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Mimetic Finite Difference Methods

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From joint works with K. Lipnikov, M. Shashkov, and V. Simoncini

INdAM Workshop on Multiscale Problems

Cortona, September 18, 2006

ESSENTIAL REFERENCES

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PLAN OF THE TALK

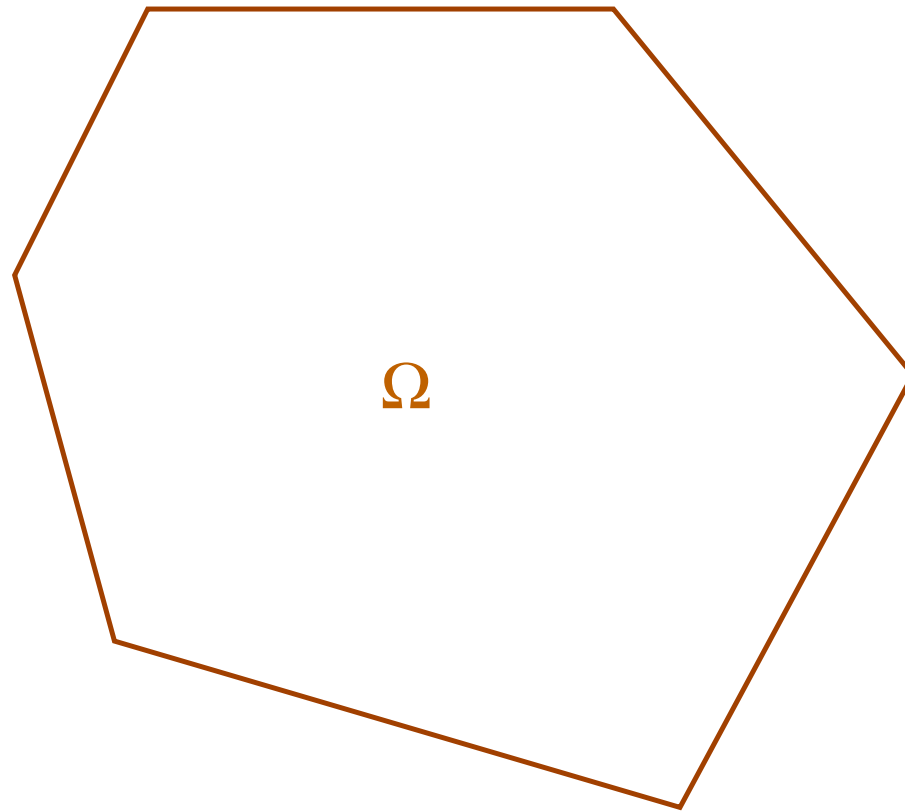
- Darcy's law
- Inner products
- Curved Faces
- Numerical Experiments
- Conclusions
- Appendix (in whatever time is left)

DARCY'S LAW

- $p =$ pressure
- $\mathbf{u} =$ velocities
- $b =$ source
- $\mathbb{K} =$ material-dependent (full) tensor
- $\mathbf{u} = -\mathbb{K}\nabla p$ (Constitutive Equation)
- $\operatorname{div} \mathbf{u} = b$ (Conservation Equation)

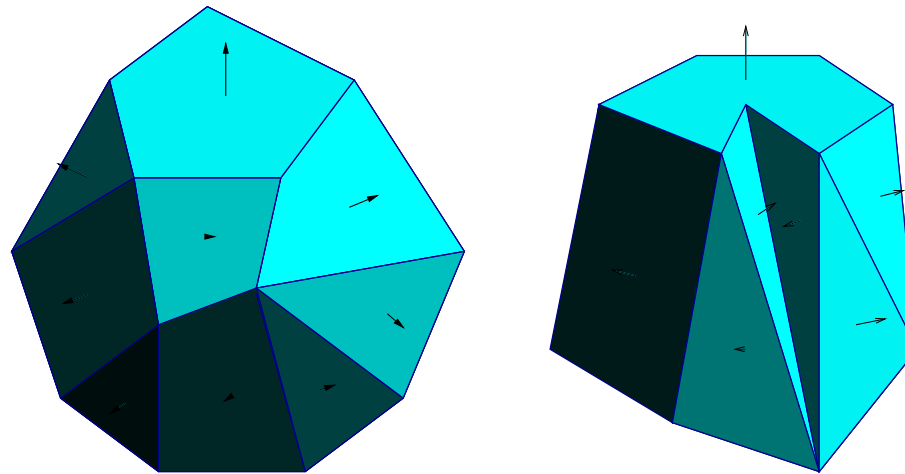
$$\begin{aligned} -\operatorname{div}(\mathbb{K}\nabla p) &= b && \text{in } \Omega, \\ p &= 0 && \text{on } \partial\Omega, \quad \text{for simplicity.} \end{aligned}$$

THE DOMAIN Ω

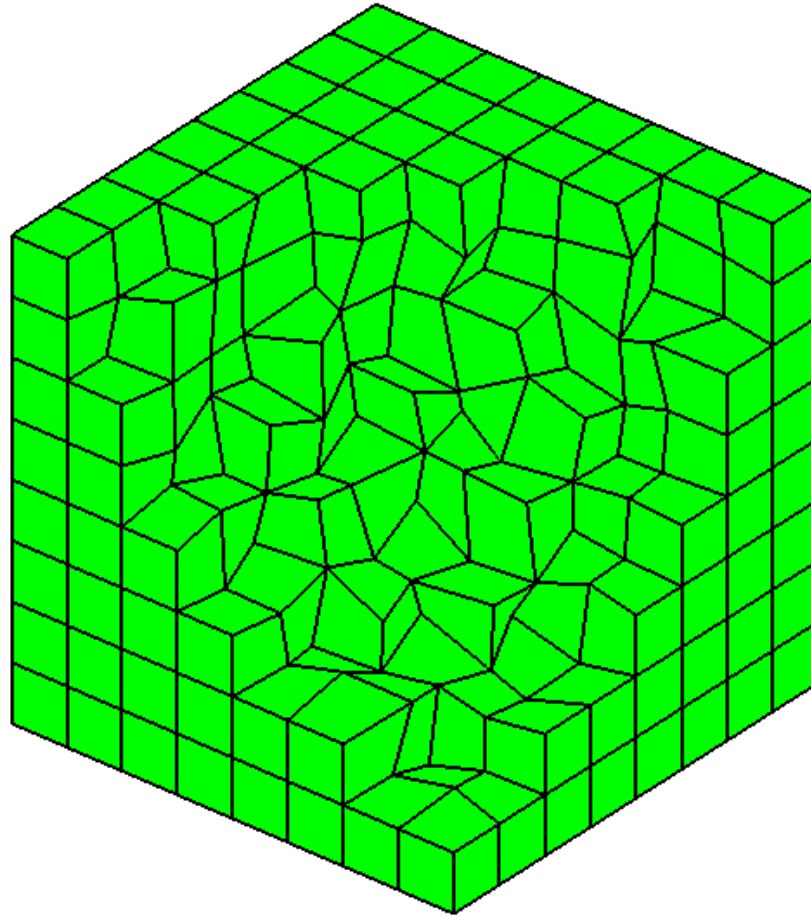


GENERALITY OF THE APPROACH

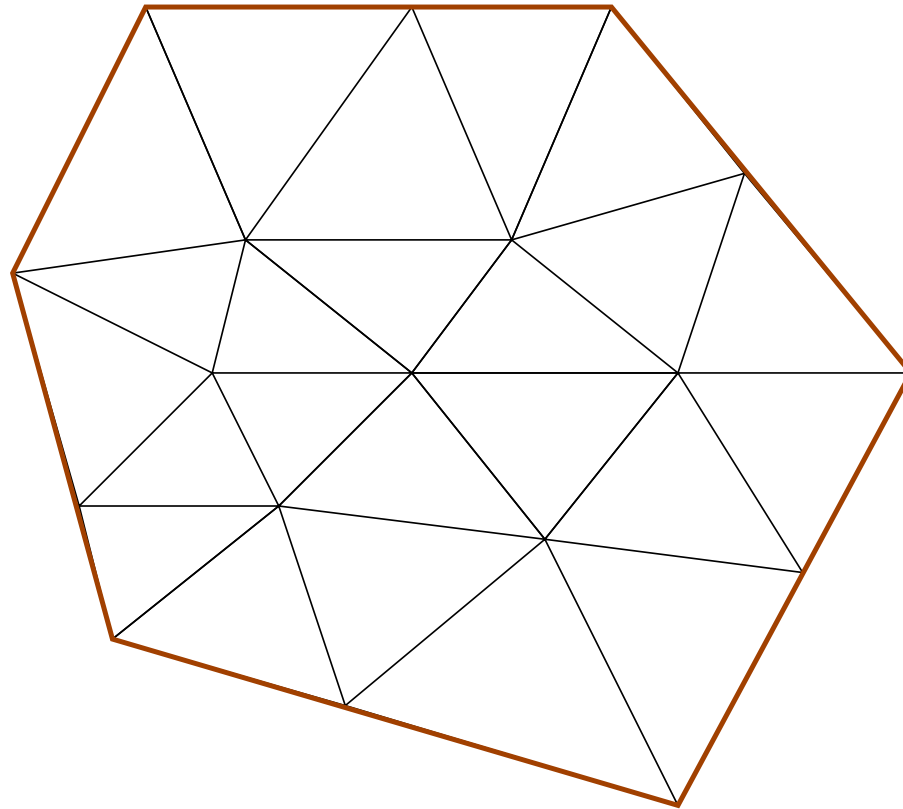
Note that, in the following discussion, many *figures* will be 2-dimensional. This corresponds to a *limitation of the speaker* and *not of the method*. Indeed, the method works in *very general* situations (including curved faces), and has actually been conceived in a three-dimensional framework.



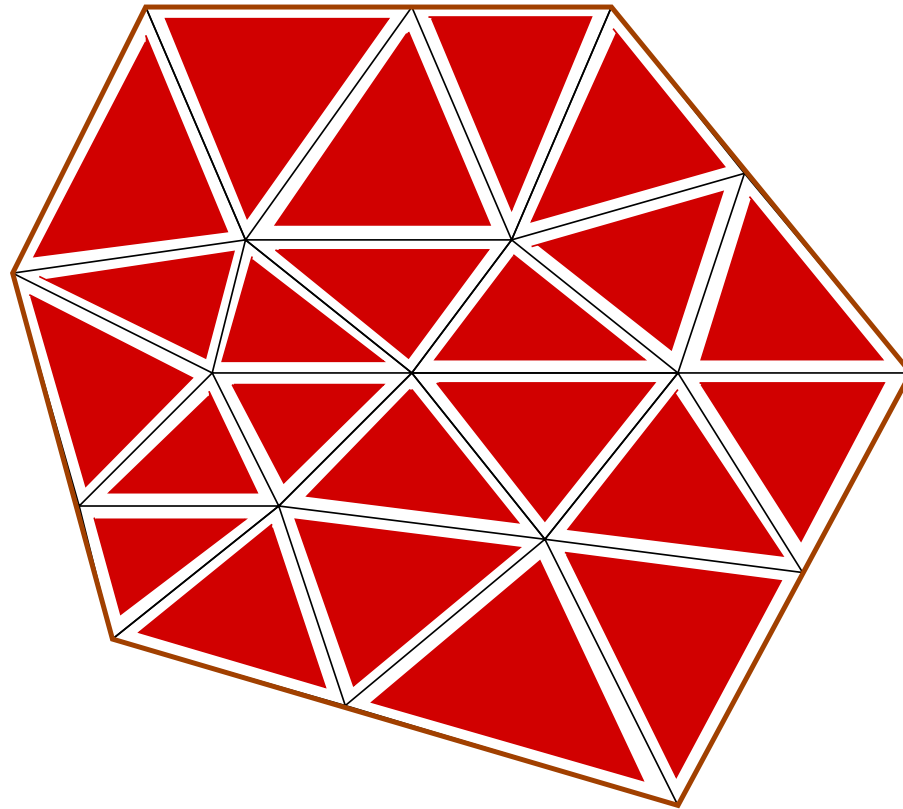
AN ACADEMIC BUT MORE REALISTIC DOMAIN



THE DECOMPOSITION OF Ω

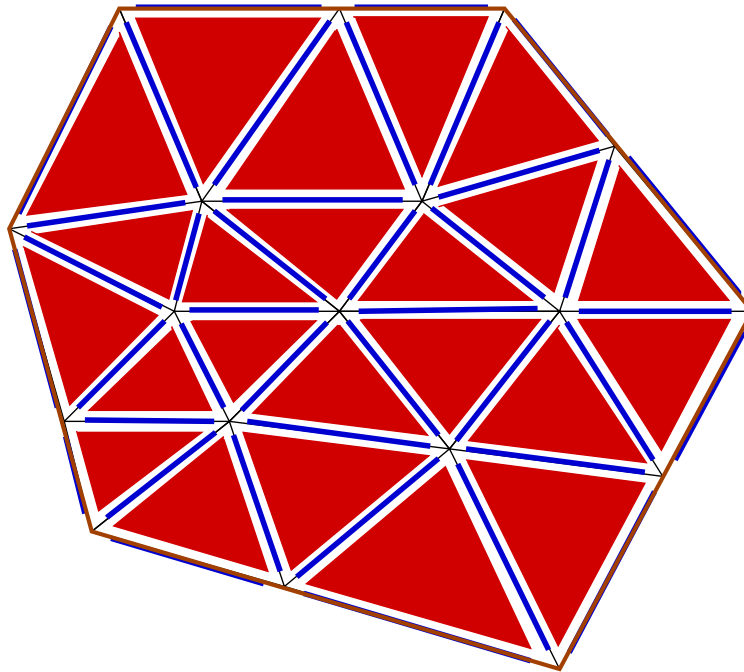


DISCRETE PRESSURES



Space Q : pressure is taken as constant in each element

DISCRETE FLUXES



Space \mathbf{X} : $\mathbf{F} = \mathbf{u} \cdot \mathbf{n}$ is taken constant on each edge (face)

Note that we *do not* discretize \mathbf{u} inside each element, but only its *normal component* on the edges (faces). The orientation of the normal is chosen once and for all.

THE DISCRETE DIVERGENCE OPERATOR

On each element P , from the knowledge of $\mathbf{F} = \mathbf{u} \cdot \mathbf{n}$ on the boundary we can deduce "the average of $\operatorname{div} \mathbf{u}$ inside" by Gauss theorem

$$\int_P \operatorname{div} \mathbf{u} \, dV = \int_{\partial P} \mathbf{u} \cdot \mathbf{n}_{ext} \, dS.$$

This suggests to introduce a *discrete divergence* operator \mathcal{DIV}^h which associates, to every $\mathbf{G} \in \mathbf{X}$, a piecewise constant $\mathcal{DIV}^h \mathbf{G} \in Q'$ defined in each P as

$$\mathcal{DIV}^h \mathbf{G}|_P := \frac{1}{|P|} \int_{\partial P} \mathbf{G}_{ext} \, ds.$$

with obvious meaning for \mathbf{G}_{ext} (equal to plus or minus \mathbf{G} , according to the orientation that was chosen for the normal).

DISCRETIZATION OF THE CONSERVATION EQUATION

In order to discretize the equation

$$\operatorname{div} \mathbf{u} = b,$$

assuming that b is piecewise constant, we can simply use our *discrete divergence* operator $\mathcal{DIV}^h : \mathbf{X} \rightarrow \mathbf{Q}'$. The conservation equation is then discretized as

$$\mathcal{DIV}^h \mathbf{F} = b,$$

or, in variational form, as

$$\mathbf{Q}' \langle \mathcal{DIV}^h \mathbf{F}, q \rangle_{\mathbf{Q}} = \int_{\Omega} b q \, dV \quad \forall q \in \mathbf{Q}.$$

This will be *the same* for Mimetic Finite Differences, Finite Volumes, and Mixed Finite Elements.

DISCRETIZATION OF THE CONSTITUTIVE EQUATION

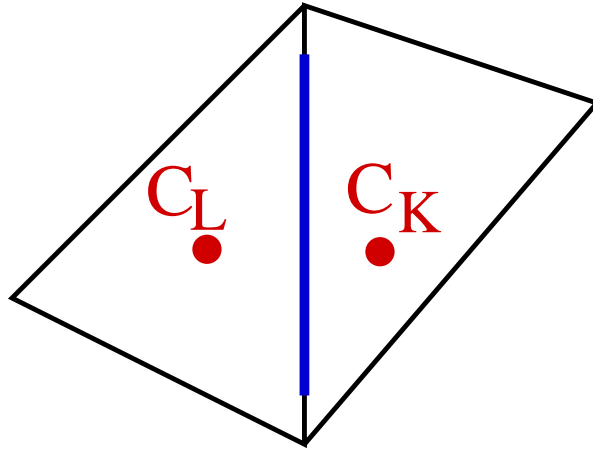
The problem is now to discretize the equation

$$\mathbf{u} = -\mathbb{K}\nabla p$$

{To be precise, we should actually define a *discrete gradient* operator $(-\mathcal{DIV}^h)^T : \mathbf{Q} \rightarrow \mathbf{X}'$ and then a (*-Hodge) operator $\mathbb{K} : \mathbf{X}' \rightarrow \mathbf{X}$.}

The discretization of the Constitutive Equation is done with *three different approaches* in Finite Volumes, in Mixed Finite Elements, and in Mimetic Finite Differences. As we shall see, Mimetic Finite Differences are somewhat *in between* Finite Volumes and Mixed Finite Elements.

FV DISCRETIZATION OF THE CONSTITUTIVE EQUATION



The points C_K and C_L are the *circumcenters* of the triangles.

Assuming for simplicity that $\mathbb{K} = \kappa \mathbb{I}$, the normal flux on the edge e_{KL} is defined, in (one of the most classical formulations of) Finite Volumes, as

$$F_{KL} := \kappa \frac{p_K - p_L}{|C_K - C_L|}$$

MFE DISCRETIZATION OF THE CONSTITUTIVE EQUATION

For Mixed Finite Elements, the passage from pressures to fluxes is made through a **suitable reconstruction** of the fluxes *inside each element*. More precisely in every triangle (tetrahedron) we consider the (**Raviart-Thomas**) space

$$RT := \{\mathbf{v} \mid v_i = a_i + bx_i \quad (i = 1, 2, (3))\}.$$

Then for each $\mathbf{G} \in \mathbf{X}$ we reconstruct $\mathcal{R}(\mathbf{G})$ in Ω such that:

- 1) **in every triangle**, $\mathcal{R}(\mathbf{G})$ belongs to RT ,
- 2) the normal components of $\mathcal{R}(\mathbf{G})$ on **edges (faces)** coincide with \mathbf{G} .
As \mathbf{G} is single valued, this will imply the continuity of the normal components of $\mathcal{R}(\mathbf{G})$.
- 3) $\text{div}\mathcal{R}(\mathbf{G})$ is a piecewise constant (hence **equal to** $\mathcal{DIV}^h \mathbf{G}$).

MFE DISCRETIZATION OF THE CONSTITUTIVE EQUATION

Once you have a reconstruction of the fluxes, you can introduce **the inner product** in \mathbf{X} , depending on \mathbb{K} , as

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{X}} := \int_{\Omega} \mathbb{K}^{-1} \mathcal{R}(\mathbf{F}) \cdot \mathcal{R}(\mathbf{G}) dV.$$

Note that $\mathcal{R}(\mathbf{F}) \sim \mathbf{u}$.

We then discretize the constitutive equation ($\sim \mathbb{K}^{-1} \mathcal{R}(\mathbf{F}) = -\nabla p$) using the variational formulation

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{X}} = \int_{\Omega} p \operatorname{div} \mathcal{R}(\mathbf{G}) dV \quad \forall \mathbf{G} \in \mathbf{X}.$$

FINAL MFE FORMULATION

Summarizing, the Mixed Finite Element formulation **can** be seen, in the present context, as: *find* $p \in Q$ and $\mathbf{F} \in \mathbf{X}$ such that:

$$\int_{\Omega} \mathbb{K}^{-1} \mathcal{R}(\mathbf{F}) \cdot \mathcal{R}(\mathbf{G}) dV = \int_{\Omega} p \operatorname{div} \mathcal{R}(\mathbf{G}) dV \quad \forall \mathbf{G} \in \mathbf{X}$$

and

$$\int_{\Omega} \operatorname{div} \mathcal{R}(\mathbf{F}) q dV = \int_{\Omega} b q dV \quad \forall q \in Q.$$

Note that, in the usual presentation of MFE, the **velocity unknown** is considered to be, directly, $\mathbf{u} \sim \mathcal{R}(\mathbf{F})$ (while its normal components \mathbf{F} on the edges are just *the degrees of freedom*). In MFD, as we shall see, we think instead that *the unknown is* \mathbf{F} , and the reconstruction $\mathbf{u} \sim \mathcal{R}(\mathbf{F})$ (if any) is used only to define the scalar product in \mathbf{X} .

MFD DISCRETIZATION OF THE CONSTITUTIVE EQUATION

We start by mimicking what is done in MFE. For this, we need a suitable *inner product* in \mathbf{X} (depending on \mathbb{K}), that allow us to write

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{X}} = \int_{\Omega} \mathbf{p} \operatorname{DIV}^h \mathbf{G} \, dV \quad \forall \mathbf{G} \in \mathbf{X}.$$

One possibility is again *to reconstruct*. For each element P , from \mathbf{G} on ∂P you reconstruct (somehow) $\mathcal{R}_P(\mathbf{G})$ in P . Then you set

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{X}} := \sum_P \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{F}) \cdot \mathcal{R}_P(\mathbf{G}) \, dV.$$

MFD IN VERY SIMPLE GEOMETRIES

If our elements are triangles (or tetrahedra), then the easiest way to define the inner product in \mathbf{X} is surely to use the Raviart-Thomas (RT) reconstruction as in MFE (although other simple choices are available). In that case, the MFD formulation will *coincide* with MFE: the method will just be *written* differently:

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{X}} \left(= \sum_P \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{F}) \cdot \mathcal{R}_P(\mathbf{G}) dV \right) = \int_{\Omega} p \mathcal{D}IV^h \mathbf{G} dV$$

for all $\mathbf{G} \in \mathbf{X}$, and

$$\int_{\Omega} \mathcal{D}IV^h \mathbf{F} q dV = \int_{\Omega} p q dV$$

for all $q \in Q$.

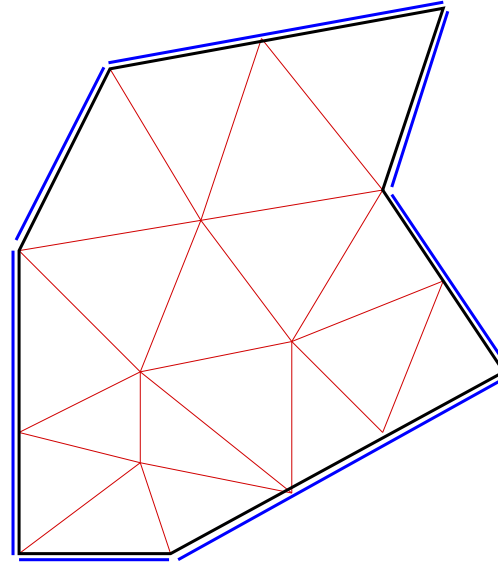
WHAT TO DO FOR COMPLEX GEOMETRIES

In more complex geometries, simple spaces (as *RT*) to be used for the reconstruction are *not available*. Hence to build a suitable reconstruction operator becomes *cumbersome*.

A very good idea to deal with the problem was proposed by Y. Kuznetsov-S. Repin (2004) and generalized by S. Christiansen (2006). It amounts to construct a *subgrid* made of triangles or tetrahedra and reconstruct the fluxes according to the following rules

- In each triangle/tetrahedron the flux is a Raviart-Thomas field.
- The divergence of the flux is *constant on the whole element*.

RECONSTRUCTION USING A SUBGRID



For each $\mathbf{G} \in \mathbf{X}$ and for each element P , we use the subgrid to construct a MFE (Raviart-Thomas) approximate solution $(\boldsymbol{\tau}_h, \phi_h)$ of the Neumann problem

$$-\operatorname{div} \mathbb{K} \nabla \phi = \operatorname{DIV}^h \mathbf{G}_P \quad \text{in } P \quad - \mathbb{K} \nabla \phi \cdot \mathbf{n}_{ext} = (\mathbf{G}_P)_{ext} \quad \text{on } \partial P$$

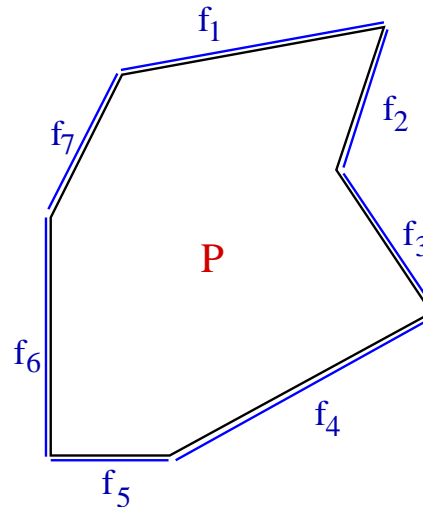
and then we set $\mathcal{R}_P(\mathbf{G}_P) := \boldsymbol{\tau}_h$.

THE BASIC IDEA

We saw that there are various possibilities to build a reconstruction operator, and then to define the scalar product in \mathbf{X} accordingly.

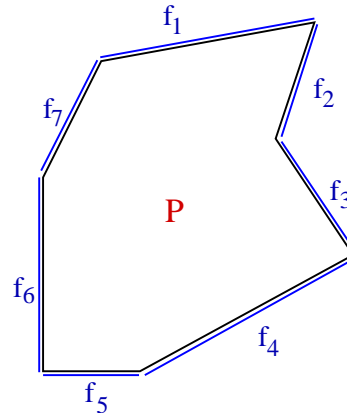
However here **the name of the game** is to **guess** how a "scalar product based on reconstruction" should be, and then **invent** a scalar product **without** actually building a reconstruction operator (and **to get away with that**).

Our scalar product in \mathbf{X} will be defined as the sum of scalar products $[\cdot, \cdot]_{\mathbf{X}_P}$ on individual elements P . To fix ideas, we assume that we are in 2 dimensions, that P has 7 edges, and that \mathbb{K}_P is constant in P .



Assume that you represent the elements of \mathbf{X}_P in the canonical basis $\mathbf{E}^{(1)}, \mathbf{E}^{(2)}, \dots, \mathbf{E}^{(7)}$ by prescribing $\mathbf{E}^{(i)}|_{f_j} = \delta_{i,j}$. Then every element \mathbf{G} in $\mathbf{X}|_P$ will be represented as an element of \mathbb{R}^7 with $\mathbf{G} = \sum_{i=1}^7 \mathbf{G}_i \mathbf{E}^{(i)}$.

ASSOCIATED MATRIX



Every possible reconstruction \mathcal{R}_P will produce a scalar product

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{x}_P} = \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{F}) \cdot \mathcal{R}_P(\mathbf{G}) dV$$

which, in turn, will be representable as a 7×7 matrix \mathbb{M}_P , namely

$$[\mathbf{F}, \mathbf{G}]_{\mathbf{x}_P} = \sum_{i,j} \mathbb{M}_{P i,j} \mathbf{F}_i \mathbf{G}_j, \text{ with}$$

$$\mathbb{M}_{P i,j} := \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{E}^{(i)}) \cdot \mathcal{R}_P(\mathbf{E}^{(j)}) dV.$$

\mathcal{P}_0 -COMPATIBLE RECONSTRUCTIONS

We shall now restrict our attention to *reasonable reconstructions* (that we shall call *\mathcal{P}_0 -compatible reconstructions*). These are linear mappings \mathcal{R}_P defined on \mathbf{X}_P and having the following properties:

- For every $\mathbf{G} \in \mathbf{X}_P$, we have that $\mathcal{R}_P(\mathbf{G}) \in L^2(P)$.
- For every $\mathbf{G} \in \mathbf{X}_P$, we have that $\operatorname{div} \mathcal{R}_P(\mathbf{G})$ is constant in P .
- For every $\mathbf{G} \in \mathbf{X}_P$ and for every face f_i of ∂P , we have $\mathcal{R}_P(\mathbf{G})|_{f_i} \cdot \mathbf{n}|_{f_i} = \mathbf{G}|_{f_i}$ (hence $\operatorname{div} \mathcal{R}_P(\mathbf{G}) = \mathcal{DI}\mathcal{V}^h \mathbf{G}$).
- For every constant vector \mathbf{c} , setting for all faces $\mathbf{G}|_{f_i}^{\mathbf{c}} := \mathbf{c} \cdot \mathbf{n}|_{f_i}$ we have that $\mathcal{R}_P(\mathbf{G}^{\mathbf{c}}) \equiv \mathbf{c}$. Note the difference between $\mathbf{c} \in \mathbb{R}^2$ and $\mathbf{G}^{\mathbf{c}} \in \mathbb{R}^7$!!!

SCALAR PRODUCTS ASSOCIATED WITH \mathcal{P}_0 -COMPATIBLE RECONSTRUCTIONS

We **claim** now that: **if** \mathbf{c} is a constant vector and \mathbf{G}^c has been constructed as before, and **if** \mathcal{R}_P is a \mathcal{P}_0 -compatible reconstruction, **then** for every $\mathbf{G} \in \mathbf{X}_P$ the result of

$$\int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{G}^c) \cdot \mathcal{R}_P(\mathbf{G}) dV$$

depends on P , \mathbb{K} , \mathbf{c} and \mathbf{G} , but **not** on the choice of the reconstruction (among all possible \mathcal{P}_0 -compatible reconstructions).

Indeed...

Set $q^1(\mathbf{x}) := (\mathbb{K}_P^{-1} \mathbf{c}) \cdot (\mathbf{x} - \mathbf{x}_B)$ (where \mathbf{x}_B is the barycenter of P).

Then we have

$$\begin{aligned}
 \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{G}^c) \cdot \mathcal{R}_P(\mathbf{G}) dV &= \\
 \int_P \mathbb{K}_P^{-1} \mathbf{c} \cdot \mathcal{R}_P(\mathbf{G}) dV &= \int_P \nabla q^1 \cdot \mathcal{R}_P(\mathbf{G}) dV = \\
 - \int_P \operatorname{div} \mathcal{R}_P(\mathbf{G}) q^1 dV + \int_{\partial P} q^1 \mathcal{R}_P(\mathbf{G}) \cdot \mathbf{n}_{ext} dS &= \\
 - \int_P \mathcal{D}IV^h \mathbf{G} q^1 dV + \int_{\partial P} q^1 \mathbf{G}_{ext} dS & \\
 &= 0 + \int_{\partial P} (\mathbb{K}_P^{-1} \mathbf{c}) \cdot (\mathbf{x} - \mathbf{x}_B) \mathbf{G}_{ext} dS.
 \end{aligned}$$

Let us summarize the previous result. We found that if the scalar product in \mathbf{X}_P is obtained through a \mathcal{P}_0 -compatible reconstruction, and if \mathbf{G}^c is associated to a *constant vector*, \mathbf{c} , then

$$[\mathbf{G}^c, \mathbf{G}]_{\mathbf{X}_P} = \int_{\partial P} (\mathbb{K}_P^{-1} \mathbf{c}) \cdot (\mathbf{x} - \mathbf{x}_B) \mathbf{G}_{ext} dS.$$

It is also simple to check that taking two constant vectors in the canonical basis of \mathbb{R}^2 , $\mathbf{e}^1 = (1, 0)$ and $\mathbf{e}^2 = (0, 1)$, then we *must* have

$$[\mathbf{G}^{e^i}, \mathbf{G}^{e^j}]_{\mathbf{X}_P} = \int_P \mathbb{K}_P^{-1} \mathcal{R}_P(\mathbf{G}^{e^i}) \cdot \mathcal{R}_P(\mathbf{G}^{e^j}) dV = \int_P \mathbb{K}_P^{-1} \mathbf{e}^i \cdot \mathbf{e}^j dV = (\mathbb{K}_P^{-1})_{i,j} |P|.$$

It seems now natural to *change the basis in \mathbf{X}_P* We take

$$\tilde{\mathbf{E}}^1 := \mathbf{G}^{\mathbf{e}^1}, \quad \tilde{\mathbf{E}}^2 := \mathbf{G}^{\mathbf{e}^2},$$

and then we complete the basis with vectors in \mathbb{R}^7

$$\tilde{\mathbf{E}}^3, \quad \tilde{\mathbf{E}}^4, \dots, \quad \tilde{\mathbf{E}}^7$$

such that

$$[\mathbf{G}^{\mathbf{e}^i}, \tilde{\mathbf{E}}^j]_{\mathbf{X}_P} = \int_{\partial P} (\mathbb{K}_P^{-1} \mathbf{e}^i) \cdot (\mathbf{x} - \mathbf{x}_B) \tilde{\mathbf{E}}_{ext}^j \, dS = 0 \quad (i = 1, 2 \quad j = 3, \dots, 7)$$

Note that all this *does not* depend on the choice of the reconstruction.

In the new basis $\widetilde{\mathbf{E}}^1, \dots, \widetilde{\mathbf{E}}^7$ the matrix associated to **any** scalar product obtained with a \mathcal{P}_0 -compatible reconstruction will then have the form

$\mathbf{K}^{-1} \mathbf{P} $	0
0	?

and the 5×5 diagonal block "?" will depend on the reconstruction.

Theorem. There exists an $\alpha_0 > 0$ such that: for every symmetric and positive definite 5×5 matrix \mathbf{S} with smallest eigenvalue $\geq \alpha_0$ there exists a \mathcal{P}_0 -compatible reconstruction whose associated scalar product corresponds, in the basis $\widetilde{\mathbf{E}}^1, \dots, \widetilde{\mathbf{E}}^7$, to the matrix

$\mathbf{K}^{-1} \mathbf{P} $	$\mathbf{0}$
$\mathbf{0}$	\mathbf{S}

In other words: I know that the matrix comes from **a** reconstruction. I don't care to know "which one".

In all our experiments we took the matrix as

$\mathbb{K}^{-1} P $	0
0	αI

with $\alpha = |P|\text{trace}(\mathbb{K}^{-1})$, and we got very good results.

ERROR ESTIMATES (B.-Lipnikov-Shashkov-Simoncini)

Assume that:

- Ω has a Lipschitz continuous boundary.
- Every element is *uniformly strictly* starshaped.
- The number of faces per element is uniformly bounded.
- Every face is *uniformly strictly* starshaped.
- The number α of the previous slide verifies $\alpha \simeq |P|$.

Then

$$\|p - p_h\|_0 + \|\mathbf{F}^I - \mathbf{F}_h\|_{\mathbf{X}} \leq C h.$$

If moreover Ω is convex and $\alpha \geq \alpha_0$, then

$$\|p^I - p_h\|_0 \leq C h^2,$$

where p^I is the (element by element) mean value of p .

CURVED FACES IN 3 DIMENSIONS

To deal with *curved faces* we must use, *inside each domain*, *three unknowns per face* (the three components of a constant vector).

The face e is called *moderately curved* if, *at every point* of e , we have

$$|\tilde{\mathbf{n}} - \mathbf{n}| \leq \frac{\sigma_*}{L} |e|^{1/2} \quad \text{where} \quad \tilde{\mathbf{n}} := \frac{\int_e \mathbf{n} dS}{|e|}.$$

Here σ_* is a (chosen) positive number and L a characteristic length. A face that is not moderately curved will be called *strongly curved*.

Clearly, the bigger we choose σ_* , the bigger will be the number of *moderately curved* faces, and the smaller that of *strongly curved* ones.

In assembling the global matrix, we require *only the continuity of the averaged normal component* on *moderately curved faces*, but we require the continuity of *all three components* on the *strongly curved faces*.

CURVED FACES

The total number of flux unknowns will then equal the number of **moderately curved** faces plus **three times** the number of **strongly curved** faces.

Hence, **strongly curved** faces are **more expensive**. From the point of view of **cost** we have then the **convenience to take σ_* very big**, so that the number of strongly curved faces will be small.

As usual, however, **you pay somewhere else**: actually the error estimates go like

$$\|p - p_h\|_0 + \|\mathbf{F}^I - \mathbf{F}_h\|_{\mathbf{X}} \leq C(1 + \sigma_*)h.$$

Finally, we point out that the **same trick** (actually, **better**) for computing the scalar products holds for curved faces.

NUMERICAL EXPERIMENTS

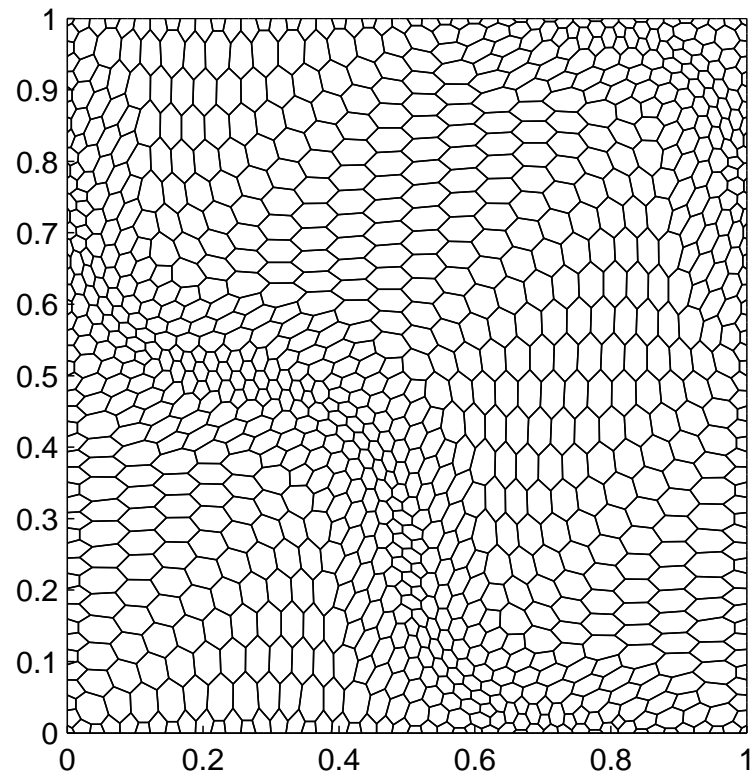
Example 1. Let us consider the Dirichlet boundary value problem in the unit square $[0, 1]^2$ with the exact solution

$$p(x, y) = x^3 y^2 + x \sin(2\pi xy) \sin(2\pi y)$$

and the full diffusion tensor

$$\mathbb{K} = \begin{pmatrix} (x+1)^2 + y^2 & -xy \\ -xy & (x+1)^2 \end{pmatrix}.$$

Polygonal mesh

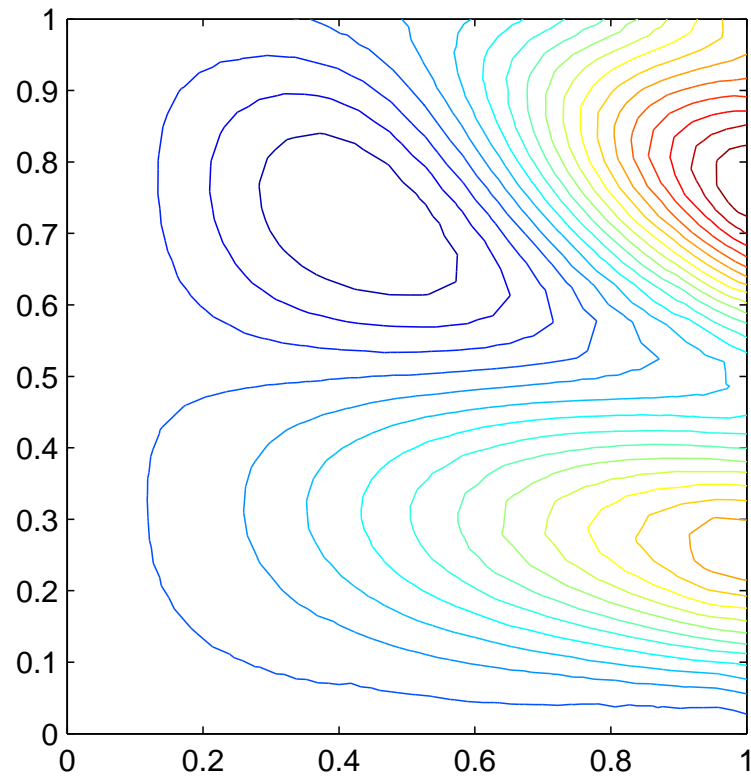


The table below shows the errors for different refinements.

Table 1: Convergence analysis on polygonal meshes.

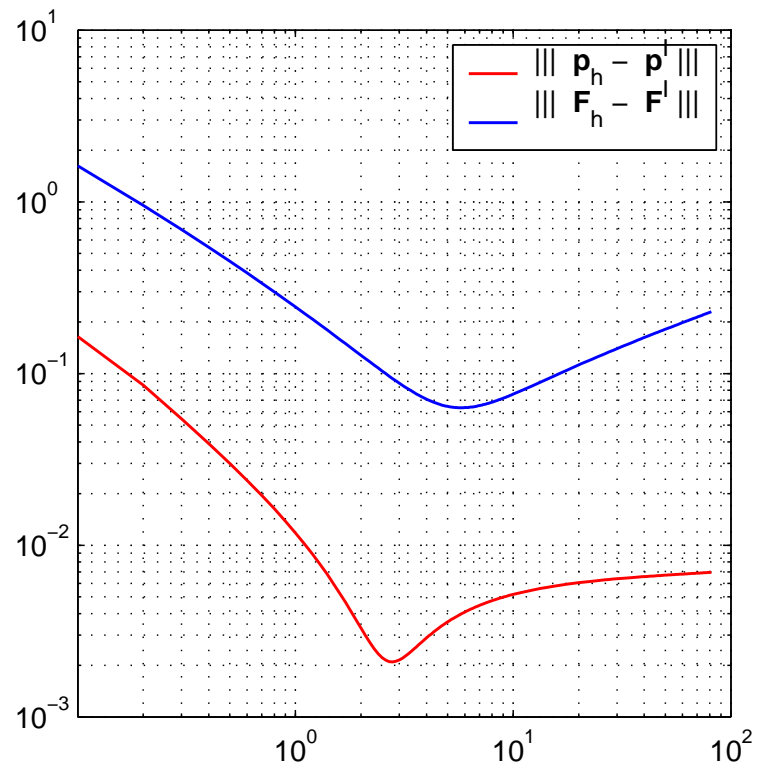
$1/h$	$\ p^I - p_h\ _0$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\mathbf{X}}$	$\ p^I - p_h\ _{\infty}$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\infty}$
16	5.17e-2	7.38e-1	1.61e-1	5.25e-0
32	1.18e-2	2.44e-1	4.54e-2	2.80e-0
64	2.76e-3	8.45e-2	1.28e-2	1.46e-0
128	6.65e-4	2.89e-2	3.06e-3	7.79e-1
rate	2.09	1.56	1.90	0.92

Isolines of the solution



Errors as a function of α

The spectral properties of the matrix \mathbb{M}_P defining the scalar product in P depend on a balance between the extreme eigenvalues of \mathbb{K} and α . The figure in the next slide shows errors (vertical axis) as functions of α^{-1} (horizontal axis) for the case $1/h = 32$. There is a quite big interval $\alpha^{-1} \in [2, 80]$ where the errors vary only 3 times. What is remarkable here is that for all values of α we observed second order convergence rate for p_h in L^2 and 1.5 convergence rate for \mathbf{F}_h in L^2 .

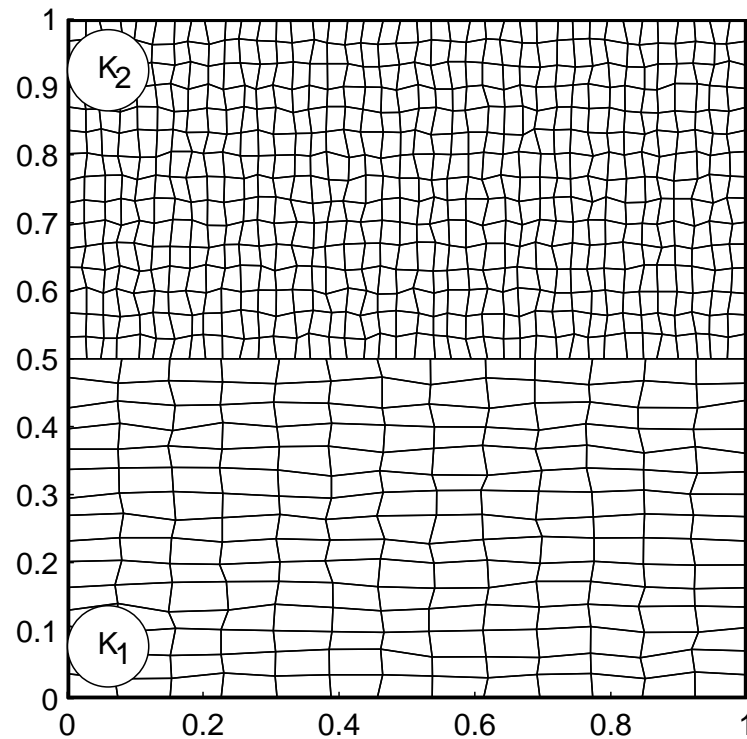


Example 2. Let us consider a problem in the unit square $[0, 1]^2$ with *mixed* boundary conditions. On the **bottom** ($y = 0$) and **top** ($y = 1$) boundaries, we impose the **Dirichlet** boundary condition. Homogeneous **Neumann** boundary conditions are set on the **other** (**vertical**) boundaries. Let the tensor \mathbb{K} be scalar and equal to $K_1\mathbb{I}$ in the region defined by $y < 0.5$ and $K_2\mathbb{I}$ in the rest of the domain. The source term is chosen in such a way that the **exact solution** is given by

$$p(x, y) = \begin{cases} \frac{7}{16} - \frac{K_2}{12K_1} + \frac{2K_2}{3K_1} y^3, & y < 0.5, \\ y - y^4, & y \geq 0.5 \end{cases}$$

and hence, in particular, **depends only on y** .

The meshes we used come from **non-matching mesh methods** but we treat them as **conformal polygonal meshes**. The mesh below is obtained as a **random perturbation** of an originally "uniform, non matching" mesh.

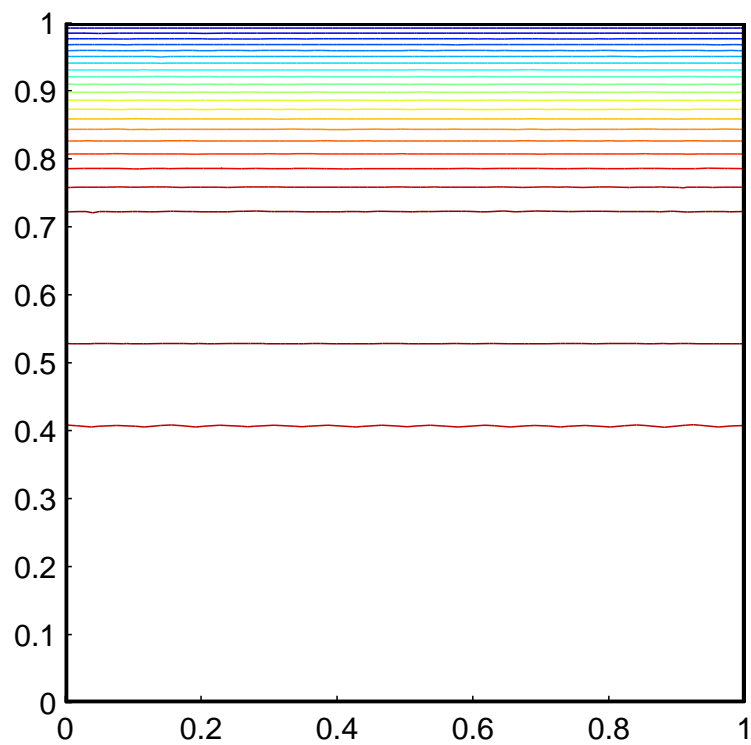


In the numerical experiments we used $K_1 = 10$, $K_2 = 1$. We observe **superconvergence** of the **scalar** variable in both norms. The **lack of flux superconvergence** is typical for random meshes and is observed in other similar discretization schemes on simplicial meshes.

Table 2: Convergence analysis on non-matching meshes.

#cells	$\ p^I - p_h\ _0$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\mathbf{x}}$	$\ p^I - p_h\ _{\infty}$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\infty}$
780	1.01e-2	1.12e-1	2.82e-2	7.80e-1
3286	2.36e-3	4.72e-2	6.70e-3	3.51e-1
13482	5.73e-4	2.24e-2	1.78e-3	1.38e-1
54610	1.41e-4	1.09e-2	4.37e-4	7.70e-2
rate	2.01	1.09	1.95	1.11

Isoline plot of the solution



Example 3. Let us consider the **Dirichlet** boundary value problem with the **exact solution**

$$p(x, y, z) = x^3 y^2 z + x \sin(2\pi xy) \sin(2\pi yz) \sin(2\pi z).$$

We take the **diffusion tensor** as:

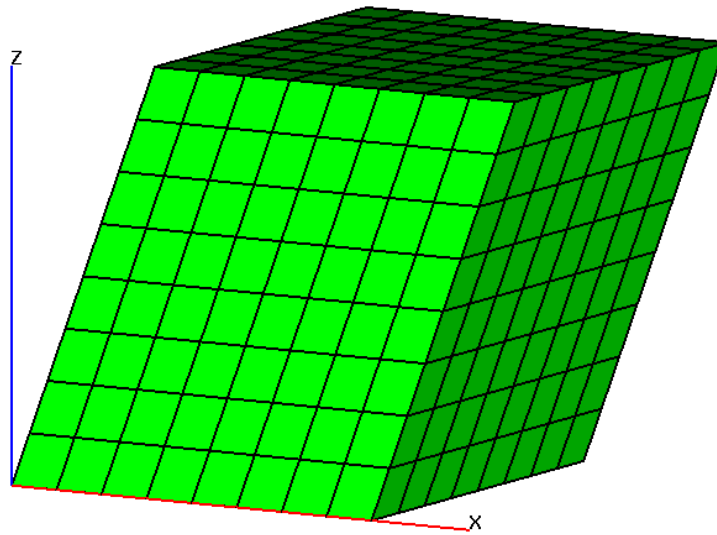
$$\mathbb{K} = \begin{pmatrix} 1 + y^2 + z^2 & -xy & -xz \\ -xy & 1 + x^2 + z^2 & -yz \\ -xz & -yz & 1 + x^2 + y^2 \end{pmatrix}.$$

It is not difficult to check that \mathbb{K} is a positive definite matrix for all values of x , y and z .

We consider a sequence of **uniform cubic meshes** in the unit cube $[0, 1]^3$ and **generate a corresponding sequence of hexahedral meshes** using the following linear transformation:

$$\tilde{x} = x + \varepsilon z, \quad \tilde{y} = y + \varepsilon z, \quad \tilde{z} = z.$$

An example of a modified mesh is shown in the next figure. It corresponds to $\varepsilon = 0.25$ and $h = 1/8$, where h is the size of a cubic element in the original mesh.

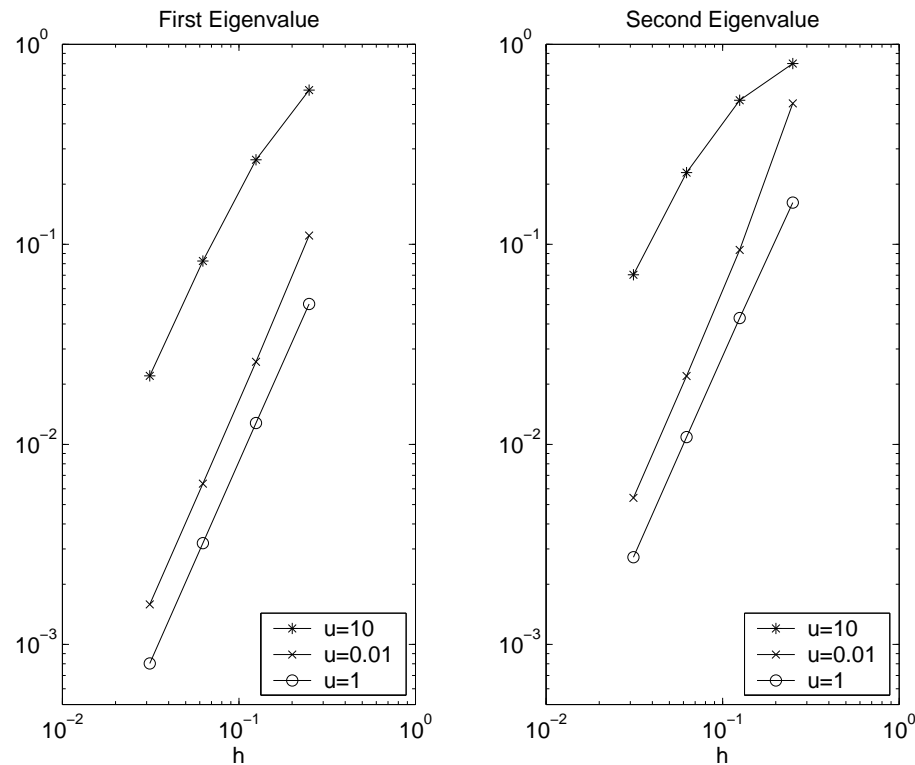


The results presented in the table below show the **superconvergence** of the scalar variable in both norms and the superconvergence of the vector variable in the discrete L_2 -norm.

Table 3: Convergence analysis on polyhedral meshes.

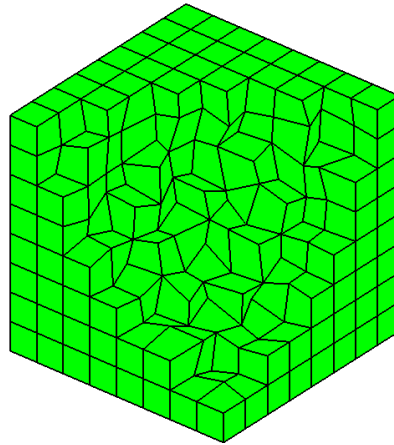
$1/h$	$\ p^I - p_h\ _0$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\mathbf{X}}$	$\ p^I - p_h\ _{\infty}$	$\ \mathbf{F}^I - \mathbf{F}_h\ _{\infty}$
8	3.83e-2	5.35e-1	1.55e-1	6.07e-0
16	1.10e-2	1.43e-1	4.83e-2	2.48e-0
32	2.86e-3	3.58e-2	1.26e-2	1.11e-0
64	7.21e-4	8.94e-3	3.28e-3	5.42e-1
rate	1.91	1.97	1.86	1.16

Approximation of eigenvalues



Approximation of the first two eigenvalues of Laplace operator on the square $]0, 1[{}^2$, on a uniform mesh of squares, for different values of $\alpha = 0.01, 1, 10$. By A. Cangiani and M. Manzini.

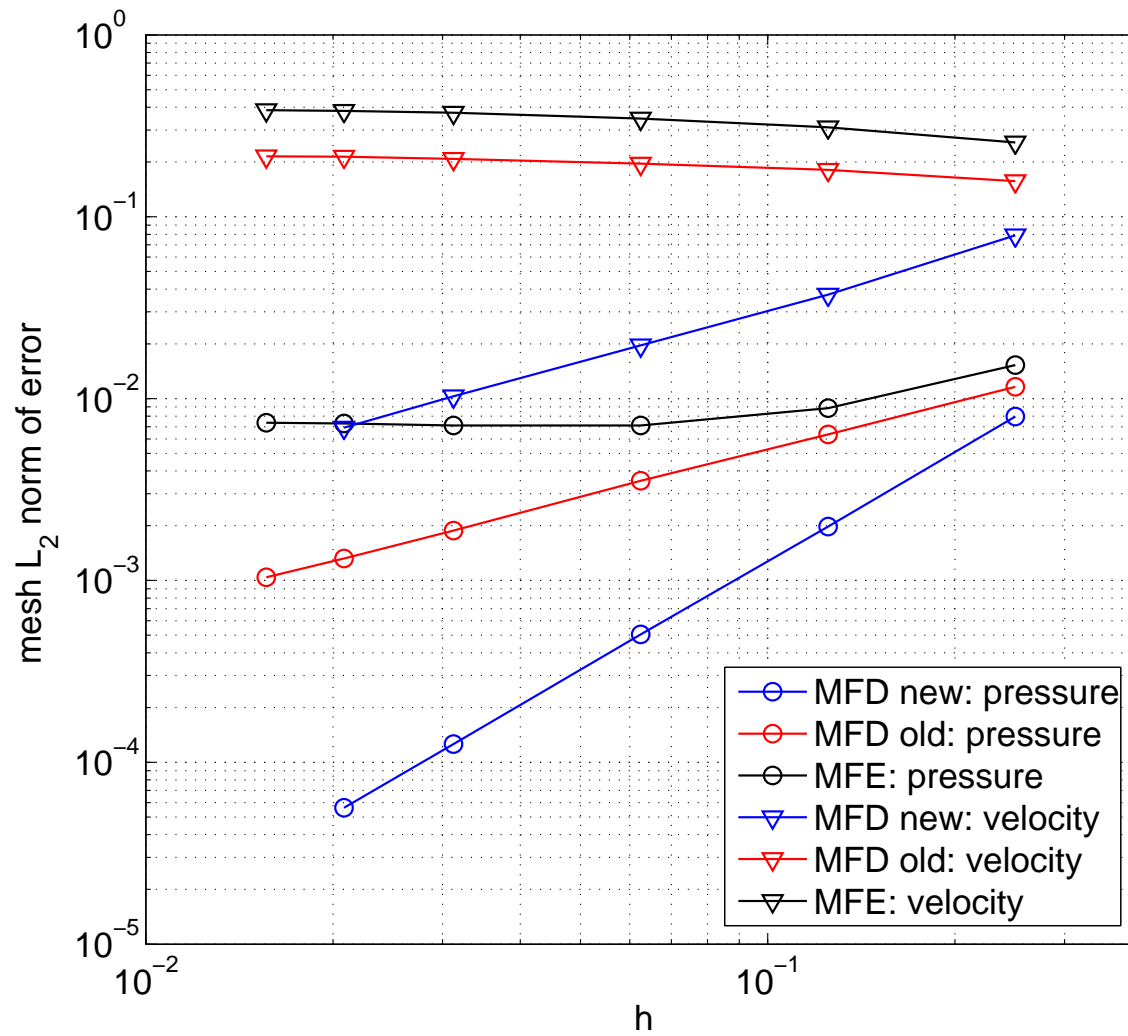
3D-Curved faces

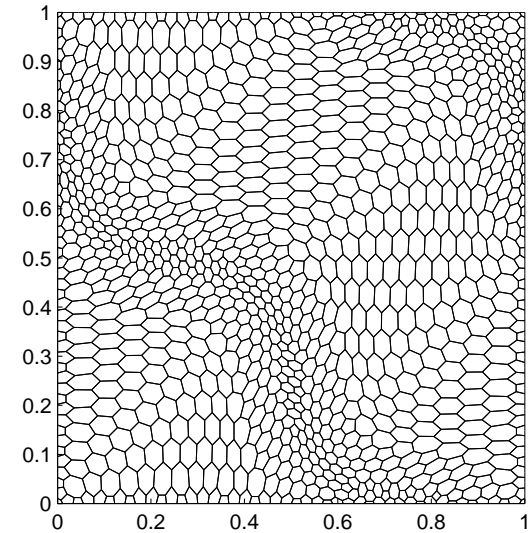
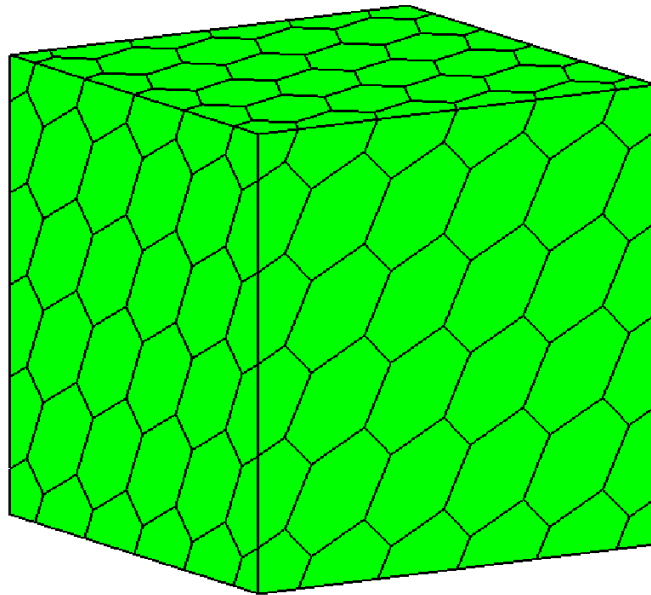


Here we have $\mathbb{K}=\text{Identity}$. The exact solution is

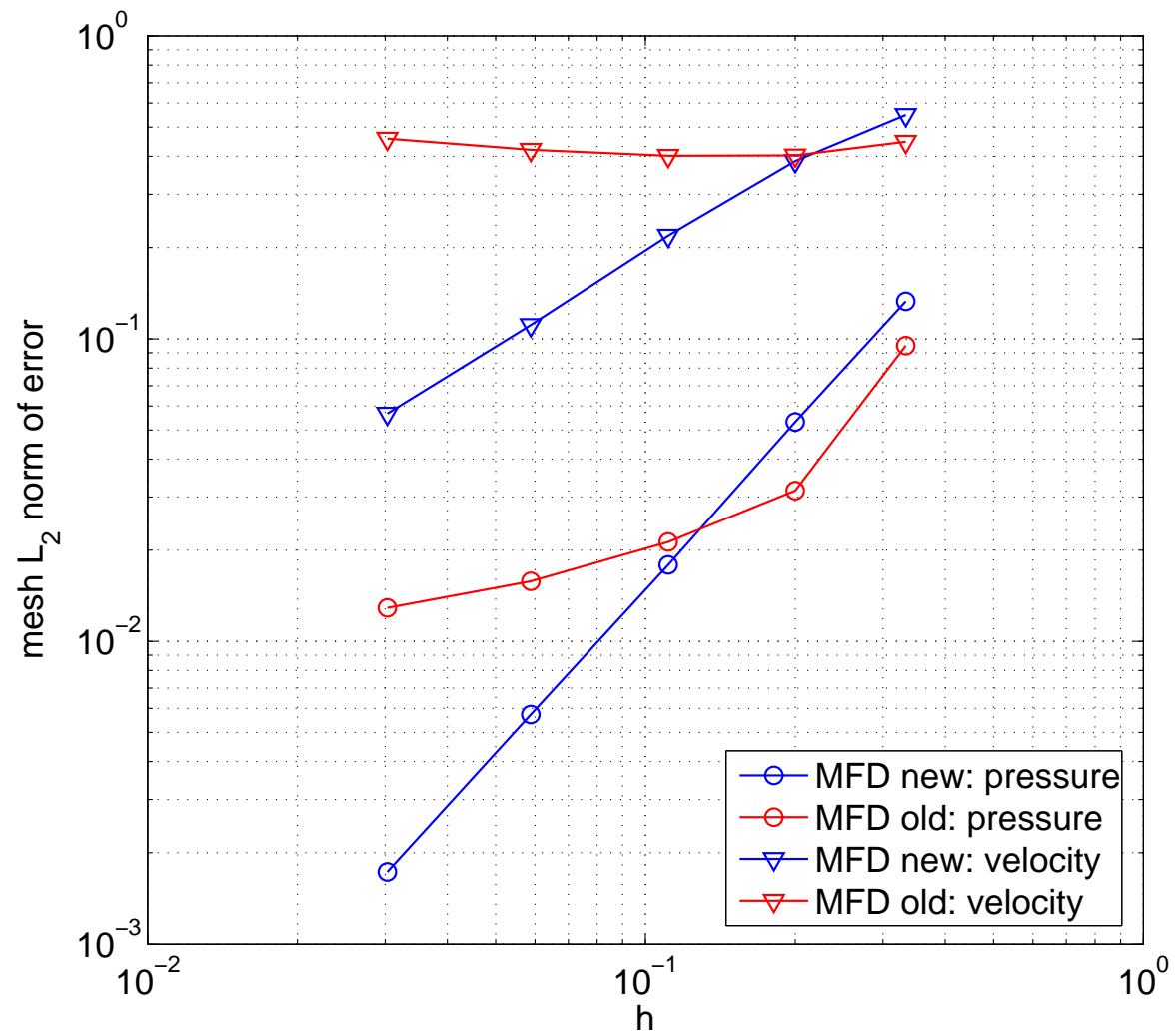
$$p(x, y, z) = x^2 y^3 z + 3x \sin(yz).$$

Now we compare: Mixed Finite Elements, Mimetic Finite Differences treating all faces as moderately curved, and Mimetic Finite Differences with $\sigma^* = 0.2$.





We do the same for a polyhedral mesh (distorted inside as in the 2-d case). Here we have no Mixed Finite Elements available for comparison.



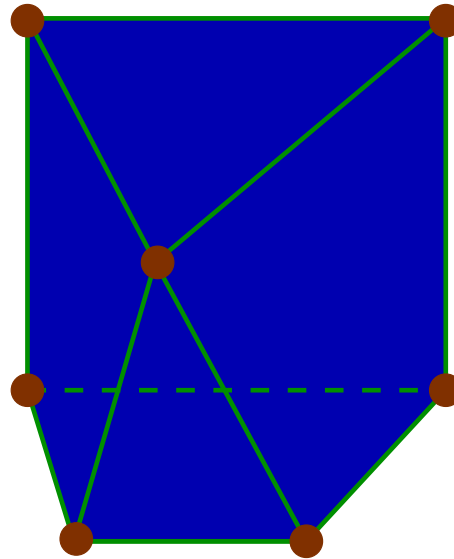
CONCLUSIONS

- Mimetic Finite Differences aim at a better representation of **fundamental physical laws**.
- The present approach allows an **enormous freedom** in the type of decomposition. The treatment of **non-matching grids** and of **curved faces** is also allowed.
- The basic ideas could be generalized to other problems, involving *gradients* and *curls* (in the framework of *cochain approximations of differential forms*).
- We still have to understand how to make the **best use** of the freedom allowed by our construction: indeed S does not need to be a diagonal matrix αI .
- For **curved faces** we need more indications on the optimal choice for the parameter σ_* .

APPENDIX

A ONE-DOMAIN SETTING

We consider a **Polyhedron**, with its **Vertices**, its **Edges**, and its **Faces**



● Vertices

— Edges

▭ Faces

7 Vertices

11 Edges

6 Faces

We consider, to start with, a single polyhedron P with vertices V_1, V_2, \dots, V_7 , edges e_1, e_2, \dots, e_{11} , and faces f_1, f_2, \dots, f_6 . The numbers (7, 11, 6) are taken from the example of the previous figure, but of course they change from one geometry to another.

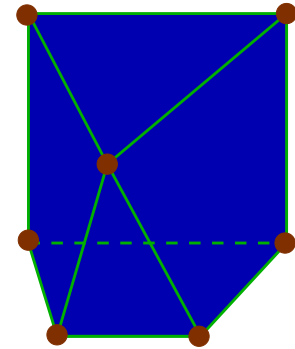
The motivation for working on a single element is that we want to discuss *scalar products*. All the scalar products that are going to be used on the whole structure (made, obviously, of several polyhedra) will be constructed *separately* in each element, and then summed over the elements.

We are going to consider four types of unknowns:

- **node unknowns** (whose values are defined only at the **vertices** and are to be interpreted as the values of a scalar function at each node)
- **edge unknowns** (whose value is defined only on the **edges** and are to be interpreted as the *integral of the tangential component* of a vector valued-function on each edge)
- **face unknowns** (whose values are defined only on the **faces** and are to be interpreted as the *integral of the outward normal component* of a vector-valued function on each face)
- **element unknowns** (whose values are defined only inside the **element** and are to be interpreted as the *integral* of a scalar function over the element)

Accordingly, we denote

- by \mathcal{N} the space of all **node unknowns**
- by \mathcal{E} the space of all **edge unknowns**
- by \mathcal{F} the space of all **face unknowns**
- by \mathcal{P} the space of all **element unknowns**



The dimensions of the above spaces will be, in our case, 7, 11, 6 and 1, respectively.

These spaces can obviously be interpreted as well as discretizations of the spaces of 0 – *forms*, 1 – *forms*, 2 – *forms*, and 3 – *forms*, respectively.

Although we did not say it explicitly, the space \mathcal{E} will require an orientation of the tangent vector to edges, and, if more than one element is considered, the space \mathbf{F} will require an orientation of faces. Here we have only one element, and we can choose the outward normal to each face, once and for all.

THE \mathcal{GRAD}^h operator

It is now natural (according to our *interpretation*) to construct three basic differential operators:

- The \mathcal{GRAD}^h operator, from \mathcal{N} to \mathcal{E} , defined as follows:

for each edge e with vertices V_1 and V_2 , and for each element $u \in \mathcal{N}$, we define $\mathcal{GRAD}^h u$ on e in the direction from V_1 to V_2 as

$$\mathcal{GRAD}^h u|_e = u(V_2) - u(V_1)$$

THE $CURL^h$ operator

- The $CURL^h$ operator, from \mathcal{E} to \mathbf{F} , defined as follows:
for each element $\varphi \in \mathcal{E}$ and for each face f we orient the edges e_1, e_2, \dots, e_n of f in the counterclockwise direction (seen from outside the element), and then we consider the corresponding values $\varphi_1, \varphi_2, \dots, \varphi_n$ of φ with the sign corresponding to the orientation just chosen. Then *the value of $CURL^h \varphi$ on the face f is defined as*

$$CURL^h \varphi|_f = \sum_{i=1}^n \varphi_i$$

THE $DI\mathcal{V}^h$ operator

- The $DI\mathcal{V}^h$ operator, from \mathbf{F} to \mathcal{P} , defined as follows:
let f_1, \dots, f_k be all the faces of our element P , and for each $\sigma \in \mathbf{F}$
let $\sigma_1, \dots, \sigma_k$ be its values on each face. Then $DI\mathcal{V}^h \sigma$ is defined as

$$DI\mathcal{V}^h \sigma = \sum_{i=1}^k \sigma_i$$

Note: in our example, $k = 6$. Of course. If we had more than one element, the same construction would be done for each element, taking care to use the values of σ corresponding to the *outward* direction.

THE ASSOCIATED MATRICES

It is interesting to note that, taking in the spaces \mathcal{N} , \mathcal{E} , \mathcal{F} , \mathcal{P} the obvious canonical basis (after choosing an orientation of the edges in an arbitrary way, but once and for all), then the matrices associated with each one of the operators $GRAD^h$, $CURL^h$, and DIV^h are just made of zeroes, ones, and minus ones.

All this can by now be considered as *classical*. I will just mention the works of Shashkov and co-authors, Kuznetsov, Bochev, Yotov, Wheeler, Christiansen, Buffa, and many others.

THE INTERPOLANTS

Along the same way of thought, for each smooth scalar function u and for each smooth vector valued function $\boldsymbol{\theta}$ we can define the following *interpolants*

- $u_{I,node} \in \mathcal{N}$ defined by

$$u_{I,node}(V) = u(V) \quad \text{for all vertex } V$$

- $\theta_{I,edge} \in \mathcal{E}$ defined by

$$\theta_{I,edge}|_e = \int_e \boldsymbol{\theta} \cdot \mathbf{t} ds \quad \text{for all edge } e$$

where the unit tangent vector \mathbf{t} indicates the orientation of e

THE INTERPOLANTS-2

- $\theta_{I,face} \in \mathbf{F}$ defined by

$$\theta_{I,face}|_f = \int_f \boldsymbol{\theta} \cdot \mathbf{n} dS \quad \text{for all face } f$$

where \mathbf{n} is the unit normal outward vector to f

- $u_{I,poly} \in \mathcal{P}$ defined by

$$u_{I,poly} = \int_P u dP.$$

THE COMMUTATION PROPERTIES

Note that the interpolation operators and the differential operators introduced above have interesting commutation properties. Namely

- $GRAD^h(u_{I,node}) = (\text{grad}u)_{I,edge}$
- $CURL^h(\boldsymbol{\theta}_{I,edge}) = (\text{curl}\boldsymbol{\theta})_{I,face}$
- $DIV^h(\boldsymbol{\theta}_{I,face}) = (\text{div}\boldsymbol{\theta})_{I,poly}$.

THE COMMUTING DIAGRAM

The above properties could be summarized saying that the following diagram

$$\begin{array}{ccccccc}
 C^\infty(P) & \xrightarrow{\text{grad}} & (C^\infty(P))^3 & \xrightarrow{\text{curl}} & (C^\infty(P))^3 & \xrightarrow{\text{div}} & C^\infty(P) \\
 \downarrow & & \downarrow & & \downarrow & & \downarrow \\
 \mathcal{N} & \xrightarrow{\underline{GRAD}^h} & \mathcal{E} & \xrightarrow{\underline{CURL}^h} & \mathbf{F} & \xrightarrow{\underline{DIV}^h} & \mathcal{P}
 \end{array}$$

(where the vertical arrows are the interpolation operators just defined), commutes.

All this *should* be related to the approximation of differential complexes intensively studied, among others, by Arnold, Falk, and Winther. See the recent review in *Acta Numerica*.

TRANSPOSED OPERATORS

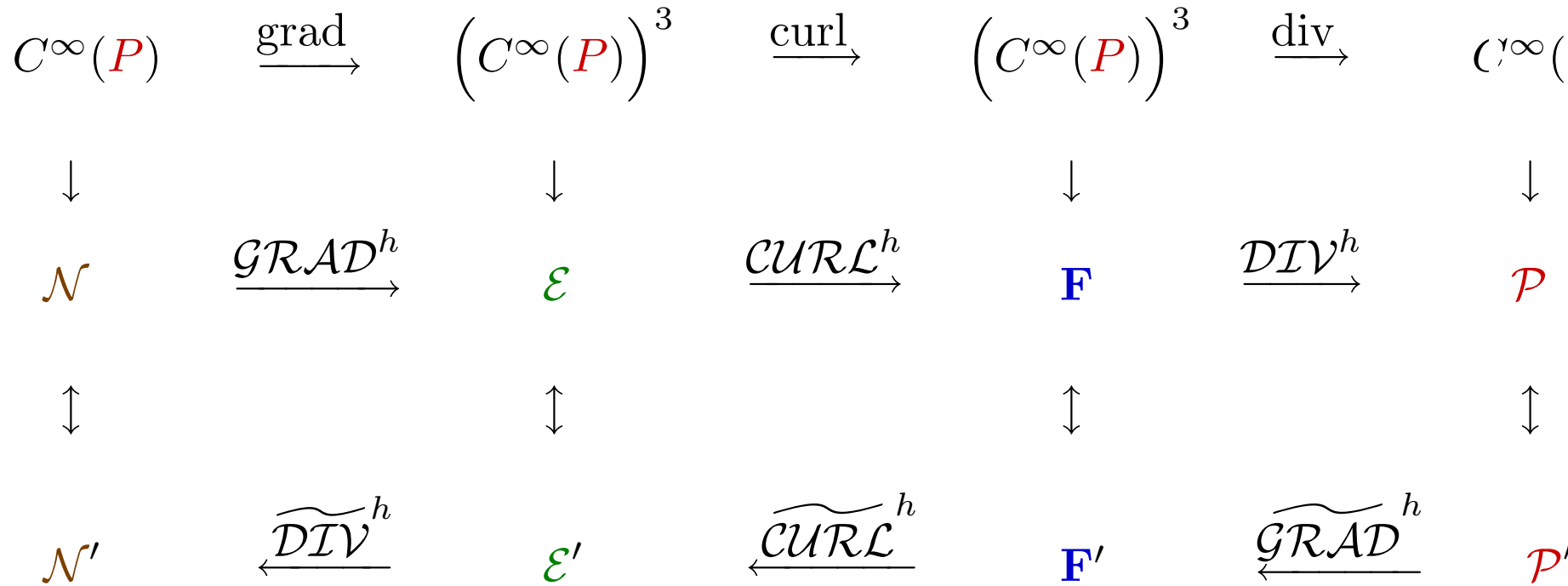
As we are dealing with finite dimensional spaces, we can introduce immediately the dual spaces of the spaces \mathcal{N} , \mathcal{E} , \mathbf{F} , \mathcal{P} , and transpose the operators $GRAD^h$, $CURL^h$, and DIV^h .

It seems rather natural (although one might question it) to denote by $\widetilde{GRAD}^h : \mathcal{P}' \rightarrow \mathbf{F}'$ the transposed of the operator $-DIV^h$, and by $\widetilde{DIV}^h : \mathcal{E}' \rightarrow \mathcal{N}'$ the transposed of the operator $-GRAD^h$. Similarly we shall denote by $\widetilde{CURL}^h : \mathbf{F}' \rightarrow \mathcal{E}'$ the transposed of $CURL^h$.

The dual spaces \mathcal{P}' , \mathbf{F}' , \mathcal{E}' , \mathcal{N}' could then be thought as (*weird*) approximations of 0 – forms, 1 – forms, 2 – forms, and 3 – forms, respectively.

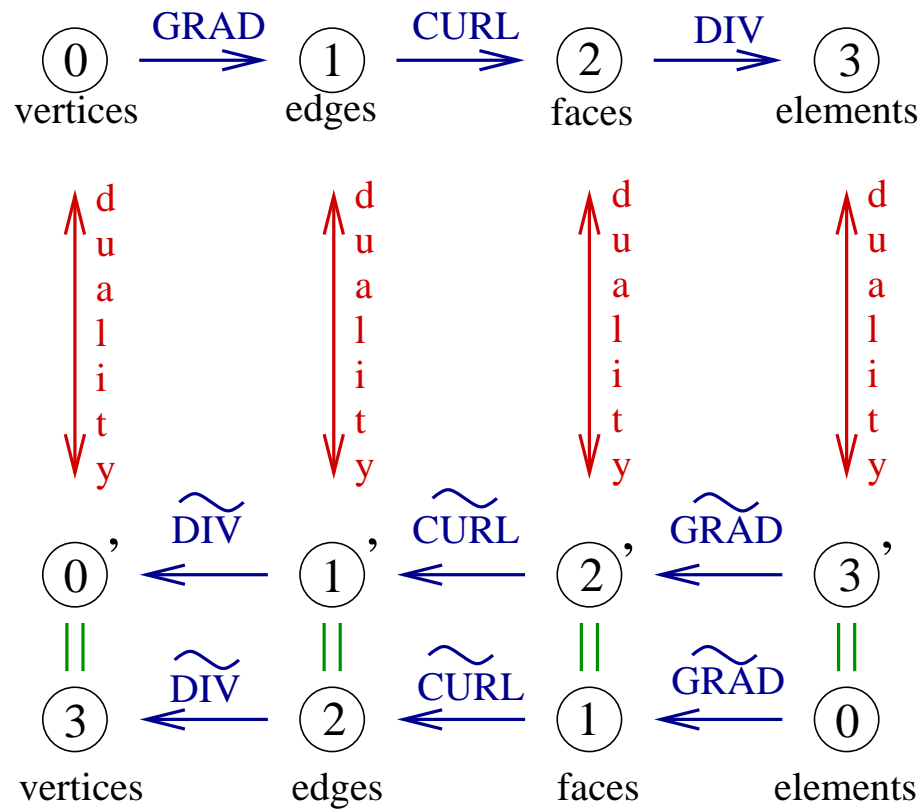
TRANSPOSED OPERATORS

In this way, we have the following diagram



DISCRETE DIFFERENTIAL FORMS

The previous diagram can be thought as made of *discretized differential forms*



INNER PRODUCTS

The previous discretizations of differential forms is quite appealing (and meaningful) on the first line. In order to transform the last two lines in a useful instrument, we need however to introduce *suitable inner products* in the spaces \mathcal{N} , \mathcal{E} , \mathbf{F} , \mathcal{P} .

If these inner products will be *consistent* with the $L^2(\Omega)$ inner product (for scalars) or with the $(L^2(\Omega))^2$ inner product (for vectors), then the adjoint operators will start to make sense in a more familiar way.

In doing this, however, it will be more convenient to insert the *material properties* into our inner products. This will lead us to a discretization of the $*$ – *Hodge* operator.

All this is vague. Let us consider a particular case in order to make it clearer.

SCALAR PRODUCTS in \mathcal{N} (L^2 -TYPE)

We already saw convenient scalar products in the spaces of 3 – *forms* (elements) and of 2 – *forms* (faces). Let us see some other scalar products.

First one might think of an L^2 -inner product. For that it seems reasonable to request a 7×7 matrix. Denoting by $M_{\mathcal{N}}$ such matrix, it seems reasonable to require that

$$\sum_{i=1}^7 \sum_{j=1}^7 (M_{\mathcal{N}})_{i,j} u(V_i) v(V_j) = \int_P u v \, dP$$

whenever u and v are the interpolant of two polynomials of degree ≤ 1 (u and v , respectively). To construct such a matrix $M_{\mathcal{N}}$ we can proceed as before (solving explicitly with a semi-positive definite matrix, and then adding a suitable correction that makes it positive definite).

SCALAR PRODUCTS in \mathcal{N} (H^1 -TYPE)

We proceed to introduce a scalar product in \mathcal{E} for gradients. The dimension of the subspace of gradients is clearly 6 (or, the number of vertices minus one). We denote by \mathcal{E}^G the subspace of \mathcal{E} made of gradients. We would like to construct a symmetric 11×11 matrix $M_{\mathcal{E}}^G$ (SPD on \mathcal{E}^G) such that: for every polynomial v of degree ≤ 1 , and for every $\varphi = \mathcal{GRAD}^h u$ (with $u \in \mathcal{N}$), setting $\psi := (\text{grad} v)_{I, \text{edge}} \equiv \mathcal{GRAD}^h v_{I, \text{node}}$ we have

$$\sum_{i=1}^{11} \sum_{j=1}^{11} (M_{\mathcal{E}}^G)_{i,j} \varphi_i \psi_j = \int_{\partial P} u \frac{\partial v}{\partial \mathbf{n}}.$$

There is a problem with the above formula: Since u is only defined at the nodes, we cannot evaluate the integral over the faces. Hence we need to pick up a numerical integration formula. In most cases we only need such formula to be exact for linear functions.

SCALAR PRODUCTS in \mathcal{N} (H^1 -TYPE)

If the matrix $M_{\mathcal{E}}^G$ satisfies the previous formula one can prove that: there exists a lifting linear mapping $r : \mathcal{N} \rightarrow C^0(\bar{P})$ such that

- i) $(r(w))_{I,node} \equiv w$ for each $w \in \mathcal{N}$,
- ii) the value of $r(u)$ on each face f depends only on the values of u at the nodes belonging to f ,
- iii) for every linear function v we have $r(v_{I,node}) = v$,
- iv) and moreover:

$$\sum_{i=1}^{11} \sum_{j=1}^{11} (M_{\mathcal{E}}^G)_{i,j} (\mathcal{GRAD}^h(u))_i (\mathcal{GRAD}^h(v_{I,node}))_j = \int_{\partial P} r(u) \frac{\partial v}{\partial \mathbf{n}}.$$

The best way to complete the scalar product in \mathcal{E} beyond the gradients is still a rather open question.

MUMBLE GRUMBLE.