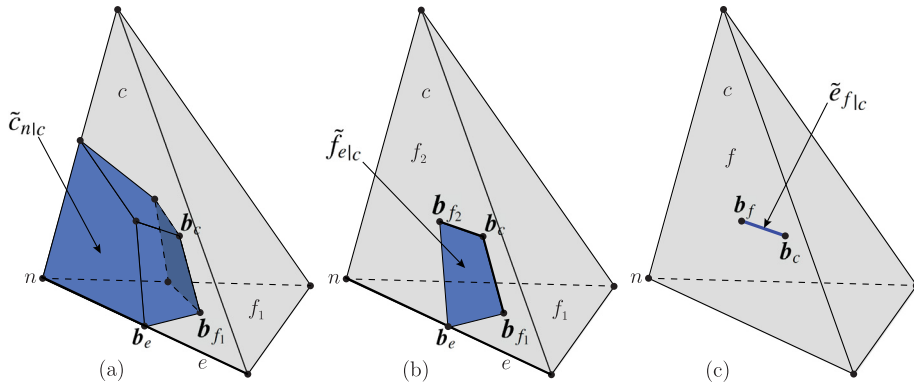
The construction of inverse Hodge operators for tetrahedral grids is not entirely new. In CM [1,3], Discrete Exterior Calculus (DEC) [19] and in the cell-centered Finite Volume literature [20], a *Voronoi dual grid* based on circumcenters is used, and the resulting mass matrices are diagonal in such a way that their inverses can be easily computed. It is important to note that, when the material is anisotropic, the mass matrices are in general not diagonal, so a sparse inverse Hodge operator cannot be easily constructed. Another solution using a Voronoi dual grid is presented in [21]. However, most commonly used mesh generators like NETGEN and GMSH produce tetrahedral grids that are not Delaunay and in this case the Voronoi dual grid cannot be defined [22].

The aim of this paper is to extend the construction of sparse inverse Hodge operators on *barycentric dual grids* which can be defined for arbitrary tetrahedral grids. The barycentric dual grids are explicitly used in DGA [6] and implicitly in the Finite Element Method (FEM) [11] and in the MFD [8]. Yet, devising a recipe to construct sparse inverse Hodge operators on a barycentric dual grid appears to be a formidable task [23,15,17,21] given that the conventional wisdom is that the barycentric dual grid “prohibits a sparse representation for their inverse operators” [24] and, consequently, only approximate constructions have been proposed [15].

By using a barycentric dual grid, in [14,25,16], inverse mass matrices that map from dual faces to primal edges are constructed by assembling local contributions inside dual cells and then computing the algebraic inverse of the resulting local matrices. Yet, we note that the time needed to compute all local inverses is not negligible, because the rank of the local matrices is twenty or more. Simultaneously with our paper, in [26] similar ideas are developed by using the differential forms formalism. Although the differential forms approach is more general, it is much more complicated to present. In addition, the theory is developed for homogeneous materials only.

In this paper we provide a unified framework for the construction of both edge and face mass matrices and their sparse inverses. The unifying principle relies on the construction of discrete Hodge operators starting from *reconstruction formulas* (that provide a representation of the discrete variables at the continuous level) defined by geometric elements of primal and dual grids. Starting from these reconstruction formulas, local mass matrices are constructed according to a well-established design strategy, as the sum of a consistent and a stabilization part [8,27]. While the consistent part is fixed, the stabilization part depends on some user-dependent parameters, which for example, offer the possibility of controlling the eigenvalues of the matrices.

The construction of the inverse mass matrix that maps from dual edges of a barycentric dual grid to primal faces of an arbitrary tetrahedral grid and possibly inhomogeneous materials proposed in this paper is entirely new. A key difference with respect to the approach proposed in [14,25,16] in the construction of the other inverse mass matrix



**Fig. 1.** Primal and dual geometric elements of a tetrahedron  $c \in C$ . (a) Dual cell. (b) Dual face. (c) Dual edge.

that maps from dual faces of the barycentric dual grid to primal edges of the simplicial grid is the *explicit* and geometric construction of the consistent part, without the necessity of computing local inverses. In addition, the recipe to construct the stabilization part adds more flexibility in the construction of the mass matrices.

In general, the major contribution of our construction is not only to have avoided the computation of the local inverses, but rather the fact that it results in symmetric expressions for both types of local mass matrices, thus highlighting the duality relationship between the pair of grids. A key requirement of our setting would be that the material parameters are constant on dual cells associated with the nodes of the primal simplicial grid. Instead, to deal with the general case of material parameters that are constant on the grid cells, but arbitrary discontinuous across the interfaces between the cells, two different approaches are proposed. The first one is a weighted averaging technique classically used in Finite Volumes Methods [20]. The second one is based on an extension of the approach proposed [14,16].

The paper is organized as follows. In Section 2, we introduce the geometric elements from which general tetrahedral grids and their barycentric duals are constructed. In addition, we present the basic building blocks of low-order compatible numerical schemes that will be used in the work. In Section 3, we detail the construction of mass matrices. In Section 4, following the same design principles of the previous section, we detail the construction of inverses of mass matrices. In addition, we describe how to deal with problems having inhomogeneous materials. In Section 5, we present an application of the newly derived inverse mass matrices to a Poisson problem and we provide a comparison among different formulations that can be used to solve the same problem on the same mesh. Finally, in Section 6, the conclusions are drawn.

## 2. Compatible discretization

A detailed presentation of compatible discretizations can be found in [28]. In what follows, we only present the main ideas that will be used throughout the paper.

### 2.1. Geometry of primal and dual grid

Let us consider a subdivision of a region  $\Omega$  of  $\mathbb{R}^3$  into a primal tetrahedral grid  $K = (N, E, F, C)$ , where the sets  $N, E, F$ , and  $C$  contain the grid *nodes*, *edges*, *faces*, and *cells* (or *tetrahedra*), respectively. We denote a *node* by  $n$ , an *edge* by  $e$ , a *face* by  $f$ , and a *cell* by  $c$ , see Fig. 1. The geometric elements of the primal grid are provided with an inner orientation [1,12].

Let  $X$  be any set among  $N, E, F$ , or  $C$ . If  $r$  is a geometric element of  $K$ , we denote by  $X(r)$  the subset defined by

$$X(r) := \{x \in X \mid r \subset x\}, \quad (1)$$

if (1) is not void, otherwise,

$$X(r) := \{x \in X \mid x \subset r\}. \quad (2)$$

For instance,  $C(e) = \{c \in C \mid e \subset c\}$  collects the cells of which  $e$  is an edge and  $E(c) = \{e \in E \mid e \subset c\}$  collects the edges of  $c$ .

Interlocked with the primal grid  $K$ , a *barycentric dual grid* [29,1,12]  $\tilde{K} = (\tilde{N}, \tilde{E}, \tilde{F}, \tilde{C})$  is introduced, where the sets  $\tilde{N}$ ,  $\tilde{E}$ ,  $\tilde{F}$ , and  $\tilde{C}$  contain *dual nodes*, *dual edges*, *dual faces*, and *dual cells*, respectively. Each geometric entity of the dual grid is in one-to-one correspondence (duality pairing) with a geometric element of the primal grid and it is constructed by means of the *barycentric subdivision* [29] of the primal grid, see Fig. 1. With symbol “ $\sim$ ” we denote geometric elements of the dual grid, thus, we denote by  $\tilde{n}$  a dual node, by  $\tilde{e}$  a dual edge, by  $\tilde{f}$  a dual face, and by  $\tilde{c}$  a dual cell. With a subscript we indicate the corresponding (unique) geometric element of the primal grid, thus, the dual of a primal node  $n$  is a dual cell  $\tilde{c}_n$ , the dual of a primal edge  $e$  is a dual face  $\tilde{f}_e$ , the dual of a primal face  $f$  is a dual edge  $\tilde{e}_f$ , and the dual of a primal cell  $c$  is the dual node denoted by  $\tilde{n}_c$ .

To describe the geometric elements of the pair of grids, we introduce a Cartesian system of coordinates with specified origin and we denote by  $\mathbf{x} = (x_1, x_2, x_3)^T$  the coordinates of its generic point. Let us first describe the construction of the *restriction of dual geometric elements to a single cell  $c$* , as the one pictured in Fig. 1. The geometric construction of the dual nodes, dual edges, dual faces, and dual cells is based on the barycentric subdivision as follows. To begin with, let us define the barycenters of geometric elements. The barycenters of an edge  $e$ , of a face  $f$ , and of a cell  $c$  are defined, respectively, as follows

$$\mathbf{b}_e = \frac{1}{|e|} \int_e \mathbf{x} \, dl, \quad \mathbf{b}_f = \frac{1}{|f|} \int_f \mathbf{x} \, dS, \quad \mathbf{b}_c = \frac{1}{|c|} \int_c \mathbf{x} \, dV, \quad (3)$$

where  $|e|$ ,  $|f|$ , and  $|c|$  denote the length of edge  $e$ , the area of face  $f$ , and the volume of cell  $c$ , respectively.

A dual node  $\tilde{n}_c$  is the barycenter  $\mathbf{b}_c$  of  $c$ . The restriction of a dual edge to  $c$ , denoted by  $\tilde{e}_{f|c}$ , is a segment which joins  $\tilde{n}_c$  with the barycenter  $\mathbf{b}_f$  of a primal face  $f$ . The restriction of a dual face to  $c$ , denoted by  $\tilde{f}_{e|c}$ , is a quadrilateral plane surface whose vertices are  $\tilde{n}_c$ , the barycenters of the pair of primal faces having a primal edge  $e$  in common, and the barycenter  $\mathbf{b}_e$  of  $e$ . The restriction of a dual cell to  $c$ , denoted by  $\tilde{c}_{n|c}$ , is a region whose boundary is made of a disjoint union of restrictions of dual faces to  $c$ , corresponding to edges of  $c$  having the node  $n$  in common, and portions of faces of  $c$  having the node  $n$  in common. The restrictions of dual edge  $\tilde{e}_{f|c}$ , dual face  $\tilde{f}_{e|c}$ , and dual cell  $\tilde{c}_{n|c}$  to  $c$  are endowed with outer orientation [1,12], in such a way that each of the pairs  $(f, \tilde{e}_{f|c})$ ,  $(e, \tilde{f}_{e|c})$ ,  $(n, \tilde{c}_{n|c})$  are oriented according to the right-hand rule.

To each of the following geometric elements  $e$ ,  $f$ ,  $\tilde{f}_{|c}$ , and  $\tilde{e}_{|c}$  of the primal or of the dual grid, we associate their corresponding vectors  $\mathbf{e}$ ,  $\mathbf{f}$ ,  $\mathbf{f}_{|c}$ , and  $\mathbf{e}_{|c}$ , respectively. Each of these vectors, will be represented with a column array of its cartesian components. Vector  $\mathbf{e}$  is the edge vector associated with edge  $e$ ; for example  $\mathbf{e} = (e_1, e_2, e_3)^T = \mathbf{n}_i - \mathbf{n}_j$ ,  $i \neq j$ , where  $\mathbf{n}_i$  and  $\mathbf{n}_j$  are the coordinates of the boundary nodes of  $e$ .  $\mathbf{t}_e$  represents the unit vector parallel to  $e$  and having the same orientation in such a way that  $\mathbf{e} = |e|\mathbf{t}_e$ . Vector  $\mathbf{f}$  is the face vector associated with face  $f$  defined as  $\mathbf{f} = |f|\mathbf{n}_f$ , where  $\mathbf{n}_f$  is the unit vector orthogonal to  $f$  and oriented according to the right-hand rule. In a similar way, vector  $\tilde{\mathbf{e}}_{f|c}$  is the edge vector associated with the dual edge  $\tilde{e}_{f|c}$ ; for instance,  $\tilde{\mathbf{e}}_{f|c} = \mathbf{b}_f - \tilde{\mathbf{n}}_c$ , see Fig. 1b. Vector  $\tilde{\mathbf{f}}_{e|c}$  is the face vector associated with the dual face  $\tilde{f}_{e|c}$ . Face vector  $\tilde{\mathbf{f}}_{e|c}$  is equal to  $\tilde{\mathbf{f}}_{e|c} = \frac{1}{2}(\tilde{\mathbf{e}}_{f_i|c} \times (\mathbf{b}_e - \tilde{\mathbf{n}}_c) - \tilde{\mathbf{e}}_{f_j|c} \times (\mathbf{b}_e - \tilde{\mathbf{n}}_c))$  with  $f_i, f_j$  the unique faces of  $c$  such that  $e = f_i \cap f_j$ , with  $i = i(e)$  and  $j = j(e)$ , and oriented in such a way that the expression induces the correct orientation on  $\tilde{\mathbf{f}}_{e|c}$ , see Fig. 1b.

Now, starting from the definition of dual geometric elements restricted to a single cell  $c$ , we introduce dual edges, dual faces, and dual cells of the barycentric dual grid. We define a dual edge  $\tilde{e}_f$  as the piecewise segment made of the union of the two segments  $\tilde{e}_{f|c_1}$  and  $\tilde{e}_{f|c_2}$  with  $c_1, c_2 \in C(f)$ . The corresponding vector  $\tilde{\mathbf{e}}_f$  is defined as follows  $\tilde{\mathbf{e}}_f := \sum_{c \in C(f)} \tilde{\mathbf{e}}_{f|c}$ . A dual face  $\tilde{f}_e$  is defined as the polyhedral surface made of the union of all  $\tilde{f}_{e|c}$  with  $c \in C(e)$ . The corresponding dual face vector  $\tilde{\mathbf{f}}_e$  is defined as follows  $\tilde{\mathbf{f}}_e := \sum_{c \in C(e)} \tilde{\mathbf{f}}_{e|c}$ . A dual cell  $\tilde{c}_n$  is defined as the region made of the union of all  $\tilde{c}_{n|c}$  with  $c \in C(n)$ , that is  $\tilde{c}_n := \bigcup_{c \in C(n)} \tilde{c}_{n|c}$ . Note that the boundary of each dual cell decomposes into dual edges and dual faces.

Let  $X$  be either  $E$  or  $F$  and let  $\tilde{X}$  be either  $\tilde{E}$  or  $\tilde{F}$ . Given a non-empty region  $\Omega' \subset \Omega$  of  $\mathbb{R}^3$ , we define the subsets of geometric elements of  $X$  and  $\tilde{X}$  restricted to  $\Omega'$  to be

$$X_{|\Omega'} := \{x \cap \Omega' \mid x \in X\}, \quad \tilde{X}_{|\Omega'} := \{\tilde{x} \cap \Omega' \mid \tilde{x} \in \tilde{X}\}. \quad (4)$$

In what follows,  $\Omega'$  will be a cell  $c \in C$ , a dual cell  $\tilde{c} \in \tilde{C}$ , or the restriction of the dual cell  $\tilde{c}$  to the cell  $c$ ,  $\tilde{c}_{|c}$ . For instance,  $E_{|\tilde{c}} := \{e \cap \tilde{c} \mid e \in E\}$  contains the restriction of primal edges to the dual cell  $\tilde{c}$ . Note that  $E_{|c}$  and  $F_{|c}$

coincide with the subsets of all edges and faces of  $c$ . Similarly,  $\tilde{E}_{|\tilde{c}}$  and  $\tilde{F}_{|\tilde{c}}$  coincide with the subsets of all dual edges and dual faces of  $\tilde{c}$ .

Finally, we introduce the corresponding matrices  $\mathbb{X}_{|\Omega'}$  and  $\tilde{\mathbb{X}}_{|\Omega'}$  whose rows collect geometric vectors associated with elements of the sets in (4). For instance, rows of  $\mathbb{E}_{|\tilde{c}}$  contain edge vectors associated with edges in  $E_{|\tilde{c}}$ .

## 2.2. Degrees of freedom and reconstruction operators

The physical variables, often referred to as *degrees of freedom* (DoFs), involved in a discrete formulation of a computational problem, have a precise association with the geometric elements of the pair of staggered primal–dual grids [1,12]. DoFs are obtained by evaluating the physical vector fields on the geometric elements of the grid, and are collected in finite dimensional arrays of real numbers endowed with a natural vector space structure. The operation of translating a vector field into the corresponding DoFs is performed by means of integration using the so-called *de Rham maps* [13,10]. We focus on DoFs attached to edges and faces of the primal and dual grid. We denote by  $\mathcal{E}$ ,  $\mathcal{F}$ ,  $\tilde{\mathcal{E}}$ , and  $\tilde{\mathcal{F}}$  the vector space of edge DoFs, face DoFs, dual edge DoFs, and dual face DoFs, respectively. Given a vector field  $\mathbf{u}$ , we denote by  $\mathbf{u}^{\mathcal{E}}$ ,  $\mathbf{u}^{\mathcal{F}}$ ,  $\mathbf{u}^{\tilde{\mathcal{E}}}$ , and  $\mathbf{u}^{\tilde{\mathcal{F}}}$  the DoFs of  $\mathbf{u}$  in  $\mathcal{E}$ ,  $\mathcal{F}$ ,  $\tilde{\mathcal{E}}$ , and  $\tilde{\mathcal{F}}$ , respectively.

We introduce the following specialized definitions of the de Rham maps, which follow as a special case of the general definition applied to the vector subspace of *constant vector fields*. Let us consider a constant vector field  $\mathbf{u}$ . Given a non-empty region  $\Omega' \subset \Omega$ , we define the corresponding subset of DoFs of  $\mathbf{u}$  *restricted to the region*  $\Omega'$  as follows

$$\mathbf{u}_{|\Omega'}^{\mathcal{E}} = \mathbb{E}_{|\Omega'} \mathbf{u}, \quad \mathbf{u}_{|\Omega'}^{\mathcal{F}} = \mathbb{F}_{|\Omega'} \mathbf{u}, \quad \mathbf{u}_{|\Omega'}^{\tilde{\mathcal{E}}} = \tilde{\mathbb{E}}_{|\Omega'} \mathbf{u}, \quad \mathbf{u}_{|\Omega'}^{\tilde{\mathcal{F}}} = \tilde{\mathbb{F}}_{|\Omega'} \mathbf{u}. \quad (5)$$

A *reconstruction operator* is defined as a formal inverse of the de Rham map restricted to the vector subspace of constant vector fields. Thus, reconstruction operators map DoFs to a constant vector field defined on a cell. In what follows, we will always refer to local reconstruction operators in such a way that the domain of definition of the reconstructed vector field is a given cell.

## 3. Mass matrices

In compatible numerical methods, it is always necessary to map DoFs attached to geometric elements of the primal grid to DoFs attached to geometric elements of the dual grid, or viceversa [1,11,12]. This process of converting one type of DoFs into another is performed by the so-called *discrete Hodge star operator* [13]. This operator is required to satisfy a set of properties. A first requirement is *consistency*, which will be specified in the next subsections, and *symmetry*. In addition, we require *positive definiteness*, which assures the stability of the resulting numerical scheme.

One possible method for explicitly converting one type of DoFs into another is to reconstruct the polynomial vector field in each cell starting from DoFs attached to a set of geometric elements, and then project this vector field on the corresponding geometric elements related by duality. We will follow this principle to design both types of local mass matrices, which give a discrete realization of the corresponding discrete Hodge operators.

In this section we show how to construct consistent mass matrices which map DoFs attached to primal geometric elements to DoFs attached to dual geometric elements. Novel proofs are given based on Stokes' theorem.

### 3.1. Primal cell reconstruction

Let us consider a pair of vector fields  $\mathbf{u}, \mathbf{w} : \mathbb{R}^3 \rightarrow \mathbb{R}^3$  along with a scalar field  $w : \mathbb{R}^3 \rightarrow \mathbb{R}$  defined on a given cell  $c$ . The following well-known integration by parts formulas hold

$$\int_c \mathbf{u} \cdot \nabla w \, dV = - \int_c \nabla \cdot \mathbf{u} \, w \, dV + \int_{\partial c} \mathbf{u} \cdot \mathbf{n} \, w \, dS, \quad (6)$$

$$\int_c \mathbf{u} \cdot \nabla \times \mathbf{w} \, dV = \int_c \nabla \times \mathbf{u} \cdot \mathbf{w} \, dV + \int_{\partial c} \mathbf{u} \times \mathbf{n} \cdot \mathbf{w} \, dS. \quad (7)$$

**Theorem 1.** *The following tensor identities hold*

$$\sum_{f \in F(c)} \tilde{\mathbf{e}}_{f|c} \otimes \mathbf{f} = |c| \mathbb{I}_3, \quad (8)$$

$$\sum_{e \in E(c)} \tilde{\mathbf{f}}_{e|c} \otimes \mathbf{e} = |c| \mathbb{I}_3. \quad (9)$$

**Proof.** Let  $\mathbf{u}, \mathbf{w} \in \mathbb{R}^3$ . Thanks to (6), we have

$$\begin{aligned} |c| \mathbf{u} \cdot \mathbf{w} &= \int_c \mathbf{u} \cdot \mathbf{w} dV = \int_c \mathbf{u} \cdot \nabla (\mathbf{w} \cdot (\mathbf{x} - \tilde{\mathbf{n}}_c)) dV \\ &= \int_{\partial c} (\mathbf{u} \cdot \mathbf{n}) (\mathbf{w} \cdot (\mathbf{x} - \tilde{\mathbf{n}}_c)) dS \\ &= \sum_{f \in F(c)} (\mathbf{u} \cdot \mathbf{f}) ((\mathbf{b}_f - \tilde{\mathbf{n}}_c) \cdot \mathbf{w}). \end{aligned} \quad (10)$$

By using the definition of  $\tilde{\mathbf{e}}_{f|c}$ , it follows that

$$|c| \mathbf{u} \cdot \mathbf{w} = \sum_{f \in F(c)} (\mathbf{u} \cdot \mathbf{f}) (\tilde{\mathbf{e}}_{f|c} \cdot \mathbf{w}). \quad (11)$$

Since  $\mathbf{w}$  is arbitrary, (8) is proved.

Thanks to (7), we have

$$\begin{aligned} 2|c| \mathbf{u} \cdot \mathbf{w} &= 2 \int_c \mathbf{u} \cdot \mathbf{w} dV = \int_c \mathbf{u} \cdot \nabla \times (\mathbf{w} \times (\mathbf{x} - \tilde{\mathbf{n}}_c)) dV \\ &= \int_{\partial c} (\mathbf{u} \times \mathbf{n}) \cdot (\mathbf{w} \times (\mathbf{x} - \tilde{\mathbf{n}}_c)) dS \\ &= \sum_{f \in F(c)} \int_f (\mathbf{u} \times \mathbf{n}_f) \cdot (\mathbf{w} \times (\mathbf{x} - \tilde{\mathbf{n}}_c)) dS \\ &= \sum_{f \in F(c)} |f| (\mathbf{u} \times \mathbf{n}_f) \cdot (\mathbf{w} \times \tilde{\mathbf{e}}_{f|c}). \end{aligned} \quad (12)$$

Now, by applying the same argument used in (10) to the vector  $\mathbf{u} \times \mathbf{n}_f$  restricted to  $f$ , we obtain

$$|f| \mathbf{u} \times \mathbf{n}_f = \sum_{e \in E(f)} (\mathbf{u} \cdot \mathbf{e}) (\mathbf{b}_e - \mathbf{p}), \quad (13)$$

where  $\mathbf{p} \in \mathbb{R}^3$  is an arbitrary point. Apply (13) to every face  $f$  in the last expression of (12), choosing the same node  $\tilde{\mathbf{n}}_c$  as the arbitrary point involved in the formula. We obtain

$$\begin{aligned} 2|c| \mathbf{u} \cdot \mathbf{w} &= \sum_{f \in F(c)} \left( \sum_{e \in E(f)} (\mathbf{u} \cdot \mathbf{e}) (\mathbf{b}_e - \tilde{\mathbf{n}}_c) \right) \cdot (\mathbf{w} \times \tilde{\mathbf{e}}_{f|c}) \\ &= \sum_{f \in F(c)} (\tilde{\mathbf{e}}_{f|c} \times \left( \sum_{e \in E(f)} (\mathbf{u} \cdot \mathbf{e}) (\mathbf{b}_e - \tilde{\mathbf{n}}_c) \right)) \cdot \mathbf{w} \\ &= \sum_{e \in E(c)} (\mathbf{u} \cdot \mathbf{e}) ((\tilde{\mathbf{e}}_{f_i|c} \times (\mathbf{b}_e - \tilde{\mathbf{n}}_c) - \tilde{\mathbf{e}}_{f_j|c} \times (\mathbf{b}_e - \tilde{\mathbf{n}}_c)) \cdot \mathbf{w}), \end{aligned} \quad (14)$$

where  $f_i, f_j$  are the unique faces of  $c$  such that  $e = f_i \cap f_j$ , with  $i = i(e)$  and  $j = j(e)$ , and oriented in such a way that they induce opposite orientations on edge  $e$ . Now, dividing by two both members of the last term in (14) and using the definition of  $\tilde{\mathbf{f}}_{e|c}$ , it follows that

$$|c| \mathbf{u} \cdot \mathbf{w} = \sum_{e \in E(c)} (\mathbf{u} \cdot \mathbf{e}) (\tilde{\mathbf{f}}_{e|c} \cdot \mathbf{w}). \quad (15)$$

Since  $\mathbf{w}$  is arbitrary, (9) is proved.  $\square$



### 3.2. Local mass matrices

Let us focus on a cell  $c$  where two pairs of constant vector fields are defined, namely,  $\mathbf{u}$ ,  $\mathbf{v}$  and  $\mathbf{w}$ ,  $\mathbf{z}$ . The two pairs are related by two constitutive relations

$$\mathbf{v} = \mathbb{K}_{1|c} \mathbf{u}, \quad (16)$$

$$\mathbf{z} = \mathbb{K}_{2|c} \mathbf{w}, \quad (17)$$

where  $\mathbb{K}_{1|c}$ ,  $\mathbb{K}_{2|c}$  are two symmetric positive definite matrices of order 3, assumed to be uniform in  $c$ .

Now, let us introduce the restriction of DoFs to  $c$ . In particular, we attach DoFs to geometric elements of the primal and dual grid as follows  $\mathbf{u}_{|c}^{\mathcal{F}} = \mathbb{F}_{|c} \mathbf{u}$ ,  $\mathbf{v}_{|c}^{\tilde{\mathcal{E}}} = \tilde{\mathbb{E}}_{|c} \mathbf{v}$ ,  $\mathbf{w}_{|c}^{\mathcal{E}} = \mathbb{E}_{|c} \mathbf{w}$ ,  $\mathbf{z}_{|c}^{\tilde{\mathcal{F}}} = \tilde{\mathbb{F}}_{|c} \mathbf{z}$ .

The local mass matrix  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$  maps DoFs of  $\mathbf{u}$  attached to faces to DoFs of  $\mathbf{v}$  attached to dual edges of the barycentric dual grid. We say that  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$  is a *consistent mass matrix* if

$$\mathbf{v}_{|c}^{\tilde{\mathcal{E}}} = \mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}} \mathbf{u}_{|c}^{\mathcal{F}} \quad (18)$$

holds exactly for any pair of constant vector fields  $\mathbf{u}$ ,  $\mathbf{v}$  satisfying (16).

Similarly, a local mass matrix  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$  maps DoFs of  $\mathbf{w}$  attached to edges to DoFs of  $\mathbf{z}$  attached to dual faces of the barycentric dual grid. We say that  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$  is a *consistent mass matrix* if

$$\mathbf{z}_{|c}^{\tilde{\mathcal{F}}} = \mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}} \mathbf{w}_{|c}^{\mathcal{E}} \quad (19)$$

holds exactly for any pair of constant vector fields  $\mathbf{w}$ ,  $\mathbf{z}$  satisfying (17).

An efficient recipe to construct consistent and symmetric matrices  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$ ,  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$  combines the geometric identities in Theorem 1 with the uniformity of the vector fields in  $c$ .

By applying Theorem 1 and the definitions of  $\mathbf{v}_{|c}^{\tilde{\mathcal{E}}}$ ,  $\mathbf{u}_{|c}^{\mathcal{F}}$ , we have that

$$\mathbf{v}_{|c}^{\tilde{\mathcal{E}}} = \tilde{\mathbb{E}}_{|c} \mathbf{v} = \tilde{\mathbb{E}}_{|c} \mathbb{K}_{1|c} \mathbf{u} = \tilde{\mathbb{E}}_{|c} \mathbb{K}_{1|c} \frac{1}{|c|} (\tilde{\mathbb{E}}_{|c}^T \mathbb{F}_{|c}) \mathbf{u} = \frac{(\tilde{\mathbb{E}}_{|c} \mathbb{K}_{1|c} \tilde{\mathbb{E}}_{|c}^T)}{|c|} \mathbf{u}_{|c}^{\mathcal{F}}, \quad (20)$$

and hence, it follows that a symmetric and consistent matrix  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$  is given by

$$\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}} = \frac{\tilde{\mathbb{E}}_{|c} \mathbb{K}_{1|c} \tilde{\mathbb{E}}_{|c}^T}{|c|}. \quad (21)$$

Similarly, by applying Theorem 1 and the definitions of  $\mathbf{z}_{|c}^{\tilde{\mathcal{F}}}$ ,  $\mathbf{w}_{|c}^{\mathcal{E}}$ , we have that

$$\mathbf{z}_{|c}^{\tilde{\mathcal{F}}} = \tilde{\mathbb{F}}_{|c} \mathbf{z} = \tilde{\mathbb{F}}_{|c} \mathbb{K}_{2|c} \mathbf{w} = \tilde{\mathbb{F}}_{|c} \mathbb{K}_{2|c} \frac{1}{|c|} (\tilde{\mathbb{F}}_{|c}^T \mathbb{E}_{|c}) \mathbf{w} = \frac{(\tilde{\mathbb{F}}_{|c} \mathbb{K}_{2|c} \tilde{\mathbb{F}}_{|c}^T)}{|c|} \mathbf{w}_{|c}^{\mathcal{E}}, \quad (22)$$

and hence, it follows that a symmetric and consistent matrix  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$  is given by

$$\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}} = \frac{\tilde{\mathbb{F}}_{|c} \mathbb{K}_{2|c} \tilde{\mathbb{F}}_{|c}^T}{|c|}. \quad (23)$$

The matrices  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$  and  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$ , defined in (21) and (23), are symmetric and consistent but are not positive definite. To achieve this, the idea, developed in Lemma 1, is to add to the consistent positive semidefinite matrices  $\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}}$  and  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$  a *stabilization matrix*, which is symmetric and positive semidefinite. The stabilization matrix coincides with the one proposed in the mimetic literature [8].

**Lemma 1.** Let  $m$  be the cardinality of either  $F(c)$  or  $E(c)$ . Let  $\mathbb{K}_{|c}$  be a symmetric and positive definite matrix of order 3. Let  $\alpha = (\alpha_1, \dots, \alpha_{m-3}) \in (\mathbb{R}^+)^{m-3}$  be any  $(m-3)$ -upla of positive real numbers and let  $\mathbb{D}_\alpha$  be the diagonal matrix whose diagonal entries are  $\alpha_1, \dots, \alpha_{m-3}$ . Denote by  $\mathbb{W}_{|c}^{\mathcal{F}}$  and  $\mathbb{W}_{|c}^{\mathcal{E}}$  the matrices whose columns form an orthonormal basis for  $\text{im}(\mathbb{F}_{|c})^\perp$  and  $\text{im}(\mathbb{E}_{|c})^\perp$ , respectively. Then, the following matrices

$$\mathbb{M}_{|c}^{\tilde{\mathcal{E}}\mathcal{F}} := \frac{1}{|c|} \tilde{\mathbb{E}}_{|c} \mathbb{K}_{|c} \tilde{\mathbb{E}}_{|c}^T + \mathbb{W}_{|c}^{\mathcal{F}} \mathbb{D}_\alpha \mathbb{W}_{|c}^{\mathcal{F}^T}, \quad (24)$$

$$\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}} := \frac{1}{|c|} \tilde{\mathbb{F}}_{|c} \mathbb{K}_{|c} \tilde{\mathbb{F}}_{|c}^T + \mathbb{W}_{|c}^{\mathcal{E}} \mathbb{D}_{\alpha} \mathbb{W}_{|c}^{\mathcal{E}T}, \quad (25)$$

are symmetric, consistent and positive definite.

**Proof.** It is sufficient to show that  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{F}}$  is positive definite. Let  $\mathbf{k} \in \mathbb{R}^m$  be such that  $\mathbf{k}^T \mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{F}} \mathbf{k} = 0$ . In order to prove that  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{F}}$  is positive-definite, we have to show that  $\mathbf{k} = \mathbf{0}$ . The condition  $\mathbf{k}^T \mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{F}} \mathbf{k} = 0$  is equivalent to require that  $(\tilde{\mathbb{E}}_{|c}^T \mathbf{k})^T \mathbb{K}_{|c} \tilde{\mathbb{E}}_{|c}^T \mathbf{k} = 0$  and  $(\mathbb{W}_{|c}^{\mathcal{F}T} \mathbf{k})^T \mathbb{D}_{\alpha} (\mathbb{W}_{|c}^{\mathcal{F}T} \mathbf{k}) = 0$ . Since  $\mathbb{K}_{|c}$  is symmetric and positive definite, and each  $\alpha_i$  is positive, the latter condition is in turn equivalent to  $\tilde{\mathbb{E}}_{|c}^T \mathbf{k} = \mathbf{0}$  and  $\mathbf{k} \in \text{im}(\mathbb{F}_{|c})$ . As a consequence,  $\mathbf{k} = \mathbb{F}_{|c} \mathbf{y}$  for some  $\mathbf{y} \in \mathbb{R}^3$  and  $\mathbf{0} = \tilde{\mathbb{E}}_{|c}^T \mathbf{k} = \tilde{\mathbb{E}}_{|c}^T \mathbb{F}_{|c} \mathbf{y} = |c| \mathbf{y}$ . Thus  $\mathbf{k} = \mathbb{F}_{|c} \mathbf{y} = \mathbf{0}$ , as desired.  $\square$

In the general case of a grid made of more than one cell, the corresponding global mass matrices  $\mathbb{M}^{\tilde{\mathcal{F}}\mathcal{F}}$  and  $\mathbb{M}^{\tilde{\mathcal{F}}\mathcal{E}}$  are obtained by assembling, cell by cell, the contributions from the local matrices  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{F}}$  and  $\mathbb{M}_{|c}^{\tilde{\mathcal{F}}\mathcal{E}}$ , respectively.

#### 4. Sparse inverse mass matrices

To derive inverse mass matrices we will mimic the reasoning of Section 3. In particular, we will derive a local mass matrix starting from reconstruction formulas and subsequently construct a global mass matrix by applying a standard assembly process. The construction of inverse mass matrices is carried out only for tetrahedral grids through the introduction of specialized geometric identities. This approach differs from the one described in Section 3. Indeed, the same arguments used in the proof of Theorem 1 cannot be applied to the barycentric dual grid since dual cells have non-planar geometric elements.

##### 4.1. Dual cell reconstruction

In this section, we give novel proofs of reconstruction formulas that provide a constant vector field defined on  $\tilde{c}$  starting from DoFs attached to dual edges and dual faces.

Let  $c$  be a tetrahedron and let  $n$  be one of its nodes. Let  $i \in \{1, 2, 3\}$ . For  $e_i \in E(n) \cap E(c)$ , consider the unique face  $f_i \in F(n) \cap F(c)$  to which  $e_i$  does not belong. In this way, for every  $i \in \{1, 2, 3\}$ , there exists, and is unique,  $r_i \in \{-1, 1\}$  such that  $r_i \mathbf{f}_i \cdot \mathbf{e}_i > 0$ .

**Lemma 2.** *The following tensor identity holds*

$$\sum_{i=1}^3 r_i (\mathbf{f}_i \otimes \mathbf{e}_i) = 3|c| \mathbb{I}_3. \quad (26)$$

**Proof.** Let  $j \in \{1, 2, 3\}$  and let  $\mathbf{n}_{f_j}$  be the unit vector of  $\mathbf{f}_j$ . Let us prove that

$$\sum_{i=1}^3 r_i (\mathbf{f}_i \otimes \mathbf{e}_i) \mathbf{n}_{f_j} = \sum_{i=1}^3 r_i (\mathbf{e}_i \cdot \mathbf{n}_{f_j}) \mathbf{f}_i = 3|c| \mathbf{n}_{f_j} = 3|c| \mathbb{I}_3 \mathbf{n}_{f_j}, \quad (27)$$

from which the claimed identity follows, since the set of vectors  $\mathbf{n}_{f_j}$  are linearly independent. Note that if  $i \neq j$ , then  $\mathbf{e}_i \cdot \mathbf{n}_{f_j} = 0$ . It follows that

$$\sum_{i=1}^3 r_i (\mathbf{e}_i \cdot \mathbf{n}_{f_j}) \mathbf{f}_i = r_j (\mathbf{e}_j \cdot \mathbf{n}_{f_j}) \mathbf{f}_j = r_j (\mathbf{e}_j \cdot \mathbf{n}_{f_j}) |\mathbf{f}_j| \mathbf{n}_{f_j} = 3|c| \mathbf{n}_{f_j}, \quad (28)$$

where we have used the expression of the volume of a tetrahedron [30].  $\square$

Now, we define the restriction of the following *oriented geometric boundary elements*  $s_{n,e|c}$  and  $l_{n,f|c}$  to a tetrahedron  $c$ , which are in bijection with edges  $e \in E(n) \cap E(c)$  and faces  $f \in F(n) \cap F(c)$ , respectively.

We first define  $s_{n,e|c}$ . It decomposes into the union of two triangles,  $t_{n,e|c}^{(1)}$  and  $t_{n,e|c}^{(2)}$ , associated with the two faces  $f_1, f_2 \in F(e) \cap F(c)$ , see Fig. 2(a). The first has the vertices  $\{\mathbf{b}_e, \mathbf{b}_{f_1}, \mathbf{p}_{n,e}\}$ , the second the vertices  $\{\mathbf{b}_e, \mathbf{b}_{f_2}, \mathbf{p}_{n,e}\}$  with node  $\mathbf{p}_{n,e}$  defined as follows

$$\mathbf{p}_{n,e} := \frac{3}{4} \mathbf{n} + \frac{1}{4} \mathbf{n}_1, \quad \mathbf{n}_1 \in N(e) \setminus \{n\}. \quad (29)$$



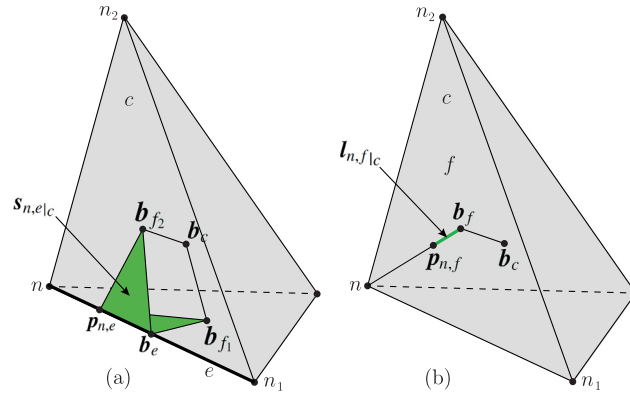


Fig. 2. (a) Geometric construction of  $s_{n,e|c}$ . (b) Geometric construction of  $l_{n,f|c}$ .

By construction, both triangles  $t_{n,e|c}^{(1)}, t_{n,e|c}^{(2)}$  are adjacent to the restriction of the dual face  $\tilde{f}_{e|c}$ . Thus, we orient them in such a way that each pair of surfaces  $t_{n,e|c}^{(1)}, \tilde{f}_{e|c}$  and  $t_{n,e|c}^{(2)}, \tilde{f}_{e|c}$  induce opposite orientations on their common edge [12]. We define  $s_{n,e|c}$  to be the surface made of the union of the two triangles  $t_{n,e|c}^{(1)}, t_{n,e|c}^{(2)}$ . We denote with  $s_{n,e|c}$  the face vector associated with  $s_{n,e|c}$ . It is defined as the sum of the face vectors  $\mathbf{t}_{n,e|c}^{(1)}$  and  $\mathbf{t}_{n,e|c}^{(2)}$ , that is  $s_{n,e|c} := \mathbf{t}_{n,e|c}^{(1)} + \mathbf{t}_{n,e|c}^{(2)}$ . We define  $s_{n,e}$  to be  $s_{n,e} := \sum_{c \in C(e)} s_{n,e|c}$ .

Next, we define  $l_{n,f|c}$  to be the segment joining  $\mathbf{b}_f$  with node  $p_{n,f}$  defined as follows

$$p_{n,f} := \frac{1}{2}\mathbf{n} + \frac{1}{4}\mathbf{n}_1 + \frac{1}{4}\mathbf{n}_2, \quad n_1, n_2 \in N(f) \setminus \{n\}, \quad (30)$$

see Fig. 2(b). By construction,  $l_{n,f|c}$  is adjacent to the restriction of the dual edge  $\tilde{e}_{f|c}$ . Thus, we orient  $l_{n,f|c}$  in such a way the two segments  $l_{n,f|c}, \tilde{e}_{f|c}$  induce opposite orientations on their common node [12]. We denote with  $l_{n,f|c}$  the edge vector associated with  $l_{n,f|c}$ . We define  $l_{n,f}$  to be  $l_{n,f} := \sum_{c \in C(f)} l_{n,f|c}$ .

By mimicking the notation introduced at the end of Section 2.1, we denote by  $\mathbb{S}_{|c}$  and  $\mathbb{L}_{|c}$  the matrices whose rows collect vectors  $s_{n,e}$  and  $l_{n,f}$  with  $e \in E(n)$  and  $f \in F(n)$ , respectively. Similarly, we denote by  $\mathbb{S}_{|\tilde{c}|c}$  and  $\mathbb{L}_{|\tilde{c}|c}$  the matrices whose rows collect vectors  $s_{n,e|c}$  and  $l_{n,f|c}$  with  $e \in E(n) \cap E(c)$  and  $f \in F(n) \cap F(c)$ , respectively.

The geometric boundary elements satisfy the following identities

$$\frac{r_i \mathbf{f}_i}{6} = \tilde{\mathbf{f}}_{e|c} + s_{n,e|c}, \quad (31)$$

$$\frac{r_i \mathbf{e}_i}{4} = \tilde{\mathbf{e}}_{f|c} + l_{n,f|c}, \quad (32)$$

which can be proven after trivial algebraic manipulations that we omit. We point out that, if we require  $s_{n,e|c}$  and  $l_{n,f|c}$  to satisfy (31) and (32), the nodes  $p_{n,e}$  and  $p_{n,f}$  are uniquely identified by (29) and (30). (31) and (32) are the main geometric identities that together with Lemma 2 allow us to prove the following result.

**Lemma 3.** *Let  $c$  be a tetrahedron and let  $n$  be one of its nodes. Let  $\tilde{c}_n$  a dual cell. The following tensor identities hold*

$$\sum_{e \in E(n) \cap E(c)} \mathbf{e}_{|c} \otimes (\tilde{\mathbf{f}}_{e|c} + s_{n,e|c}) = |\tilde{c}_n| \mathbb{I}_3, \quad (33)$$

$$\sum_{f \in F(n) \cap F(c)} \mathbf{f}_{|c} \otimes (\tilde{\mathbf{e}}_{f|c} + l_{n,f|c}) = |\tilde{c}_n| \mathbb{I}_3. \quad (34)$$

**Proof.** Let us apply Lemma 2 to  $c$ . We rewrite (26) as follows

$$\frac{1}{12} \sum_{i=1}^3 r_i (\mathbf{e}_i \otimes \mathbf{f}_i) = |\tilde{c}_n| \mathbb{I}_3, \quad (35)$$

where we have used the fact that  $4|\tilde{c}|_c| = |c|$ , as a result of barycentric subdivision [30]. Now, using (31) it follows that

$$\sum_{i=1}^3 \frac{\mathbf{e}_i}{2} \otimes (\tilde{\mathbf{f}}_{e|_c} + \mathbf{s}_{n,e|_c}) = \sum_{i=1}^3 \frac{\mathbf{e}_i}{2} \otimes \frac{r_i \mathbf{f}_i}{6} = \frac{1}{12} \sum_{i=1}^3 \mathbf{e}_i \otimes (r_i \mathbf{f}_i) = |\tilde{c}|_c| \mathbb{I}_3. \quad (36)$$

This proves (33), since the length of each edge in  $E|_c$  is half of the length a primal edge.

Similarly, using (32) it follows that

$$\sum_{i=1}^3 \frac{\mathbf{f}_i}{3} \otimes (\tilde{\mathbf{e}}_{f|_c} + \mathbf{l}_{n,f|_c}) = \sum_{i=1}^3 \frac{\mathbf{f}_i}{3} \otimes \frac{r_i \mathbf{e}_i}{4} = \frac{1}{12} \sum_{i=1}^3 \mathbf{f}_i \otimes (r_i \mathbf{e}_i) = |\tilde{c}|_c| \mathbb{I}_3. \quad (37)$$

This proves (34), since the area of each face in  $F|_c$  is a third of the area of a primal face.  $\square$

Now we are ready to prove the main result of this section, which is the dual counterpart of Theorem 1. This result is remarkable since dual cells have non-planar geometric elements.

**Theorem 2.** *Let  $\tilde{c}_n$  be a dual cell. The following two assertions hold.*

(1) *If  $\tilde{c}$  does not intersect with  $\partial\Omega$ , then*

$$\sum_{e \in E(n)} \mathbf{e}_{|c} \otimes \tilde{\mathbf{f}}_e = |\tilde{c}| \mathbb{I}_3, \quad (38)$$

$$\sum_{f \in F(n)} \mathbf{f}_{|c} \otimes \tilde{\mathbf{e}}_f = |\tilde{c}| \mathbb{I}_3. \quad (39)$$

(2) *If  $\tilde{c}$  does intersect with  $\partial\Omega$ , then*

$$\sum_{e \in E(n)} \mathbf{e}_{|c} \otimes (\tilde{\mathbf{f}}_e + \mathbf{s}_{n,e}) = |\tilde{c}| \mathbb{I}_3, \quad (40)$$

$$\sum_{f \in F(n)} \mathbf{f}_{|c} \otimes (\tilde{\mathbf{e}}_f + \mathbf{l}_{n,f}) = |\tilde{c}| \mathbb{I}_3. \quad (41)$$

**Proof.** Proof of (38) and (40). Let us apply (33) to every tetrahedron  $c \in C(n)$ . By summing over the set  $C(n)$  we have

$$\sum_{c \in C(n)} \left( \sum_{e \in E(n) \cap E(c)} \mathbf{e}_{|c} \otimes (\tilde{\mathbf{f}}_{e|_c} + \mathbf{s}_{n,e|_c}) \right) = \sum_{c \in C(n)} |\tilde{c}|_c| \mathbb{I}_3 = |\tilde{c}| \mathbb{I}_3. \quad (42)$$

We can rewrite (42) as a sum over the set  $E(n)$  as follows

$$\sum_{c \in C(n)} \left( \sum_{e \in E(n) \cap E(c)} \mathbf{e}_{|c} \otimes (\tilde{\mathbf{f}}_{e|_c} + \mathbf{s}_{n,e|_c}) \right) = \sum_{e \in E(n)} \mathbf{e}_{|c} \otimes \left( \sum_{c \in C(e)} \tilde{\mathbf{f}}_{e|_c} + \mathbf{s}_{n,e|_c} \right). \quad (43)$$

By using the definition of  $\tilde{\mathbf{f}}_e$ ,  $\mathbf{s}_{n,e}$ , and combining the two expressions (42), (43) we obtain

$$\sum_{e \in E(n)} \mathbf{e}_{|c} \otimes (\tilde{\mathbf{f}}_e + \mathbf{s}_{n,e}) = |\tilde{c}| \mathbb{I}_3. \quad (44)$$

This proves (40). To prove (38), note that if  $\tilde{c}$  does not intersect with  $\partial\Omega$ , then

$$\mathbf{s}_{n,e} = \sum_{e \in C(e)} \mathbf{s}_{n,e|_c} = \mathbf{0}, \quad (45)$$

since each  $\mathbf{s}_{n,e|_c}$  decomposes into the union of two triangles and each of them appears in the sum exactly two times and with opposite orientation.

Proof of (39) and (41). Let us apply (34) to every tetrahedron  $c \in C(n)$ . By summing over the set  $C(n)$  we have

$$\sum_{c \in C(n)} \left( \sum_{f \in F(n) \cap F(c)} \mathbf{f}_{|c} \otimes (\tilde{\mathbf{e}}_{f|_c} + \mathbf{l}_{n,f|_c}) \right) = \sum_{c \in C(n)} |\tilde{c}|_c| \mathbb{I}_3 = |\tilde{c}| \mathbb{I}_3. \quad (46)$$

We can rewrite (46) as a sum over the set  $F(n)$  as follows

$$\sum_{c \in C(n)} \left( \sum_{f \in F(n) \cap F(c)} \mathbf{f}_{|c} \otimes (\tilde{\mathbf{e}}_{f|c} + \mathbf{l}_{n,f|c}) \right) = \sum_{f \in F(n)} \mathbf{f}_{|c} \otimes \left( \sum_{c \in C(f)} \tilde{\mathbf{e}}_{f|c} + \mathbf{l}_{n,f|c} \right). \quad (47)$$

By using the definition of  $\tilde{\mathbf{e}}_f$ ,  $\mathbf{l}_{n,f}$ , and combining the two expressions (46), (47) we obtain

$$\sum_{f \in F(n)} \mathbf{f}_{|c} \otimes (\tilde{\mathbf{e}}_f + \mathbf{l}_{n,f}) = |\tilde{c}| \mathbb{I}_3. \quad (48)$$

This proves (41). To prove (39), note that if  $\tilde{c}$  does not intersect with  $\partial\Omega$ , then

$$\mathbf{l}_{n,f} = \sum_{c \in C(f)} \mathbf{l}_{n,f|c} = \mathbf{0}, \quad (49)$$

since each  $\mathbf{l}_{n,f|c}$  appears in the sum exactly two times and with opposite orientation.  $\square$

The results of Theorem 2 cannot be extended to arbitrary polyhedral grids. To see this, it is sufficient to consider a generic 3D polyhedral grid and check that the relations in Theorem 2 are not satisfied. Simple random examples provide readily a counterexample.

#### 4.2. Local inverse mass matrices

Let us focus on a dual cell  $\tilde{c}_n$ . For the sake of clarity, we suppose that  $\tilde{c}$  does not intersect with  $\partial\Omega$ . Otherwise, it is sufficient to repeat the same reasoning using  $\tilde{\mathbb{F}}_{|\tilde{c}} + \mathbb{S}_{|\tilde{c}}$  and  $\tilde{\mathbb{E}}_{|\tilde{c}} + \mathbb{L}_{|\tilde{c}}$  in place of  $\tilde{\mathbb{F}}_{|\tilde{c}}$  and  $\tilde{\mathbb{E}}_{|\tilde{c}}$ , respectively. In  $\tilde{c}$ , two pairs of constant vector fields are defined, namely,  $\mathbf{u}$ ,  $\mathbf{v}$  and  $\mathbf{w}$ ,  $\mathbf{z}$ . The two pairs are related by two constitutive relations

$$\mathbf{v} = \tilde{\mathbb{K}}_{1|\tilde{c}} \mathbf{u}, \quad (50)$$

$$\mathbf{z} = \tilde{\mathbb{K}}_{2|\tilde{c}} \mathbf{w}, \quad (51)$$

where  $\tilde{\mathbb{K}}_{1|\tilde{c}}$ ,  $\tilde{\mathbb{K}}_{2|\tilde{c}}$  are two symmetric positive definite matrices of order 3, assumed to be uniform in  $\tilde{c}$ .

Now, let us introduce the restriction of DoFs to  $\tilde{c}$ . In particular, we attach DoFs to geometric elements of the primal and dual grid as follows  $\mathbf{u}_{|\tilde{c}}^{\tilde{\mathcal{F}}} = \tilde{\mathbb{F}}_{|\tilde{c}} \mathbf{u}$ ,  $\mathbf{v}_{|\tilde{c}}^{\mathcal{E}} = \mathbb{E}_{|\tilde{c}} \mathbf{v}$ ,  $\mathbf{w}_{|\tilde{c}}^{\tilde{\mathcal{E}}} = \tilde{\mathbb{E}}_{|\tilde{c}} \mathbf{w}$ ,  $\mathbf{z}_{|\tilde{c}}^{\mathcal{F}} = \mathbb{F}_{|\tilde{c}} \mathbf{z}$ .

The local mass matrix  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$  maps DoFs of  $\mathbf{u}$  attached to dual faces to DoFs of  $\mathbf{v}$  attached to edges of the primal grid. We say that  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$  is a *consistent mass matrix* if

$$\mathbf{v}_{|\tilde{c}}^{\mathcal{E}} = \mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}} \mathbf{u}_{|\tilde{c}}^{\tilde{\mathcal{F}}} \quad (52)$$

holds exactly for any pair of constant vector fields  $\mathbf{u}$ ,  $\mathbf{v}$  satisfying (50).

Similarly, a local mass matrix  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  maps DoFs of  $\mathbf{w}$  attached to dual edges to DoFs of  $\mathbf{z}$  attached to faces of the primal grid. We say that  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  is a *consistent mass matrix* if

$$\mathbf{z}_{|\tilde{c}}^{\mathcal{F}} = \mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}} \mathbf{w}_{|\tilde{c}}^{\tilde{\mathcal{E}}} \quad (53)$$

holds exactly for any pair of constant vector fields  $\mathbf{w}$ ,  $\mathbf{z}$  satisfying (51).

An efficient recipe to construct a consistent and symmetric matrices  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$ ,  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  combines the geometric identities in Theorem 2 with the uniformity of the vector fields in  $\tilde{c}$ .

By applying Theorem 2, and the definitions of  $\mathbf{v}_{|\tilde{c}}^{\mathcal{E}}$ ,  $\mathbf{u}_{|\tilde{c}}^{\tilde{\mathcal{F}}}$ , we have that

$$\mathbf{v}_{|\tilde{c}}^{\mathcal{E}} = \mathbb{E}_{|\tilde{c}} \mathbf{v} = \mathbb{E}_{|\tilde{c}} \tilde{\mathbb{K}}_{1|\tilde{c}} \mathbf{u} = \mathbb{E}_{|\tilde{c}} \tilde{\mathbb{K}}_{1|\tilde{c}} \frac{1}{|\tilde{c}|} (\mathbb{E}_{|\tilde{c}}^T \tilde{\mathbb{F}}_{|\tilde{c}}) \mathbf{u} = \frac{(\mathbb{E}_{|\tilde{c}} \tilde{\mathbb{K}}_{1|\tilde{c}} \mathbb{E}_{|\tilde{c}}^T)}{|\tilde{c}|} \mathbf{u}_{|\tilde{c}}^{\tilde{\mathcal{F}}}, \quad (54)$$

and hence, it follows that a symmetric and consistent matrix  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$  is given by

$$\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}} = \frac{\mathbb{E}_{|\tilde{c}} \tilde{\mathbb{K}}_{1|\tilde{c}} \mathbb{E}_{|\tilde{c}}^T}{|\tilde{c}|}. \quad (55)$$

Similarly, by applying [Theorem 2](#), and the definitions of  $\mathbf{z}_{|\tilde{c}}^{\mathcal{F}}$ ,  $\mathbf{w}_{|\tilde{c}}^{\tilde{\mathcal{E}}}$ , we have that

$$\mathbf{z}_{|\tilde{c}}^{\mathcal{F}} = \mathbb{F}_{|\tilde{c}} \mathbf{z} = \mathbb{F}_{|\tilde{c}} \tilde{\mathbb{K}}_{2|\tilde{c}} \mathbf{w} = \mathbb{F}_{|\tilde{c}} \tilde{\mathbb{K}}_{2|\tilde{c}} \frac{1}{|\tilde{c}|} (\mathbb{F}_{|\tilde{c}}^T \tilde{\mathbb{E}}_{|\tilde{c}}) \mathbf{w} = \frac{(\mathbb{F}_{|\tilde{c}} \tilde{\mathbb{K}}_{1|\tilde{c}} \mathbb{F}_{|\tilde{c}}^T)}{|\tilde{c}|} \mathbf{w}_{|\tilde{c}}^{\tilde{\mathcal{E}}}, \quad (56)$$

and hence, it follows that a symmetric and consistent matrix  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  is given by

$$\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}} = \frac{\mathbb{F}_{|\tilde{c}} \tilde{\mathbb{K}}_{2|\tilde{c}} \mathbb{F}_{|\tilde{c}}^T}{|\tilde{c}|}. \quad (57)$$

The matrices  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$  and  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  defined in (55) and (57) are symmetric and consistent but are not positive definite. To achieve this, we add a stabilization matrix just as we did at the end of [Section 3.2](#).

**Lemma 4.** *Let  $m$  be the cardinality of either  $F(n)$  or  $E(n)$ . Let  $\tilde{\mathbb{K}}_{|\tilde{c}}$  be a symmetric and positive definite matrix of order 3. Let  $\alpha = (\alpha_1, \dots, \alpha_{m-3}) \in (\mathbb{R}^+)^{m-3}$  be any  $(m-3)$ -upla of positive real numbers and let  $\mathbb{D}_\alpha$  be the diagonal matrix whose diagonal entries are  $\alpha_1, \dots, \alpha_{m-3}$ . Denote by  $\mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{F}}}$  and  $\mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{E}}}$  the matrices whose columns form an orthonormal basis for  $\text{im}(\mathbb{F}_{|\tilde{c}} + \mathbb{S}_{|\tilde{c}})^\perp$  and  $\text{im}(\tilde{\mathbb{E}}_{|\tilde{c}} + \mathbb{L}_{|\tilde{c}})^\perp$ , respectively. Then, the following matrices*

$$\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}} := \frac{1}{|\tilde{c}|} \mathbb{E}_{|\tilde{c}} \tilde{\mathbb{K}}_{|\tilde{c}} \mathbb{E}_{|\tilde{c}}^T + \mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{F}}} \mathbb{D}_\alpha \mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{F}T}}, \quad (58)$$

$$\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}} := \frac{1}{|\tilde{c}|} \mathbb{F}_{|\tilde{c}} \tilde{\mathbb{K}}_{|\tilde{c}} \mathbb{F}_{|\tilde{c}}^T + \mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{E}}} \mathbb{D}_\alpha \mathbb{W}_{|\tilde{c}}^{\tilde{\mathcal{E}T}}, \quad (59)$$

are symmetric, consistent and positive definite.

**Proof.** It is a direct consequence of [Lemma 1](#).  $\square$

In the general case of a dual grid made of more than one dual cell, the corresponding global inverse mass matrices  $\mathbb{M}^{\mathcal{E}\tilde{\mathcal{F}}}$  and  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  are obtained by assembling, dual cell by dual cell, the contributions from the local matrices  $\mathbb{M}_{|\tilde{c}}^{\mathcal{E}\tilde{\mathcal{F}}}$  and  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$ , respectively. However, we need also to take into account the case where  $\tilde{c}$  does intersect with  $\partial\Omega$ . In this case, (40) and (41) imply that, in order to obtain a valid reconstruction, we have to attach DoFs of vector fields also to geometric boundary elements defined by  $s_{n,e}$  and  $l_{n,f}$ . In the numerical examples we show how the proposed sparse inverse mass matrices are used in formulations to solve boundary value problems. In this case, the additional DoFs on the boundary are either known because of boundary conditions or not used in the laws enforced by the formulation. We emphasize that the additional DoFs on the boundary appear as a canonical and necessary construction to fulfill the geometrical identities (31) and (32).

#### 4.3. Handling material parameter discontinuities inside dual cells

In this section we are interested in the discretization of problems with discontinuous material parameters, which may appear for example when inhomogeneous Poisson or wave propagation problems are considered. In order to apply the setting detailed in [Section 4.2](#), a key requirement is that the material parameters are uniform inside dual cells. In this case, discontinuities of the material parameters do not create any complication if the grid is chosen in such a way that the interfaces between different materials are aligned with the boundaries of the dual cells. Hence, in this case, the assignment of the material parameters must be based on the dual cells. This is a common practice in the Finite Volume (FE) literature [20]. We observe that such an assignment is correctly defined since dual cells provide a partition of the computational domain.

However, in most cases, it is desired that the material parameters are constant on cells of the primal grid and that they can be arbitrary discontinuous across interfaces between cells. To each  $c \in C$  is assigned a possibly different material parameter  $\mathbb{K}_{|c}$ . As a result, the material parameters of each dual cell may be no longer uniform. In order to handle this kind of situations, we propose two different strategies which differ on how the vector field is reconstructed inside dual cells. The first strategy is to use a *weighted average* of the material parameters. The reconstructed vector field is uniform and the construction detailed in [Section 4.2](#) can be applied. The second strategy consists of using a *piecewise constant vector field* representation inside dual cells, which is suitable to represent discontinuities of the vector field components due to material parameters discontinuities.

#### 4.3.1. Weighted average

Let us consider a dual cell  $\tilde{c}_n \in \tilde{C}$ . To motivate the definition of the new material parameter associated with  $\tilde{c}$ , let us consider a uniform vector field  $\mathbf{u}$  defined in  $\tilde{c}$ . Next, we define the vector field  $\mathbf{v}$  whose restriction to each cell  $c \in C(n)$  is  $\mathbf{v}|_c := \mathbb{K}_{|c} \mathbf{u}|_c$ . It follows that  $\mathbf{v}$  is a piecewise constant vector field that is spatially constant in each  $c \in C(n)$ . The weighted average of  $\mathbf{v}$  over  $\tilde{c}$  is

$$\frac{1}{|\tilde{c}|} \sum_{c \in C(n)} |\tilde{c}|_c \mathbb{K}_{|c} \mathbf{u} = \left( \frac{1}{|\tilde{c}|} \sum_{c \in C(n)} |\tilde{c}|_c \mathbb{K}_{|c} \right) \mathbf{u}, \quad (60)$$

where we have used the fact that  $\mathbf{u}$  is uniform in  $\tilde{c}$ . To each dual cell  $\tilde{c}$  we assign a new material parameter defined by a weighted average as follows

$$\tilde{\mathbb{K}}_{|\tilde{c}} := \left( \frac{1}{|\tilde{c}|} \sum_{c \in C(n)} |\tilde{c}|_c \mathbb{K}_{|c} \right)^{-1}. \quad (61)$$

We use (61) as the material parameter appearing in the expressions of the local mass matrices in Lemma 4. As a consequence, if the material parameter is constant over the whole computational domain, the same material parameter is assigned to every dual cell.

#### 4.3.2. Piecewise constant vector field

Let us consider a dual cell  $\tilde{c}_n \in \tilde{C}$ . Since material parameters may differ on each  $c \in C(n)$ , we use a piecewise constant vector field representation that is spatially constant on each  $c$ . Let us consider two pairs of piecewise constant vector fields defined on  $\tilde{c}$ , namely,  $\mathbf{u}, \mathbf{v}$  and  $\mathbf{w}, \mathbf{z}$ . The restriction of the vector fields  $\mathbf{u}, \mathbf{v}, \mathbf{w}, \mathbf{z}$  to a cell  $c \in C(n)$  satisfy the following constitutive relations

$$\mathbf{v}|_c = \mathbb{K}_{1|c} \mathbf{u}|_c, \quad (62)$$

$$\mathbf{z}|_c = \mathbb{K}_{2|c} \mathbf{w}|_c, \quad (63)$$

where  $\mathbb{K}_{1|c}, \mathbb{K}_{2|c}$  are two symmetric positive definite matrices of order 3. In addition, let us suppose that  $\mathbf{u}, \mathbf{z}$  and  $\mathbf{v}, \mathbf{w}$  preserve the *normal* and *tangential* components over the boundaries of  $\tilde{c}|_c$ , respectively. Thus, the restriction of DoFs associated with the vector fields defined in  $\tilde{c}$  are well defined. In particular, we attach DoFs to geometric elements of the primal and dual grid as follows  $\mathbf{u}|_{\tilde{c}}^{\mathcal{F}} = \mathbb{F}_{|\tilde{c}} \mathbf{u}, \mathbf{v}|_{\tilde{c}}^{\tilde{\mathcal{E}}} = \tilde{\mathbb{E}}_{|\tilde{c}} \mathbf{v}, \mathbf{w}|_{\tilde{c}}^{\mathcal{E}} = \mathbb{E}_{|\tilde{c}} \mathbf{w}, \mathbf{z}|_{\tilde{c}}^{\tilde{\mathcal{F}}} = \tilde{\mathbb{F}}_{|\tilde{c}} \mathbf{z}$ .

The idea underlying the construction is to mimic the reasoning proposed in Sections 3.2 and 4.2 to construct global mass matrices starting from local mass matrices defined on cells and dual cells. To be more precise, “local” contributions from matrices  $\mathbb{M}_{|\tilde{c}|_c}^{\tilde{\mathcal{F}}\mathcal{E}}$  and  $\mathbb{M}_{|\tilde{c}|_c}^{\mathcal{F}\tilde{\mathcal{E}}}$  restricted to each region  $\tilde{c}|_c$  are assembled to construct “global” mass matrices  $\mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}$  and  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  on the dual cell  $\tilde{c}$ , respectively. The analogy is well defined since the regions  $\tilde{c}|_c$  provide a partition of the cell  $\tilde{c}$  and the material parameters are uniform on each of them. These assumptions allow us to apply to each region  $\tilde{c}|_c$  the construction of a consistent part using the recipe detailed in Section 4.2.

Thus, by applying Lemma 3, a *consistent* mass matrix  $\mathbb{M}_{|\tilde{c}|_c}^{\tilde{\mathcal{F}}\mathcal{E}}$  restricted to  $\tilde{c}|_c$  is given by

$$\mathbb{M}_{|\tilde{c}|_c}^{\tilde{\mathcal{F}}\mathcal{E}} = \frac{(\tilde{\mathbb{F}}_{|\tilde{c}|_c} + \mathbb{S}_{|\tilde{c}|_c}) \mathbb{K}_{|c} (\tilde{\mathbb{F}}_{|\tilde{c}|_c} + \mathbb{S}_{|\tilde{c}|_c})^T}{|\tilde{c}|_c}. \quad (64)$$

A local mass matrix  $\mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}$  is obtained by assembling the local contributions of each  $\mathbb{M}_{|\tilde{c}|_c}^{\tilde{\mathcal{F}}\mathcal{E}}$ .  $\mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}$  constructed in this way is symmetric and positive definite since each matrix (64) has rank 3. Moreover,  $\mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}$  satisfy the following *piecewise consistency property*, that is

$$\mathbf{z}|_{\tilde{c}}^{\tilde{\mathcal{F}}} = \mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}} \mathbf{w}|_{\tilde{c}}^{\mathcal{E}} \quad (65)$$

holds exactly for any pair of piecewise constant vector fields  $\mathbf{w}, \mathbf{z}$  satisfying (63). This follows from the fact that the normal component of  $\mathbf{z}$  is preserved over the boundaries of  $\tilde{c}|_c$  and hence the geometric boundary elements cancel out, as shown in the proof of Theorem 2. We observe that the piecewise consistency property implies the consistency property (52). A local mass matrix  $\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}}$  is obtained by computing the algebraic inverse of  $\mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}$

$$\mathbb{M}_{|\tilde{c}}^{\mathcal{F}\tilde{\mathcal{E}}} = \mathbb{M}_{|\tilde{c}}^{\tilde{\mathcal{F}}\mathcal{E}}^{-1}. \quad (66)$$

In a similar way, by applying [Lemma 3](#), a *consistent* mass matrix  $\mathbb{M}_{|\tilde{c}_{lc}}^{\tilde{\mathcal{E}}\mathcal{F}}$  restricted to  $\tilde{c}_{lc}$  is given by

$$\mathbb{M}_{|\tilde{c}_{lc}}^{\tilde{\mathcal{E}}\mathcal{F}} = \frac{(\tilde{\mathbb{E}}_{|\tilde{c}_{lc}} + \mathbb{I}_{|\tilde{c}_{lc}})\mathbb{K}_{lc}(\tilde{\mathbb{E}}_{|\tilde{c}_{lc}} + \mathbb{I}_{|\tilde{c}_{lc}})^T}{|\tilde{c}_{lc}|}. \quad (67)$$

A local mass matrix  $\mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}$  is obtained by assembling the local contributions of each  $\mathbb{M}_{|\tilde{c}_{lc}}^{\tilde{\mathcal{E}}\mathcal{F}}$ .  $\mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}$  constructed in this way is symmetric and positive definite since each matrix (67) has rank 3. Moreover,  $\mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}$  satisfy the following *piecewise consistency property*, that is

$$\mathbf{v}_{lc}^{\tilde{\mathcal{E}}} = \mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}} \mathbf{u}_{lc}^{\mathcal{F}} \quad (68)$$

holds exactly for any pair of piecewise constant vector fields  $\mathbf{u}, \mathbf{v}$  satisfying (62). This follows from the fact that the tangential component of  $\mathbf{v}$  is preserved over the boundaries of  $\tilde{c}_{lc}$  and hence the geometric boundary elements cancel out, as shown in the proof of [Theorem 2](#). We observe that the piecewise consistency property implies the consistency property (53). A local mass matrix  $\mathbb{M}_{lc}^{\mathcal{F}\tilde{\mathcal{E}}}$  is obtained by computing the algebraic inverse of  $\mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}$

$$\mathbb{M}_{lc}^{\mathcal{F}\tilde{\mathcal{E}}} = \mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}^{-1}. \quad (69)$$

We observe that even if the material parameters are uniform inside  $\tilde{c}$ , using this approach it is in any case necessary to compute the inverses of matrices  $\mathbb{M}_{lc}^{\tilde{\mathcal{F}}\mathcal{E}}$  and  $\mathbb{M}_{lc}^{\tilde{\mathcal{E}}\mathcal{F}}$ . Instead, in [Section 4.2](#) explicit expressions of  $\mathbb{M}_{lc}^{\tilde{\mathcal{F}}\mathcal{E}}$  and  $\mathbb{M}_{lc}^{\mathcal{F}\tilde{\mathcal{E}}}$  are derived.

#### 4.3.3. Hybrid approach to handle material discontinuities inside dual cells

The proposed approach to handle discontinuities of the material parameters inside dual cells is the following. If material parameters are uniform inside a dual cell, then we can apply the construction detailed in [Section 4.2](#). Otherwise, we apply one of the two approaches described in [Sections 4.3.1](#) and [4.3.2](#). We point out that these constructions are applied only to cells which lie at the intersection between regions enclosing different materials. Thus, the required computational effort to compute local inverses in the approach described in [Section 4.3.2](#) is negligible in practice.

It is important to note that, if we want to satisfy exactly a *multi-material patch test* (see [Section 5](#)), it is necessary in our hybrid approach to exploit the construction proposed in [Section 4.3.2](#). This is because, being the reconstructed field piecewise constant, it can represent exactly the discontinuities of the vector field components that appear at interfaces between different materials. Instead, in [Section 4.3.1](#), the reconstructed field is constant.

## 5. Numerical results

Beside all the other applications of sparse inverse mass matrices, to validate the construction proposed in this paper we concentrate on the solution of Poisson boundary value problems formulated with one unknown per element.

We first verified that the proposed technique is able to pass a patch test on grids made by general tetrahedra. Then, a stationary current conduction problem representing a square resistor is solved. The stationary current conduction is a Poisson problem in region  $\Omega$  of the 3-D Euclidean space

$$\nabla \times \mathbf{E} = \mathbf{0}, \quad (70a)$$

$$\nabla \cdot \mathbf{J} = 0, \quad (70b)$$

$$\mathbf{J} = \sigma \mathbf{E}, \quad (70c)$$

where  $\sigma$  is the material parameter electric conductivity,  $\mathbf{E}$  and  $\mathbf{J}$  are the *conservative* electrostatic field and the current density vectors, respectively. The electric conductivity  $\sigma$  is assumed to be a positive scalar value which is piecewise uniform in each material region. The region boundary  $\partial\Omega$  is partitioned into a set of  $N^i$  surfaces of perfect insulators  $\partial\Omega_k^i$ , and a set of  $N^c + 1$  disjoint equipotential surfaces of perfect conductors  $\partial\Omega_k^c$  (usually called electrodes):

$$\partial\Omega = \sum_{k=1}^{N^i} \partial\Omega_k^i + \sum_{k=0}^{N^c} \partial\Omega_k^c. \quad (71)$$



Electrode  $\partial\Omega_0^c$  is considered as reference for all the voltages of the remaining electrodes, that are supposed to be assigned.  $\mathbf{J} \cdot \mathbf{n}_{\partial\Omega} = 0$  is set as boundary conditions (b.c.) on each  $\partial\Omega_k^i$ , where  $\mathbf{n}_{\partial\Omega}$  is the outwards oriented normal unit vector of  $\partial\Omega$ .

Since  $\nabla \times \mathbf{E} = \mathbf{0}$  and the electrostatic field is a conservative field, we introduce the scalar potential  $U$  such that  $\mathbf{E} = -\nabla U$ . Similarly, since  $\nabla \cdot \mathbf{J} = 0$ , we introduce a vector potential  $\mathbf{T}$  such that  $\mathbf{J} = \nabla \times \mathbf{T}$ .

### 5.1. Survey of standard formulations

There are several formulations to numerically solve Poisson problems like (70a)–(70b)–(70c). The most common one is the FEM formulation based on the scalar potential  $SP$  expanded with the classical nodal basis functions. With this formulation the unknowns  $U^{\mathcal{N}}$ , are the DoFs of the scalar potentials  $U$  sampled on the grid nodes. The voltages  $\mathbf{E}^{\mathcal{E}} = -\mathbb{G}U^{\mathcal{N}}$ , are associated with the grid edges, where  $\mathbb{G}$  is the edge-node incidence matrix. Finally, the DoFs of  $\mathbf{J}$  are attached to dual faces as [11] shows. Faraday's law (70a) is enforced implicitly by the scalar potential, whereas the solenoidality of the current (70b) is enforced on the boundary of each dual cell by the linear system, see for example [31].

Other possibilities emerge as we exchange the association of physical variables between the primal and dual grids. In *complementary* formulations the DoFs of  $\mathbf{J}$  are attached to faces of the primal grid, whereas the scalar potential DoFs are attached to dual nodes. Consequently, the voltages  $\mathbf{E}^{\tilde{\mathcal{E}}} = -\mathbb{D}^T U^{\tilde{\mathcal{N}}}$ , are associated with dual edges. There are two methods to obtain a complementary formulation. The first *complementary* formulation  $VP$  fulfill (70b) by using the electric vector potential  $\mathbf{T}$  [31] such that  $\mathbf{J}^{\mathcal{F}} = \mathbb{C}\mathbf{T}^{\mathcal{E}}$  (We remark that in general the so-called *relative cohomology theory* is required to present this formulation, see [31]). We note that in the  $VP$  formulation the role of physical laws are exchanged w.r.t. the  $SP$  formulation since Faraday's law (70a) is enforced with a linear system in the  $VP$  formulation.

There is a second method to produce complementary formulations that we call *complementary-dual*: they are complementary formulation but still use the scalar potential  $U^{\tilde{\mathcal{N}}}$ , which is sampled on grid dual nodes, one-to-one with grid cells. An effective complementary-dual formulation, which is algebraically equivalent to the  $VP$  formulation, is the mixed-hybrid  $MH$  formulation [32,33].

### 5.2. Formulations with one unknown per element

There are other complementary-dual formulations, much less explored in the literature, that feature one unknown per element [20,34,22]. As pointed out in [34], there has been a long-standing interest to reduce the system to one potential value per element, to reduce unknowns and obtain a positive-definite system.

In [34] two approaches are proposed. The first one requires the solution of local problems on patches of elements to produce local flux expressions. The second one is based on the use of the algebraic inverse of the mass matrix and has been deemed as impossible in [34] because the “the inverse of a mass matrix is not sparse”. We follow the latter approach enabled by the efficient construction of sparse inverse of mass matrices proposed for the first time in this paper. To show the details, let us consider the dual scalar potential formulations  $DSP$  obtained by writing the problem in the geometric framework [11] as

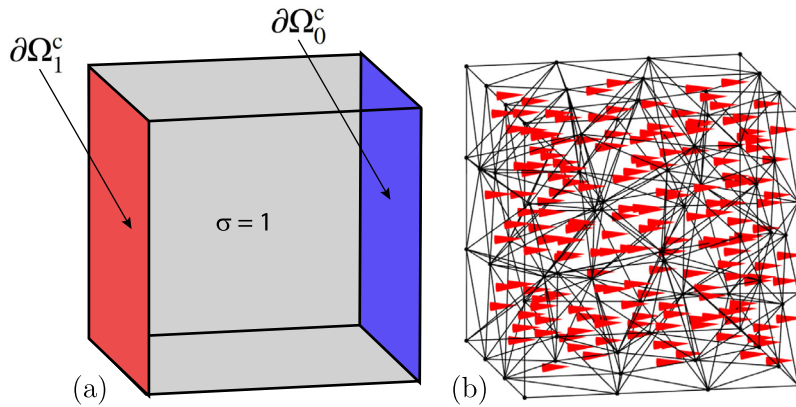
$$\mathbb{C}^T \mathbf{E}^{\tilde{\mathcal{E}}} = \mathbf{0}, \quad (72)$$

$$\mathbb{D} \mathbf{J}^{\mathcal{F}} = \mathbf{0}, \quad (73)$$

$$\mathbf{J}^{\mathcal{F}} = \mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}} \mathbf{E}^{\tilde{\mathcal{E}}}, \quad (74)$$

where  $\mathbb{D}$  is the cell-face incidence matrix,  $\mathbb{C}$  is the face-edge incidence matrix and  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  is the inverse mass matrix of  $\mathbb{M}^{\tilde{\mathcal{E}}\mathcal{F}}$ , which maps DoFs of  $\mathbf{E}$  attached to dual edges to DoFs of  $\mathbf{J}$  attached to primal faces. Thus, we can interpret  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  as a *dual conductance matrix*. To implicitly satisfy (72), the scalar potential  $U^{\tilde{\mathcal{N}}}$  in the dual nodes is introduced through

$$\mathbf{E}^{\tilde{\mathcal{E}}} = -\mathbb{D}^T U^{\tilde{\mathcal{N}}} + \mathbf{E}_s^{\tilde{\mathcal{E}}}, \quad (75)$$



**Fig. 3.** (a) The definition of a patch test as the solution of an electric conduction problem inside a planar resistor. (b) All mentioned formulations are able to retrieve the analytical solution up to machine precision or iterative solver tolerance.

where  $\mathbf{E}_s^{\tilde{\mathcal{E}}}$  is introduced to take into account Dirichlet b.c., so that  $\mathbb{C}^T \mathbf{E}_s^{\tilde{\mathcal{E}}} = \mathbf{0}$  [23,33]. Its construction is straightforward and detailed in [33]. By substituting (74) and (75) into (73), one gets

$$(\mathbb{D}\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}\mathbb{D}^T)U^{\tilde{\mathcal{N}}} = \mathbb{D}\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}\mathbf{E}_s^{\tilde{\mathcal{E}}}, \quad (76)$$

having the scalar potential on dual nodes as unknowns.

A symmetric and positive-definite  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  may be computed as  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}} = \mathbb{M}^{\tilde{\mathcal{E}}\mathcal{F}}^{-1}$ , where the Raviart–Thomas mass matrix  $\mathbb{M}^{\tilde{\mathcal{E}}\mathcal{F}}$  can be interpreted as a *resistance mass matrix*. Yet, the obtained matrix  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  using this recipe is fully populated, hence not usable in practice [34]. Another solution is enabled by the geometric construction of  $\mathbb{M}^{\mathcal{F}\tilde{\mathcal{E}}}$  by using (59), which produces a symmetric and positive definite matrix for any tetrahedral grid given as input. We point out that the additional voltages DoFs on the boundary are zero on the electrodes, because electrodes are equipotential by hypothesis. The other additional voltages DoFs on the boundary are not needed because the boundary conditions are applied on the currents of the corresponding boundary faces.

We remark that another formulation arises that we may call dual vector potential *DVP*, which is the dual of the *VP* formulation. I.e. one obtains a system in which the unknowns are the DoFs of the vector potential attached on dual edges. This formulation is not presented in detail since it has been verified that it is far to be competitive with the others in terms of computational efficiency.

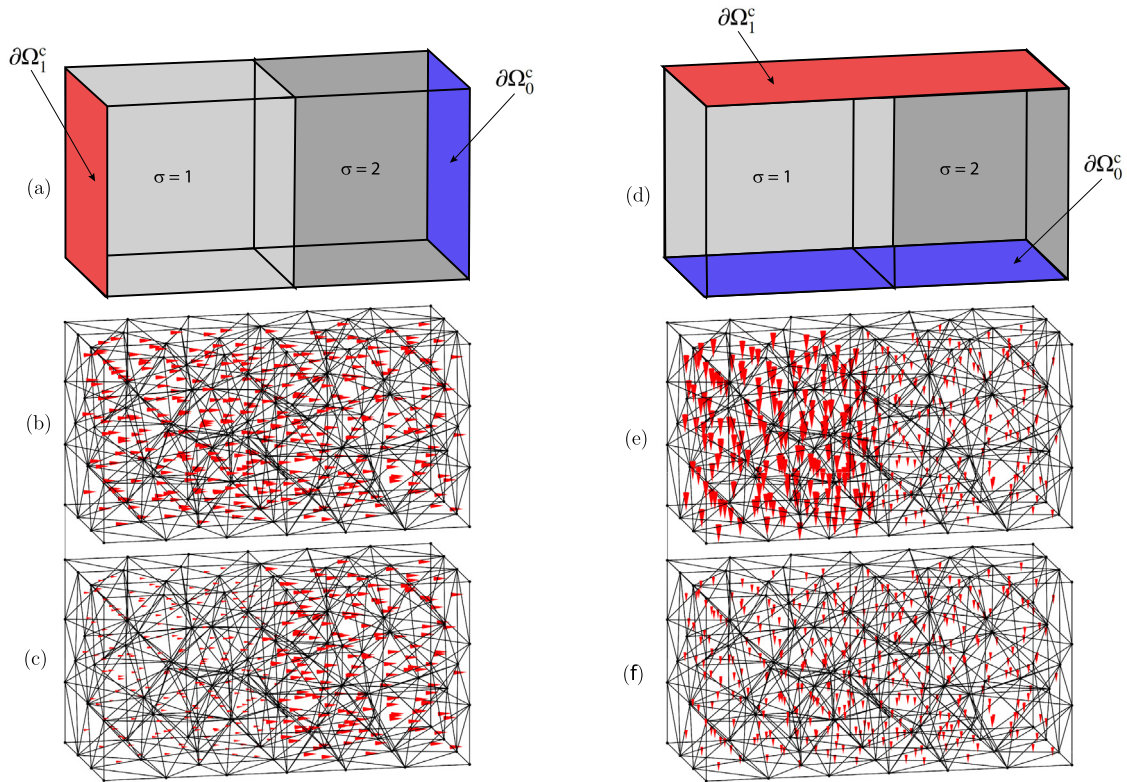
### 5.3. Classical patch tests

The patch tests are Poisson problems designed in such a way that their analytical solution is uniform. By interpreting the Poisson problems as direct current conduction problems, a simple way to produce a patch test is to consider a planar resistor, as described in Fig. 3(a). It has been verified that the *SP*, *VP* = *MH*, and *DSP* formulations produce the analytical solution as represented in Fig. 3(b).

### 5.4. Multi-material patch tests

The multi-material patch tests are direct current conduction problems designed in such a way that their solution is *piecewise uniform*. A simple way to produce multi-material patch test is to consider a resistor with two conductors with different material properties placed in *series* or in *parallel* as described in Figs. 4(a) and 4(d). A voltage of  $u = 1$  V is enforced between the reference electrode  $\partial\Omega_0^c$  and the electrode  $\partial\Omega_1^c$ .

In the first multi-material patch test, two different materials with different material properties are placed in *series*. From the result represented in Fig. 4(b–c), we conclude that the tangential component of the electric field  $\mathbf{E}$  is conserved across the material interface, whereas the tangential component of the current density field  $\mathbf{J}$  jumps. This result holds for all formulations.



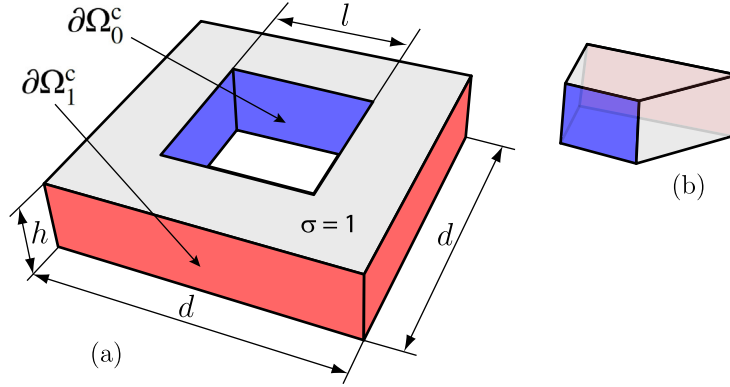
**Fig. 4.** (a) The *series* multi-material patch test. (b) Electrostatic field  $\mathbf{E}$  obtained with the series. (c) Current density field  $\mathbf{J}$  obtained with the series. (d) The *parallel* multi-material patch test. (e) Electric field  $\mathbf{E}$  obtained with the parallel. (f) Current density field  $\mathbf{J}$  obtained with the parallel.

In the second multi-material patch test, two different materials with different material properties are placed in *parallel*. From the result represented in Fig. 4(e–f), we conclude that the normal component of the current density  $\mathbf{J}$  is conserved across the material interface, whereas the normal component of the electric field  $\mathbf{E}$  jumps. This result holds for all formulations.

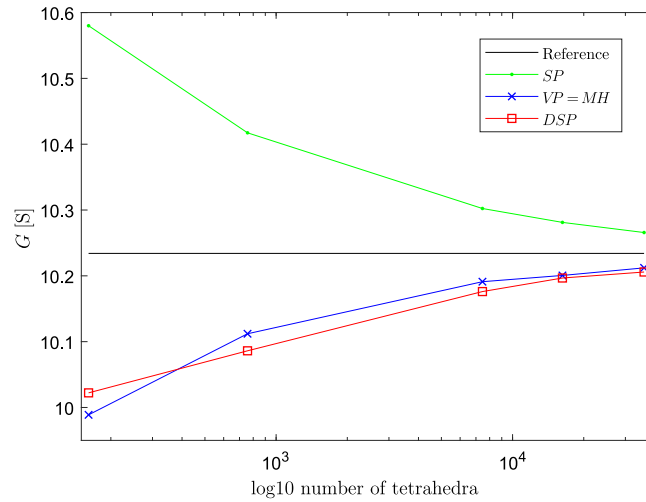
### 5.5. Square resistor benchmark

As a more complicated example, we compute the conductance  $G$  of a square resistor [33], see Fig. 5(a). A voltage of  $u = 1$  V is enforced between the reference electrode  $\partial\Omega_0^c$  and the electrode  $\partial\Omega_1^c$ , which are placed in the two lateral surfaces of the solid square torus. The conductor placed inside the solid torus has an electrical conductivity  $\sigma = 1$  S/m. This problem, like most industrial problems, exhibits a singular analytical solution. Yet, the analytical solution is available ( $G = 10.23409256$  S). The conductance  $G$  is extracted by computing the total dissipated power  $P = Gu^2 = \int_{\Omega} \frac{|\mathbf{J}|^2}{\sigma}$  or alternatively by using the Ohm's law  $G = i/u$ , where  $i$  is the current that flows between the two electrodes.

We recall that the *VP* and *MH* formulations produce the same results given that they are algebraically equivalent [33]. The conductance of the square resistor has been evaluated by the *SP*, *VP* = *MH*, and *DSP* formulations on refined grids. The results are collected in Fig. 6. First, it is interesting to note that the results relative to the scalar potential *SP* and to the vector potential *VP* or mixed-hybrid *MH* formulations provide, respectively, the upper and lower bounds for the conductance [35,31]. This property is called *complementarity* in the computational electromagnetics community [10,33].



**Fig. 5.** (a) The geometry of the square resistor benchmark ( $h = 1$  m,  $d = 4$  m, and  $l = 2$  m). The two electrodes are depicted in red and blue. (b) Thanks to the symmetry, the computational domain has been reduced to one in eight of the resistor. (For interpretation of the references to color in this figure legend, the reader is referred to the web version of this article.)



**Fig. 6.** Results for the square resistor benchmark.

## 6. Conclusions

In this paper, we have proposed a new construction of sparse inverse edge and face mass matrices for arbitrary tetrahedral grids. The proposed framework is based on novel geometric reconstruction formulas, from which, local mass matrices are defined as the sum of a consistent and a stabilization part. The consistent part is geometrically and explicitly defined, thus providing a sensible speedup and an easier implementation, whereas the stabilization part is constructed according to a well-established design strategy. The resulting expressions of both types of local mass matrices are highly symmetric, thus highlighting the duality relationship between the pair of grids. Global inverse mass matrices are constructed using a standard assembly process. In this way, sparse matrix representations of the inverse mass matrices are obtained. A key aspect of the proposed construction is the usage of the barycentric dual grid, which can be defined for arbitrary tetrahedral grids. Two different strategies are proposed to deal with the case of problems involving discontinuous material parameters for the case where the interfaces are aligned with the tetrahedral grid, and a hybrid approach between the two is employed in practical applications. We tested the newly derived inverse mass matrices by solving a Poisson problem and we verified that patch test is passed, even when material discontinuities are present. A real case problem involving a square resistor benchmark is analyzed. Given that reconstruction formulas in each dual cell provide a first-order approximation of a vector field, we expect that first-order error estimates can be derived for the numerical scheme.

## Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

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