

## LA-UR-12-24326

Approved for public release; distribution is unlimited.

Title:	From the Mimetic Finite Difference Method to the Virtual Element Method
Author(s):	Manzini, Gianmarco
Intended for:	NSF Workshop on Barycentric Coordinates in Geometry Processing and Finite/Boundary Element Methods, 2012-07-25/2012-07-27 (New York, New York, United States)



### Disclaimer:

Los Alamos National Laboratory, an affirmative action/equal opportunity employer, is operated by the Los Alamos National Security, LLC for the National Nuclear Security Administration of the U.S. Department of Energy under contract DE-AC52-06NA25396. By approving this article, the publisher recognizes that the U.S. Government retains nonexclusive, royalty-free license to publish or reproduce the published form of this contribution, or to allow others to do so, for U.S. Government purposes. Los Alamos National Laboratory requests that the publisher identify this article as work performed under the auspices of the U.S. Department of Energy. Los Alamos National Laboratory strongly supports academic freedom and a researcher's right to publish; as an institution, however, the Laboratory does not endorse the viewpoint of a publication or guarantee its technical correctness.

# From the Mimetic Finite Difference method to the Virtual Element Method

Gianmarco Manzini

Los Alamos National Laboratory  
New Mexico, USA

July 27, 2012 - Columbia University

# Outline

- 1 the Virtual Element Method (VEM) for the Laplace operator:
  - the degrees of freedom and the local Virtual Element (VE) space;
  - the abstract VE formulation;
  - the convergence theorem; consistency, stability;
  - the mimetic approximation of the VE bilinear form;
  - high-order and high-regular extensions.
2. A numerical experiment.
3. Final remarks, future work.

# The linear diffusion problem

- Differential formulation:

$$\begin{aligned}-\nabla u &= f \quad \text{in } \Omega, \\ u &= g \quad \text{on } \Gamma,\end{aligned}$$

- Variational formulation:

*Find  $u \in H_g^1(\Omega)$  such that:*

$$\int_{\Omega} \nabla u \cdot \nabla v \, dV = \int_{\Omega} f v \, dV \quad \forall v \in H_0^1(\Omega),$$

# People and References

- People:

- ▶ the "Pavia team": L. Beirão da Veiga, F. Brezzi, A. Cangiani, D. Marini, A. Russo;
- ▶ the "Los Alamos team": K. Lipnikov, D. Svyatskiy, M. Shashkov;

- Papers:

1. *F. Brezzi, A. Buffa, K. Lipnikov*, M2AN (2009): the low-order node-based MFD;
2. *L. Beirão da Veiga, K. Lipnikov, G. Manzini*, SINUM (2011): the arbitrary-order node-based MFD;
3. *the "six-name paper"*, M3AS (to appear in January 2013): basic principles of VEM; abstract formulation

# People and References

- People:

- ▶ the "Pavia team": L. Beirão da Veiga, F. Brezzi, A. Cangiani, D. Marini, A. Russo;
- ▶ the "Los Alamos team": K. Lipnikov, D. Svyatskiy, M. Shashkov;

- Papers:

1. *F. Brezzi, A. Buffa, K. Lipnikov*, M2AN (2009): the low-order node-based MFD;
2. *L. Beirão da Veiga, K. Lipnikov, G. Manzini*, SINUM (2011): the arbitrary-order node-based MFD;
3. *the "six-name paper"*, M3AS (to appear in January 2013): basic principles of VEM; abstract formulation

# People and References

- People:

- ▶ the "Pavia team": L. Beirão da Veiga, F. Brezzi, A. Cangiani, D. Marini, A. Russo;
- ▶ the "Los Alamos team": K. Lipnikov, D. Svyatskiy, M. Shashkov;

- Papers:

1. *F. Brezzi, A. Buffa, K. Lipnikov*, M2AN (2009): the low-order node-based MFD;
2. *L. Beirão da Veiga, K. Lipnikov, G. Manzini*, SINUM (2011): the arbitrary-order node-based MFD;
3. *the "six-name paper"*, M3AS (to appear in January 2013): basic principles of VEM; abstract formulation

# People and References

- People:

- ▶ the "Pavia team": L. Beirão da Veiga, F. Brezzi, A. Cangiani, D. Marini, A. Russo;
- ▶ the "Los Alamos team": K. Lipnikov, D. Svyatskiy, M. Shashkov;

- Papers:

1. *F. Brezzi, A. Buffa, K. Lipnikov*, M2AN (2009): the low-order node-based MFD;
2. *L. Beirão da Veiga, K. Lipnikov, G. Manzini*, SINUM (2011): the arbitrary-order node-based MFD;
3. *the "six-name paper"*, M3AS (to appear in January 2013): basic principles of VEM; abstract formulation



# People and References

- People:

- ▶ the "Pavia team": L. Beirão da Veiga, F. Brezzi, A. Cangiani, D. Marini, A. Russo;
- ▶ the "Los Alamos team": K. Lipnikov, D. Svyatskiy, M. Shashkov;

- Papers:

1. *F. Brezzi, A. Buffa, K. Lipnikov*, M2AN (2009): the low-order node-based MFD;
2. *L. Beirão da Veiga, K. Lipnikov, G. Manzini*, SINUM (2011): the arbitrary-order node-based MFD;
3. ***the "six-name paper"***, M3AS (to appear in January 2013): basic principles of VEM; abstract formulation

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^V$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space
    - $\mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- *Remarks:*
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to *compute* the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^{\mathcal{V}}$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space
    - $\mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- *Remarks:*
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to compute the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^V$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space  
 $\Rightarrow \mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- *Remarks:*
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to compute the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^\mathcal{V}$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space  
 $\Rightarrow \mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- *Remarks:*
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to compute the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^\mathcal{V}$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space  
 $\Rightarrow \mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- **Remarks:**
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to *compute* the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^\mathcal{V}$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space  
 $\Rightarrow \mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- **Remarks:**
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to *compute* the functions of  $\mathcal{V}_{h,P}$ !

# The Virtual Element approach

- The **Virtual Element** approach for the **Mimetic Finite Difference** (MFD) method is based on a **local finite element space**  $\mathcal{V}_{h,P}$  on  $P$  such that:
  - ▶ the degrees of freedom are the vertex values;  $\dim \mathcal{V}_{h,P} = N_P^\mathcal{V}$ ;
  - ▶ on triangles  $\mathcal{V}_{h,P}$  must be the linear Galerkin finite element space  
 $\Rightarrow \mathcal{V}_{h,P}$  must contain the linear polynomials  $1, x, y$ ;
  - ▶ the local spaces  $\mathcal{V}_{h,P}$  *glue gracefully* to give a conformal global finite element space  $\mathcal{V}_h$ .
- **Remarks:**
  - ▶ we will specify the behavior of the functions of  $\mathcal{V}_{h,P}$  on  $\partial P$ , the boundary of  $P$ ;
  - ▶ we will not ask to be able to *compute* the functions of  $\mathcal{V}_{h,P}$ !



# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}.$

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}$ .

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}$ .

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}.$

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}$ .

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}$ .

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:

- ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
- ▶  $\delta_i$  is continuous;
- ▶  $\delta_i$  is linear on each edge.

2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;

3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:

⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$   
(we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N_P}\}.$

# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:
  - ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
  - ▶  $\delta_i$  is continuous;
  - ▶  $\delta_i$  is linear on each edge.
2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;
3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:
  - ⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$  (we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N^P}\}.$



# The local finite element space

We define the local finite element space  $\mathcal{V}_{h,P}$  through a basis.

For each vertex  $v_i$  we define a function  $\varphi_i \in H^1(P)$ :

1. let  $\delta_i$  be the function defined on  $\partial P$  such that:
  - ▶  $\delta_i(v_j) = 1$  if  $i = j$ , and 0 otherwise;
  - ▶  $\delta_i$  is continuous;
  - ▶  $\delta_i$  is linear on each edge.
2. we set:  $\varphi_i|_{\partial P} = \delta_i$ ;
3. we **formally** extend  $\varphi_i|_{\partial P}$  inside  $P$  by the **harmonic lifting**:
  - ⇒ the functions  $\varphi_i$  are uniquely determined by the corresponding  $\delta_i$  (we can prove the unisolvency!)

Eventually, we set:  $\mathcal{V}_{h,P} := \text{span}\{\varphi_1, \dots, \varphi_{N^P}\}$ .

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N^P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N^P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N_P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N^P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N_P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N_P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N_P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.



# The harmonic lifting

- $\varphi_i$  is the harmonic function on  $P$  having  $\delta_i$  as boundary value:

$$\begin{cases} -\Delta\varphi_i = 0 & \text{in } \Omega \\ \varphi_i = \delta_i & \text{on } \partial\Omega. \end{cases}$$

- ▶ the functions  $\{\varphi_i\}$  are linearly independent;
- ▶ if  $w_h \in \mathcal{V}_{h,P}$ , then  $w_h = \sum_{i=1}^{N_P} w_h(v_i) \varphi_i$ ;
- ▶  $1, x, y \in \mathcal{V}_{h,P}$ ;
- ▶ the local spaces  $\mathcal{V}_{h,P}$  glue together giving a conformal finite element space  $\mathcal{V}_h \subset H_0^1(\Omega)$ .

- *Remarks:*

- ▶ if  $P$  is a triangle, we recover the  $\mathbb{P}_1$  Galerkin elements;
- ▶ if  $P$  is a parallelogram, we recover the  $\mathbb{Q}_1$  bilinear elements.

# The Harmonic Finite Element Method

The Harmonic Finite Element approximation of our elliptic problems is formally given by:

*Find  $u_h \in \mathcal{V}_h$  such that*

$$\mathcal{A}(u_h, v_h) = F_h(v_h) \quad \text{for all } v_h \in \mathcal{V}_h$$

where (as usual)

$$\mathcal{A}(u_h, v_h) = \int_{\Omega} \nabla u_h \cdot \nabla v_h$$

and  $F_h(v_h)$  is a suitable (and computable!) approximation of  $\int_{\Omega} f v$  (that uses only the vertex values of  $v_h$  and  $f$ ).

# The Harmonic Finite Element Method

The Harmonic Finite Element approximation of our elliptic problems is formally given by:

*Find  $u_h \in \mathcal{V}_h$  such that*

$$\mathcal{A}(u_h, v_h) = F_h(v_h) \quad \text{for all } v_h \in \mathcal{V}_h$$

where (as usual)

$$\mathcal{A}(u_h, v_h) = \int_{\Omega} \nabla u_h \cdot \nabla v_h$$

and  $F_h(v_h)$  is a suitable (and computable!) approximation of  $\int_{\Omega} f v$  (that uses only the vertex values of  $v_h$  and  $f$ ).

# The Harmonic Finite Element Method

The Harmonic Finite Element approximation of our elliptic problems is formally given by:

*Find  $u_h \in \mathcal{V}_h$  such that*

$$\mathcal{A}(u_h, v_h) = F_h(v_h) \quad \text{for all } v_h \in \mathcal{V}_h$$

where (as usual)

$$\mathcal{A}(u_h, v_h) = \int_{\Omega} \nabla u_h \cdot \nabla v_h$$

and  $F_h(v_h)$  is a suitable (and computable!) approximation of  $\int_{\Omega} f v$  (that uses only the vertex values of  $v_h$  and  $f$ ).

# The Harmonic Finite Element Method

Now, we are very happy, because...

- ...under *reasonable assumptions on the mesh*, the harmonic finite element approximation of an elliptic problem using the harmonic space  $\mathcal{V}_h$  enjoys the usual convergence properties!
- *Which assumptions?*
  - ▶ all geometric objects must scale properly:  $|P| \simeq h^2$ ,  $|e| \simeq h$ ;
  - ▶ each polygon is star-shaped (or the union of a *Uniformly bounded* number of star-shaped subcells) with respect to an internal ball of points (see Brenner-Scott, etc);
  - ▶ ...

# The Harmonic Finite Element Method

Now, we are very happy, because...

- ...under *reasonable assumptions on the mesh*, the harmonic finite element approximation of an elliptic problem using the harmonic space  $\mathcal{V}_h$  enjoys the usual convergence properties!

- *Which assumptions?*

- ▶ all geometric objects must scale properly:  $|P| \simeq h^2$ ,  $|e| \simeq h$ ;
- ▶ each polygon is star-shaped (or the union of a *Uniformly bounded* number of star-shaped subcells) with respect to an internal ball of points (see Brenner-Scott, etc);
- ▶ ...

# The Harmonic Finite Element Method

Now, we are very happy, because...

- ...under *reasonable assumptions on the mesh*, the harmonic finite element approximation of an elliptic problem using the harmonic space  $\mathcal{V}_h$  enjoys the usual convergence properties!
- *Which assumptions?*
  - ▶ all geometric objects must scale properly:  $|P| \simeq h^2$ ,  $|e| \simeq h$ ;
  - ▶ each polygon is star-shaped (or the union of a *uniformly bounded* number of star-shaped subcells) with respect to an internal ball of points (see Brenner-Scott, etc);
  - ▶ ...

# The Harmonic Finite Element Method

Now, we are very happy, because...

- ...under *reasonable assumptions on the mesh*, the harmonic finite element approximation of an elliptic problem using the harmonic space  $\mathcal{V}_h$  enjoys the usual convergence properties!
- *Which assumptions?*
  - ▶ all geometric objects must scale properly:  $|P| \simeq h^2$ ,  $|e| \simeq h$ ;
  - ▶ each polygon is star-shaped (or the union of a *uniformly bounded* number of star-shaped subcells) with respect to an internal ball of points (see Brenner-Scott, etc);
  - ▶ ...



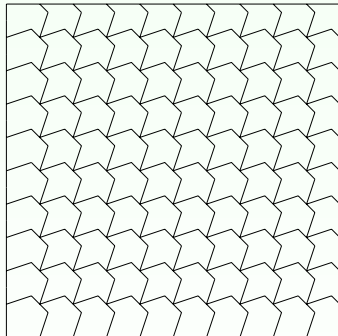
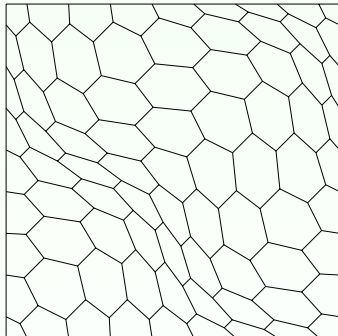
# The Harmonic Finite Element Method

Now, we are very happy, because...

- ...under *reasonable assumptions on the mesh*, the harmonic finite element approximation of an elliptic problem using the harmonic space  $\mathcal{V}_h$  enjoys the usual convergence properties!
- *Which assumptions?*
  - ▶ all geometric objects must scale properly:  $|P| \simeq h^2$ ,  $|e| \simeq h$ ;
  - ▶ each polygon is *star-shaped* (or the union of a *uniformly bounded* number of *star-shaped subcells*) with respect to an internal ball of points (see Brenner-Scott, etc);
  - ▶ ...

# Polygonal meshes

Examples: convex and non-convex polygonal cells



# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- Here, the mimetic technology comes into play!

# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- Here, the mimetic technology comes into play!

# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- Here, the mimetic technology comes into play!

# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- Here, the mimetic technology comes into play!

# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- Here, the mimetic technology comes into play!

# The Virtual Element Method

- So, we have a very nice method that works on polygonal meshes with very general shapes (also non-convex cells) and with a solid mathematical foundation (a priori error estimates, etc);
- we can also extend it to higher order polynomials (considering additional degrees of freedom)...

...BUT...

- ...if we do not know how to compute explicitly the basis functions...
- ...we don't know how to compute the stiffness matrix

$$\mathcal{A}(\varphi_i, \varphi_j) = \int_{\Omega} \nabla \varphi_i \cdot \nabla \varphi_j$$

and the right-hand side  $F_h(\mathbf{v}_h)$ !

- **Here, the mimetic technology comes into play!**



# Mimetic approximation of the bilinear form $\mathcal{A}(\varphi_i, \varphi_j)$

- Let  $\mathcal{A}_h$  be such approximation, i.e.,  $\mathcal{A}_h(\varphi_i, \varphi_j) \approx \mathcal{A}(\varphi_i, \varphi_j)$ .
- If  $\mathcal{A}_P$  is the restriction of  $\mathcal{A}$  to the polygon  $P$

$$\mathcal{A}(v_h, w_h) = \sum_P \mathcal{A}_P(v|_P, w|_P) = \sum_P \int_P \nabla v \cdot \nabla w$$

it is natural to assume that  $\mathcal{A}_h$  can be split in the same way:

$$\mathcal{A}_h(v_h, w_h) = \sum_P \mathcal{A}_{h,P}(v_h|_P, w_h|_P).$$

- Now, we give two conditions on  $\mathcal{A}_{h,P}$  that will guarantee the convergence: **consistency** and **stability**.

# Mimetic approximation of the bilinear form $\mathcal{A}(\varphi_i, \varphi_j)$

- Let  $\mathcal{A}_h$  be such approximation, i.e.,  $\mathcal{A}_h(\varphi_i, \varphi_j) \approx \mathcal{A}(\varphi_i, \varphi_j)$ .
- If  $\mathcal{A}_P$  is the restriction of  $\mathcal{A}$  to the polygon  $P$

$$\mathcal{A}(v_h, w_h) = \sum_P \mathcal{A}_P(v|_P, w|_P) = \sum_P \int_P \nabla v \cdot \nabla w$$

it is natural to assume that  $\mathcal{A}_h$  can be split in the same way:

$$\mathcal{A}_h(v_h, w_h) = \sum_P \mathcal{A}_{h,P}(v_h|_P, w_h|_P).$$

- Now, we give two conditions on  $\mathcal{A}_{h,P}$  that will guarantee the convergence: **consistency** and **stability**.

# Mimetic approximation of the bilinear form $\mathcal{A}(\varphi_i, \varphi_j)$

- Let  $\mathcal{A}_h$  be such approximation, i.e.,  $\mathcal{A}_h(\varphi_i, \varphi_j) \approx \mathcal{A}(\varphi_i, \varphi_j)$ .
- If  $\mathcal{A}_P$  is the restriction of  $\mathcal{A}$  to the polygon  $P$

$$\mathcal{A}(v_h, w_h) = \sum_P \mathcal{A}_P(v|_P, w|_P) = \sum_P \int_P \nabla v \cdot \nabla w$$

it is natural to assume that  $\mathcal{A}_h$  can be split in the same way:

$$\mathcal{A}_h(v_h, w_h) = \sum_P \mathcal{A}_{h,P}(v_h|_P, w_h|_P).$$

- Now, we give two conditions on  $\mathcal{A}_{h,P}$  that will guarantee the convergence: **consistency** and **stability**.

# Mimetic approximation of the bilinear form $\mathcal{A}(\varphi_i, \varphi_j)$

- Let  $\mathcal{A}_h$  be such approximation, i.e.,  $\mathcal{A}_h(\varphi_i, \varphi_j) \approx \mathcal{A}(\varphi_i, \varphi_j)$ .
- If  $\mathcal{A}_P$  is the restriction of  $\mathcal{A}$  to the polygon  $P$

$$\mathcal{A}(v_h, w_h) = \sum_P \mathcal{A}_P(v|_P, w|_P) = \sum_P \int_P \nabla v \cdot \nabla w$$

it is natural to assume that  $\mathcal{A}_h$  can be split in the same way:

$$\mathcal{A}_h(v_h, w_h) = \sum_P \mathcal{A}_{h,P}(v_{h|P}, w_{h|P}).$$

- Now, we give two conditions on  $\mathcal{A}_{h,P}$  that will guarantee the convergence: **consistency** and **stability**.

# Consistency and Stability $\Rightarrow$ Convergence

Six-name paper: *Basic Principles of Virtual Elements*, M3AS, to appear

**Theorem.** Assume that for each polygonal cell  $P$  the bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  satisfies the following properties:

- **Consistency:** for all  $q \in \mathbb{P}_1(P)$  and for all  $v_h \in \mathcal{V}_{h,P}$ :

$$\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}(v_h, q)$$

(an exactness property on linear polynomials).

- **Stability:** there exist two positive constants  $\alpha^*$  and  $\alpha_*$  independent of  $P$ , such that

$$\alpha_* \mathcal{A}_P(v_h, v_h) \leq \mathcal{A}_{h,P}(v_h, v_h) \leq \alpha^* \mathcal{A}_P(v_h, v_h).$$

# Consistency and Stability $\Rightarrow$ Convergence

Six-name paper: *Basic Principles of Virtual Elements*, M3AS, to appear

**Theorem.** Assume that for each polygonal cell  $P$  the bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  satisfies the following properties:

- **Consistency:** for all  $q \in \mathbb{P}_1(P)$  and for all  $v_h \in \mathcal{V}_{h,P}$ :

$$\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}(v_h, q)$$

(an exactness property on linear polynomials).

- **Stability:** there exist two positive constants  $\alpha^*$  and  $\alpha_*$  independent of  $P$ , such that

$$\alpha_* \mathcal{A}_P(v_h, v_h) \leq \mathcal{A}_{h,P}(v_h, v_h) \leq \alpha^* \mathcal{A}_P(v_h, v_h).$$

# Consistency and Stability $\Rightarrow$ Convergence

Six-name paper: *Basic Principles of Virtual Elements*, M3AS, to appear

**Theorem.** Assume that for each polygonal cell  $P$  the bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  satisfies the following properties:

- **Consistency:** for all  $q \in \mathbb{P}_1(P)$  and for all  $v_h \in \mathcal{V}_{h,P}$ :

$$\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}(v_h, q)$$

(an exactness property on linear polynomials).

- **Stability:** there exist two positive constants  $\alpha^*$  and  $\alpha_*$  independent of  $P$ , such that

$$\alpha_* \mathcal{A}_P(v_h, v_h) \leq \mathcal{A}_{h,P}(v_h, v_h) \leq \alpha^* \mathcal{A}_P(v_h, v_h).$$

# Consistency and Stability $\Rightarrow$ Convergence

Six-name paper: *Basic Principles of Virtual Elements*, M3AS, to appear

**Theorem.** Assume that for each polygonal cell  $P$  the bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  satisfies the following properties:

- **Consistency:** for all  $q \in \mathbb{P}_1(P)$  and for all  $v_h \in \mathcal{V}_{h,P}$ :

$$\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}(v_h, q)$$

(an exactness property on linear polynomials).

- **Stability:** there exist two positive constants  $\alpha^*$  and  $\alpha_*$  independent of  $P$ , such that

$$\alpha_* \mathcal{A}_P(v_h, v_h) \leq \mathcal{A}_{h,P}(v_h, v_h) \leq \alpha^* \mathcal{A}_P(v_h, v_h).$$

Let  $u_h \in \mathcal{V}_h$  be such that  $\mathcal{A}_h(u_h, v_h) = F_h(v_h)$  for all  $v_h \in \mathcal{V}_h$ .



# Consistency and Stability $\Rightarrow$ Convergence

Six-name paper: *Basic Principles of Virtual Elements*, M3AS, to appear

**Theorem.** Assume that for each polygonal cell  $P$  the bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  satisfies the following properties:

- **Consistency:** for all  $q \in \mathbb{P}_1(P)$  and for all  $v_h \in \mathcal{V}_{h,P}$ :

$$\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}(v_h, q)$$

(an exactness property on linear polynomials).

- **Stability:** there exist two positive constants  $\alpha^*$  and  $\alpha_*$  independent of  $P$ , such that

$$\alpha_* \mathcal{A}_P(v_h, v_h) \leq \mathcal{A}_{h,P}(v_h, v_h) \leq \alpha^* \mathcal{A}_P(v_h, v_h).$$

Then:

$$\|u - u_h\|_{H^1(\Omega)} \leq Ch \|u\|_{H^2(\Omega)}.$$

## A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ ).
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

- $\overline{\nabla v_h}$  is a constant vector in  $\mathbb{R}^2$ .

## A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ ).
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

- $\overline{\nabla v_h}$  is a constant vector in  $\mathbb{R}^2$ .

## A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ )
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

$$\underbrace{\int_P \nabla v_h = \int_{\partial P} v_h \mathbf{n}_P}_{(\text{Gauss--Green})}$$

## A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ )
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

$$\int_P \nabla v_h = \underbrace{\int_{\partial P} v_h \mathbf{n}_P}_{\text{split the boundary integral}} = \sum_{i=1}^{N_P} \left( \int_{e_i} v_h \right) \mathbf{n}_{P,i}$$

# A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ )
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

$$\int_P \nabla v_h = \int_{\partial P} v_h \mathbf{n}_P = \sum_{i=1}^{N^P} \underbrace{\left( \int_{\mathbf{e}_i} v_h \right)}_{v_h|_{\mathbf{e}} \in \mathbb{P}_1(\mathbf{e})} \mathbf{n}_{P,i} = \sum_{i=1}^{N^P} \underbrace{\frac{v_h(\mathbf{v}_i) + v_h(\mathbf{v}_{i+1})}{2} |\mathbf{e}_i|}_{\text{trapezoidal rule}} \mathbf{n}_{P,i}$$

# A crucial remark

- How can we define a local bilinear form  $\mathcal{A}_{h,P}(\cdot, \cdot)$  with the properties of consistency and stability? (Remember that we know the functions  $v_h$  of  $\mathcal{V}_{h,P}$  only on the boundary of  $P$ )
- If  $v_h \in \mathcal{V}_{h,P}$ , we can compute the following quantity

$$\overline{\nabla v_h} := \frac{1}{|P|} \int_P \nabla v_h$$

using only the vertex values.

- In fact,

$$\int_P \nabla v_h = \int_{\partial P} v_h \mathbf{n}_P = \sum_{i=1}^{N^P} \left( \int_{\mathbf{e}_i} v_h \right) \mathbf{n}_{P,i} = \sum_{i=1}^{N^P} \frac{v_h(\mathbf{v}_i) + v_h(\mathbf{v}_{i+1})}{2} |\mathbf{e}_i| \mathbf{n}_{P,i}$$

- $\overline{\nabla v_h}$  is a constant vector in  $\mathbb{R}^2$ .

# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

Why not? If  $P$  is a triangle, we get the stiffness matrix of the linear Galerkin FEM!

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- approximates the gradients using only the vertex values:

$$\nabla (\Pi_{h,P} v_h) = \nabla v_h$$

- and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$



# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\mathcal{A}_P(\varphi_i, \varphi_j) := \int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j} =: \mathcal{A}_{h,P}(\varphi_i, \varphi_j)$$

Why not? If  $P$  is a triangle, we get the stiffness matrix of the linear Galerkin FEM!

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- approximates the gradients using only the vertex values:

$$\nabla(\Pi_{h,P} v_h) = \overline{\nabla v_h}$$

- and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$

# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\mathcal{A}_P(\varphi_i, \varphi_j) := \int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j} =: \mathcal{A}_{h,P}(\varphi_i, \varphi_j)$$

But  $\mathcal{A}_{h,P}(\varphi_i, \varphi_j)$  would have **rank 2** for any kind of polygons, thus leading to a **singular** approximation for  $\mathcal{A}_h$ !

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- approximates the gradients using only the vertex values:

$$\nabla (\Pi_{h,P} v_h) = \nabla \bar{v}_h$$

- and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$

# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\mathcal{A}_P(\varphi_i, \varphi_j) := \int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j} =: \mathcal{A}_{h,P}(\varphi_i, \varphi_j)$$

But  $\mathcal{A}_{h,P}(\varphi_i, \varphi_j)$  would have rank 2 for any kind of polygons, thus leading to a singular approximation for  $\mathcal{A}_h$ !

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- approximates the gradients using only the vertex values:

$$\nabla (\Pi_{h,P} v_h) = \overline{\nabla v_h}$$

- and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$

# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\mathcal{A}_P(\varphi_i, \varphi_j) := \int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j} =: \mathcal{A}_{h,P}(\varphi_i, \varphi_j)$$

But  $\mathcal{A}_{h,P}(\varphi_i, \varphi_j)$  would have rank 2 for any kind of polygons, thus leading to a singular approximation for  $\mathcal{A}_h$ !

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- ▶ approximates the gradients using only the vertex values:

$$\nabla (\Pi_{h,P} v_h) = \overline{\nabla v_h}$$

- ▶ and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$

# The local projector $\Pi_{h,P}$

- Now, we are really tempted to say that

$$\mathcal{A}_P(\varphi_i, \varphi_j) := \int_P \nabla \varphi_i \cdot \nabla \varphi_j \approx \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j} =: \mathcal{A}_{h,P}(\varphi_i, \varphi_j)$$

But  $\mathcal{A}_{h,P}(\varphi_i, \varphi_j)$  would have rank 2 for any kind of polygons, thus leading to a singular approximation for  $\mathcal{A}_h$ !

- Key idea: define a **local projection operator** for each polygonal cell  $P$

$$\Pi_{h,P} : \mathcal{V}_{h,P} \longrightarrow \mathbb{P}_1(P)$$

that

- approximates the gradients using only the vertex values:

$$\nabla (\Pi_{h,P} v_h) = \overline{\nabla v_h}$$

- and preserves the linear polynomials:

$$\Pi_{h,P} q = q \quad \text{for all } q \in \mathbb{P}_1(P).$$

# The mimetic bilinear form $\mathcal{A}_{h,P}$

We start writing that

$$\mathcal{A}_{h,P}(u_h, v_h) = \mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) + \mathcal{A}_{h,P}(u_h - \Pi_{h,P} u_h, v_h).$$

With an easy computation it can be shown that

$$\mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) = \mathcal{A}_P(\Pi_{h,P} u_h, \Pi_{h,P} v_h) := \mathcal{A}_{h,P}^0(u_h, v_h)$$

and

$$\mathcal{A}_{h,P}((I - \Pi_{h,P})u_h, v_h) = \mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) \longrightarrow \mathcal{A}_{h,P}^1(u_h, v_h)$$

We will set:

$$\mathcal{A}_{h,P} = \mathcal{A}_{h,P}^0 + \mathcal{A}_{h,P}^1 = \text{CONSISTENCY} + \text{STABILITY}$$

# The mimetic bilinear form $\mathcal{A}_{h,P}$

We start writing that

$$\mathcal{A}_{h,P}(u_h, v_h) = \mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) + \mathcal{A}_{h,P}(u_h - \Pi_{h,P} u_h, v_h).$$

With an easy computation it can be shown that

$$\mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) = \mathcal{A}_P(\Pi_{h,P} u_h, \Pi_{h,P} v_h) := \mathcal{A}_{h,P}^0(u_h, v_h)$$

and

$$\mathcal{A}_{h,P}((I - \Pi_{h,P})u_h, v_h) = \mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) \longrightarrow \mathcal{A}_{h,P}^1(u_h, v_h)$$

We will set:

$$\mathcal{A}_{h,P} = \mathcal{A}_{h,P}^0 + \mathcal{A}_{h,P}^1 = \text{CONSISTENCY} + \text{STABILITY}$$

# The mimetic bilinear form $\mathcal{A}_{h,P}$

We start writing that

$$\mathcal{A}_{h,P}(u_h, v_h) = \mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) + \mathcal{A}_{h,P}(u_h - \Pi_{h,P} u_h, v_h).$$

With an easy computation it can be shown that

$$\mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) = \mathcal{A}_P(\Pi_{h,P} u_h, \Pi_{h,P} v_h) := \mathcal{A}_{h,P}^0(u_h, v_h)$$

and

$$\mathcal{A}_{h,P}((I - \Pi_{h,P})u_h, v_h) = \mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) \longrightarrow \mathcal{A}_{h,P}^1(u_h, v_h)$$

We will set:

$$\mathcal{A}_{h,P} = \mathcal{A}_{h,P}^0 + \mathcal{A}_{h,P}^1 = \text{CONSISTENCY} + \text{STABILITY}$$



# The mimetic bilinear form $\mathcal{A}_{h,P}$

We start writing that

$$\mathcal{A}_{h,P}(u_h, v_h) = \mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) + \mathcal{A}_{h,P}(u_h - \Pi_{h,P} u_h, v_h).$$

With an easy computation it can be shown that

$$\mathcal{A}_{h,P}(\Pi_{h,P} u_h, \Pi_{h,P} v_h) = \mathcal{A}_P(\Pi_{h,P} u_h, \Pi_{h,P} v_h) := \mathcal{A}_{h,P}^0(u_h, v_h)$$

and

$$\mathcal{A}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) = \mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) \longrightarrow \mathcal{A}_{h,P}^1(u_h, v_h)$$

We will set:

$$\mathcal{A}_{h,P} = \mathcal{A}_{h,P}^0 + \mathcal{A}_{h,P}^1 = \text{CONSISTENCY} + \text{STABILITY}$$

# The mimetic bilinear form $\mathcal{A}_{h,P}$

We start writing that

$$\mathcal{A}_{h,P}(u_h, v_h) = \mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) + \mathcal{A}_{h,P}(u_h - \Pi_{h,P} u_h, v_h).$$

With an easy computation it can be shown that

$$\mathcal{A}_{h,P}(\Pi_{h,P} u_h, v_h) = \mathcal{A}_P(\Pi_{h,P} u_h, \Pi_{h,P} v_h) := \mathcal{A}_{h,P}^0(u_h, v_h)$$

and

$$\mathcal{A}_{h,P}((I - \Pi_{h,P})u_h, v_h) = \mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h) \longrightarrow \mathcal{A}_{h,P}^1(u_h, v_h)$$

We will set:

$$\mathcal{A}_{h,P} = \mathcal{A}_{h,P}^0 + \mathcal{A}_{h,P}^1 = \text{CONSISTENCY} + \text{STABILITY}$$

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P} \quad \text{and} \quad \Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P).$

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:
- $\mathcal{A}_{h,P}^0$  ensures the consistency condition:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned}\mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q).\end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P} \quad \text{and} \quad \Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P).$

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j)$$

- $\mathcal{A}_{h,P}^0$  ensures the consistency condition:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q). \end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P} \quad \text{and} \quad \Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P).$

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j$$

- $\mathcal{A}_{h,P}^0$  ensures the consistency condition:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q). \end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the consistency condition:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q). \end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\mathcal{A}_{h,P}^0(v_h, q) = \int_P \overline{\nabla v_h} \cdot \overline{\nabla q}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .



# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\mathcal{A}_{h,P}^0(v_h, q) = \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\mathcal{A}_{h,P}^0(v_h, q) = \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} \end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q \end{aligned}$$

- the remaining term is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the **consistency condition**:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q). \end{aligned}$$

- the second term  $\mathcal{A}_{h,P}^1$  is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The consistency term $\mathcal{A}_{h,P}^0$

Recall that:  $\nabla \Pi_{h,P} v_h = \overline{\nabla v_h} \quad \forall v_h \in \mathcal{V}_{h,P}$  and  $\Pi_{h,P} q = q \quad \forall q \in \mathbb{P}_1(P)$ .

- $\mathcal{A}_{h,P}^0$  is the “constant gradient approximation” of the stiffness matrix:

$$\mathcal{A}_{h,P}^0(\varphi_i, \varphi_j) = \mathcal{A}_P(\Pi_{h,P} \varphi_i, \Pi_{h,P} \varphi_j) = \int_P \nabla \Pi_{h,P} \varphi_i \cdot \nabla \Pi_{h,P} \varphi_j = \int_P \overline{\nabla \varphi_i} \cdot \overline{\nabla \varphi_j}$$

- $\mathcal{A}_{h,P}^0$  ensures the consistency condition:  $\mathcal{A}_{h,P}(v_h, q) = \mathcal{A}_P(v_h, q)$  for all  $q \in \mathbb{P}_1(P)$ ; in fact,

$$\begin{aligned} \mathcal{A}_{h,P}^0(v_h, q) &= \int_P \overline{\nabla v_h} \cdot \overline{\nabla q} = |P| \overline{\nabla v_h} \cdot \overline{\nabla q} = \left( \int_P \nabla v_h \right) \cdot \overline{\nabla q} \\ &= \int_P \nabla v_h \cdot \overline{\nabla q} = \int_P \nabla v_h \cdot \nabla q = \mathcal{A}_P(v_h, q). \end{aligned}$$

- the second term  $\mathcal{A}_{h,P}^1$  is zero because  $(I - \Pi_{h,P})q = 0$  if  $q \in \mathbb{P}_1(P)$ .

# The stability term $\mathcal{A}_{h,P}^1$

- We need to correct  $\mathcal{A}_{h,P}^0$  in such a way that:
  - ▶ consistency is not upset;
  - ▶ we get stability;
  - ▶ we can compute the correction!
- In the six-name paper we show that we can substitute the (non computable!) term  $\mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$  with

$$\mathcal{A}_{h,P}^1(u_h, v_h) := \mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$$

where  $\mathcal{S}_{h,P}$  can be **any symmetric and positive definite bilinear form** that behaves (asymptotically) like  $\mathcal{A}_P$  on the kernel of  $\Pi_{h,P}$ .

- Hence:

$$\mathcal{A}_{h,P}(u_h, v_h) := \boxed{\mathcal{A}_P(\Pi_{h,P}u_h, \Pi_{h,P}v_h)} + \boxed{\mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)}$$

CONSISTENCY

STABILITY

# The stability term $\mathcal{A}_{h,P}^1$

- We need to correct  $\mathcal{A}_{h,P}^0$  in such a way that:
  - ▶ consistency is not upset;
  - ▶ we get stability;
  - ▶ we can compute the correction!
- In the six-name paper we show that we can substitute the (non computable!) term  $\mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$  with

$$\mathcal{A}_{h,P}^1(u_h, v_h) := \mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$$

where  $\mathcal{S}_{h,P}$  can be **any symmetric and positive definite bilinear form** that behaves (asymptotically) like  $\mathcal{A}_P$  on the kernel of  $\Pi_{h,P}$ .

- Hence:

$$\mathcal{A}_{h,P}(u_h, v_h) := \boxed{\mathcal{A}_P(\Pi_{h,P}u_h, \Pi_{h,P}v_h)} + \boxed{\mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)}$$

CONSISTENCY

STABILITY



# The stability term $\mathcal{A}_{h,P}^1$

- We need to correct  $\mathcal{A}_{h,P}^0$  in such a way that:
  - ▶ consistency is not upset;
  - ▶ we get stability;
  - ▶ we can compute the correction!
- In the six-name paper we show that we can substitute the (non computable!) term  $\mathcal{A}_P((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$  with

$$\mathcal{A}_{h,P}^1(u_h, v_h) := \mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)$$

where  $\mathcal{S}_{h,P}$  can be **any symmetric and positive definite bilinear form** that behaves (asymptotically) like  $\mathcal{A}_P$  on the kernel of  $\Pi_{h,P}$ .

- Hence:

$$\mathcal{A}_{h,P}(u_h, v_h) := \boxed{\mathcal{A}_P(\Pi_{h,P}u_h, \Pi_{h,P}v_h)} + \boxed{\mathcal{S}_{h,P}((I - \Pi_{h,P})u_h, (I - \Pi_{h,P})v_h)}$$

CONSISTENCY

STABILITY

# Arbitrary-order polynomials

Let us integrate by parts on cell P:

$$\int_P \nabla u \cdot \nabla v = - \int_P \Delta u v + \sum_{e \in \partial e} \int_e \nabla u \cdot \mathbf{n}_{P,e} v.$$

# Arbitrary-order polynomials

Let us integrate by parts on cell P:

$$\int_P \nabla u \cdot \nabla v = - \int_P \underbrace{\Delta u}_{\text{not zero!}} v + \sum_{e \in \partial e} \int_e \nabla u \cdot \mathbf{n}_{P,e} v.$$

If  $u$  is a polynomial of degree  $m$  on P:

- $\Delta u$  is a polynomial of degree  $m - 2$ ;

# Arbitrary-order polynomials

Let us integrate by parts on cell P:

$$\int_P \nabla u \cdot \nabla v = - \int_P \underbrace{\Delta u}_{\text{not zero!}} v + \sum_{e \in \partial e} \int_e \underbrace{\nabla u \cdot \mathbf{n}_{P,e}}_{\text{not constant!}} v.$$

If  $u$  is a polynomial of degree  $m$  on P:

- $\Delta u$  is a polynomial of degree  $m - 2$ ;
- $\nabla u \cdot \mathbf{n}_{P,e}$  is a polynomial of degree  $m - 1$ ;

# Divergence term: internal degrees of freedom

1. We use the **moments of  $\mathbf{v}$**  to express the integral over  $P$ :

if

$$\Delta u = a_0 \mathbf{1} + a_1 \mathbf{x} + a_2 \mathbf{y} + \dots \in \mathbb{P}_{m-2}(P)$$

# Divergence term: internal degrees of freedom

1. We use the **moments of  $\mathbf{v}$**  to express the integral over  $P$ :

if

$$\Delta u = a_0 \mathbf{1} + a_1 \mathbf{x} + a_2 \mathbf{y} + \dots \in \mathbb{P}_{m-2}(P)$$

then

$$\begin{aligned} \int_P \Delta u \, \mathbf{v} &= a_0 \underbrace{\int_P \mathbf{1} \, \mathbf{v}}_{\hat{\mathbf{v}}_{P,0}} + a_1 \underbrace{\int_P \mathbf{x} \, \mathbf{v}}_{\hat{\mathbf{v}}_{P,1,\mathbf{x}}} + a_2 \underbrace{\int_P \mathbf{y} \, \mathbf{v}}_{\hat{\mathbf{v}}_{P,1,\mathbf{y}}} + \dots \\ &= a_0 \hat{\mathbf{v}}_{P,0} + a_1 \hat{\mathbf{v}}_{P,1,\mathbf{x}} + a_2 \hat{\mathbf{v}}_{P,1,\mathbf{y}} + \dots \end{aligned}$$

# Divergence term: internal degrees of freedom

1. We use the **moments of  $\mathbf{v}$**  to express the integral over  $P$ :

if

$$\Delta u = a_0 \mathbf{1} + a_1 \mathbf{x} + a_2 \mathbf{y} + \dots \in \mathbb{P}_{m-2}(P)$$

then

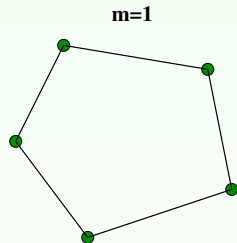
$$\begin{aligned} \int_P \Delta u \mathbf{v} &= a_0 \underbrace{\int_P \mathbf{1} \mathbf{v}}_{\hat{\mathbf{v}}_{P,0}} + a_1 \underbrace{\int_P \mathbf{x} \mathbf{v}}_{\hat{\mathbf{v}}_{P,1,\mathbf{x}}} + a_2 \underbrace{\int_P \mathbf{y} \mathbf{v}}_{\hat{\mathbf{v}}_{P,1,\mathbf{y}}} + \dots \\ &= a_0 \hat{\mathbf{v}}_{P,0} + a_1 \hat{\mathbf{v}}_{P,1,\mathbf{x}} + a_2 \hat{\mathbf{v}}_{P,1,\mathbf{y}} + \dots \end{aligned}$$

This choice suggests us to define

- $m(m-1)/2$  **internal** degrees of freedom  $\approx \hat{\mathbf{v}}_{P,0}, \hat{\mathbf{v}}_{P,1,\mathbf{x}}, \hat{\mathbf{v}}_{P,1,\mathbf{y}}, \dots$

# $C^0$ high-order approximations

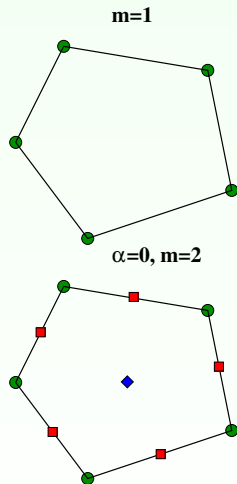
- The “ $C^0 - \mathbb{P}_1$ ” approximation requires:
  - one real number per mesh vertex  $v$ ;





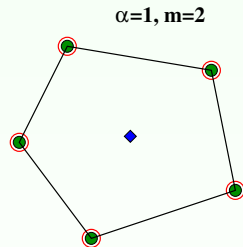
# $C^0$ high-order approximations

- The “ $C^0 - \mathbb{P}_1$ ” approximation requires:
  - one real number per mesh vertex  $v$ ;
- the “ $C^0 - \mathbb{P}_m$ ” approximations for  $m > 1$  require
  - one real number per mesh vertex  $v$ ;
  - $(m - 1)$  real numbers per mesh edge  $e$ ;
  - $m(m - 1)/2$  real numbers per mesh cell  $P$ ;



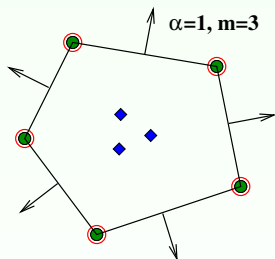
# Approximations with high regularity

- The “ $C^1 - \mathbb{P}_2$ ” approximation requires:
  - vertex dofs  $\rightarrow$  solution and derivatives at each vertex;
  - cell dofs  $\rightarrow$  solution moments inside the cells;



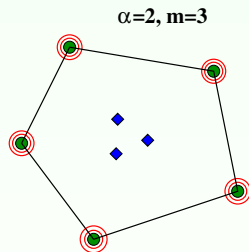
# Approximations with high regularity

- The “ $C^1 - \mathbb{P}_3$ ” approximation requires:
  - vertex dofs  $\rightarrow$  **solution** and **derivatives** at each vertex;
  - cell dofs  $\rightarrow$  **solution moments** inside the cells;
  - edge dofs  $\rightarrow$  *solution* and *normal derivatives* along the edges;



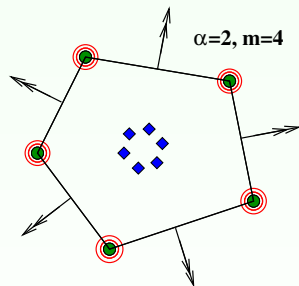
# Approximations with high regularity

- The “ $C^2 - \mathbb{P}_3$ ” approximation requires:
  - vertex dofs  $\rightarrow$  solution and derivatives at each vertex;
  - cell dofs  $\rightarrow$  solution moments inside the cells;



# Approximations with high regularity

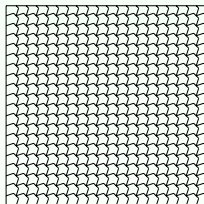
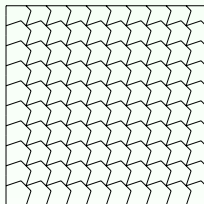
- The “ $C^2 - \mathbb{P}_4$ ” approximation requires:
  - vertex dofs  $\rightarrow$  **solution** and **derivatives** at each vertex;
  - cell dofs  $\rightarrow$  **solution moments** inside the cells;
  - edge dofs  $\rightarrow$  *solution* and *normal derivatives* along the edges;



# Numerical experiments

## Meshes with non-convex polygons

- **Meshes:**



- **Exact solution:**  $u(x, y) = e^{-2\pi y} \sin(2\pi x)$

- **Diffusion tensor**

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

# Continuous approximations

$\alpha = 0$ , non-convex polygons,  $\|\cdot\|_{1,h}$  errors, non-constant K

		<b>m = 1</b>		<b>m = 2</b>	
n	<i>h</i>	Error	Rate	Error	Rate
0	$1.458 \cdot 10^{-1}$	3.544	—	3.007	—
1	$7.289 \cdot 10^{-2}$	3.046	0.22	$8.081 \cdot 10^{-1}$	1.89
2	$3.644 \cdot 10^{-2}$	1.887	0.69	$2.071 \cdot 10^{-1}$	1.96
3	$1.822 \cdot 10^{-2}$	1.000	0.92	$5.303 \cdot 10^{-2}$	1.97
4	$9.111 \cdot 10^{-3}$	$5.154 \cdot 10^{-1}$	<b>0.98</b>	$1.348 \cdot 10^{-2}$	<b>1.98</b>

# High-regular approximations

$\alpha = 1, 2$ ; non-convex polygons,  $\|\cdot\|_{1,h}$  errors, non-constant  $K$

		$\alpha = 1, \mathbf{m} = 2$		$\alpha = 2, \mathbf{m} = 3$	
n	$h$	Error	Rate	Error	Rate
0	$1.458 \cdot 10^{-1}$	$8.901 \cdot 10^{-2}$	—	$1.054 \cdot 10^{-2}$	—
1	$7.289 \cdot 10^{-2}$	$1.983 \cdot 10^{-2}$	2.26	$4.543 \cdot 10^{-4}$	4.72
2	$3.644 \cdot 10^{-2}$	$4.815 \cdot 10^{-3}$	2.08	$4.663 \cdot 10^{-5}$	3.36
3	$1.822 \cdot 10^{-2}$	$1.198 \cdot 10^{-3}$	<b>2.03</b>	$5.528 \cdot 10^{-6}$	<b>3.11</b>



# Summary

- **VEM is a family of schemes on polygonal meshes:** new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $P_k(P)$  polynomials for the local VE space;
  - ▶ the stabilization term is chosen so that the scheme is stable;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- **VEM works for any order of accuracy:**
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- VEM works for any order of regularity:
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- **VEM works for any order of regularity:**
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.



# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- **VEM works for any order of regularity:**
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- VEM is a *family of schemes* on polygonal meshes: new schemes are generated by changing the stabilization term;
- VEM works for any order of accuracy:
  - ▶ we can use  $\mathbb{P}_k(P)$  polynomials for the local VE space;
  - ▶ "dofs" are vertex values, nodal values on the edges of  $\partial P$  and moments inside  $P$ ;
  - ▶ the behavior on  $\partial P$  is given by a polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.
- **VEM works for any order of regularity:**
  - ▶ we use also derivatives as degrees of freedom at vertices and edge nodes
  - ▶ the behavior on  $\partial P$  is given by a Hermite-like polynomial interpolation;
  - ▶ optimal error estimates in the energy norm are confirmed by experiments.

# Summary

- **VEM works on degenerate meshes (experiments):**
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶ VEM works on arbitrary polyhedral meshes (e.g. hexahedra, tetrahedra, prisms, pyramids, ...).
  - ▶  $C^0 - P_m$  ( $m \geq 1$ ) requires vortex values and moments on edges, faces, and inside  $P$ ;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- **VEM works on degenerate meshes (experiments):**
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶ VEM works on arbitrary polyhedra, including non-convex polyhedra;
  - ▶  $C^0 - P_m$  ( $m \geq 1$ ) requires vertex values and moments on edges, faces, and inside  $P$ ;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- **VEM works on degenerate meshes (experiments):**
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶ VEM works on arbitrary polyhedra (tetrahedra, hexahedra, etc.);
  - ▶  $C^0 - P_m$  ( $m \geq 1$ ) requires vertex values and moments on edges, faces, and inside  $P$ ;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- **VEM works on degenerate meshes (experiments):**
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - P_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - P_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside  $P$ ;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- **VEM works on degenerate meshes (experiments):**
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - P_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - P_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside  $P$ ;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- VEM works on degenerate meshes (experiments):
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - \mathbb{P}_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - \mathbb{P}_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside P;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.



# Summary

- VEM works on degenerate meshes (experiments):
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - \mathbb{P}_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - \mathbb{P}_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside P;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- VEM works on degenerate meshes (experiments):
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - \mathbb{P}_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - \mathbb{P}_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside P;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- VEM works on degenerate meshes (experiments):
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - \mathbb{P}_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - \mathbb{P}_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside P;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Summary

- VEM works on degenerate meshes (experiments):
  - ▶ meshes with convex and non-convex elements;
  - ▶ meshes with very stretched elements;
  - ▶ meshes with hanging nodes;
  - ▶ meshes with collapsing nodes.
- VEM can be generalized to 3-D polyhedral mesh (in progress):
  - ▶  $C^0 - \mathbb{P}_1$  works in 3-D just using vertex values as degrees of freedom (dofs);
  - ▶  $C^0 - \mathbb{P}_m$  ( $m > 1$ ) requires vertex values and moments on edges, faces, and inside P;
  - ▶ no need of numerical integration, VEM does not use the basis functions explicitly;
  - ▶ no need of isoparametric mappings, VEM works in the physical domain.
- There is no difference between VEM and the mimetic finite difference method, the two families of schemes coincide.

# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.

# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.

# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.

# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.



# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.

# Current/future developments

- full extension to three dimensional problems;
- other differential equations: elasticity, advection-diffusion, Stokes, etc;
- understand the role of the mimetic stabilization;
- justify the numerical results for degenerate meshes (not covered by the theory);

Thank for your attention.