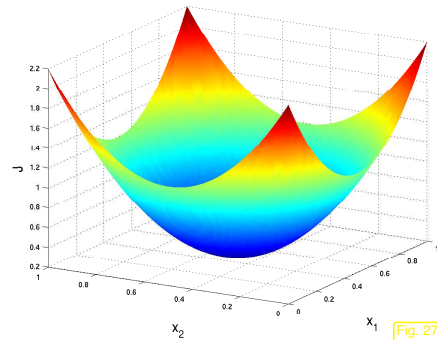


Example 31 (Quadratic functionals).

Analogy \rightarrow parabola:

$$\begin{array}{c} J(v) = \frac{1}{2}a(v, v) - f(v) \\ \uparrow \qquad \qquad \uparrow \\ f(x) = ax^2 + bx \end{array}$$

quadratic functional $\mathbb{R}^2 \mapsto \mathbb{R}$ \triangleright



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3.1.1 Galerkin discretization

Abstract discussion: start from **linear variational problem** (see Sect. 2.6, (2.6.1))

$$u \in V: \quad a(u, v) = f(v) \quad \forall v \in V, \quad (2.6.1)$$

V = Hilbert space with norm $\|\cdot\|_V$, $a(\cdot, \cdot)$ continuous bilinear form, f continuous linear form.

Norm of $a(\cdot, \cdot)$:

$$C_A := \sup_{v \in V \setminus \{0\}} \sup_{u \in V \setminus \{0\}} \frac{|a(u, v)|}{\|u\|_V \|v\|_V} < \infty$$

Assumption: V -ellipticity of $a(\cdot, \cdot)$, see Def. 2.6.6:

$$\exists \gamma > 0: \quad |a(u, u)| \geq \gamma \|u\|_V^2 \quad \forall u \in V. \quad (2.8.1)$$

Remark. If $a(\cdot, \cdot)$ symmetric (\rightarrow inner product, see Def. 2.6.3) and $\|\cdot\|_V$ = energy norm $\|\cdot\|_A$

$$\blacktriangleright \quad \gamma, C_A = 1 \quad (\text{prove with Cauchy-Schwarz inequality})$$

Idea of **Galerkin discretization**

Replace V in (2.6.1) with a **finite dimensional subspace** V_N (discrete trial/test space).

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Notation: N = formal index, tagging “discrete entities” (\rightarrow “finite amount of information”)

Discrete variational problem

$$u_N \in V_N: \quad a(u_N, v_N) = f(v_N) \quad \forall v_N \in V_N. \quad (3.1.1)$$

Lax-Milgram Lemma Thm. 2.6.7 \Rightarrow Existence & Uniqueness of solution $u_N \in V_N$, stability

$$\|u_N\|_V \leq \frac{1}{\gamma} \sup_{v_N \in V_N \setminus \{0\}} \frac{|f(v_N)|}{\|v_N\|_V}.$$

Issues:

1. How “accurate” is the **Galerkin solution** u_N ?
 - (a) What measure for accuracy?
 - (b) How to assess accuracy?
2. How to convert (3.1.1) into (linear) system of equations?

Ad 1(a): Focus on norm $\|\cdot\|_V$ (and $\|\cdot\|_A$, if $a(\cdot, \cdot)$ inner product)

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3

The Finite Element Method (FEM)

Problem : scalar second-order elliptic boundary value problem
 Perspective : **variational** interpretation in Sobolev spaces \rightarrow Sect. 2.6.2
 Objective : algorithm for the computation of an **approximate numerical solution**

3.1 Fundamentals

Moot point (\rightarrow Sect. 1.2): any computer can only handle a finite amount of information (reals)

Variational boundary value problem

DISCRETIZATION \rightarrow

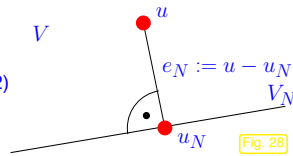
System of a finite number of equations for real unknowns

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Galerkin orthogonality

$$a(u - u_N, v_N) = 0 \quad \forall v_N \in V_N. \quad (3.1.2)$$

[Geometric meaning for inner product $a(\cdot, \cdot) \rightarrow$]



► **Discretization error** $e_N := u - u_N$ “ $a(\cdot, \cdot)$ -orthogonal” to discrete trial/test space V_N

Remark 32. If $a(\cdot, \cdot)$ is inner product on V : “Pythagoras’ theorem” \rightarrow Fig. 28

$$\|u - u_N\|_A^2 = \|u\|_A^2 - \|u_N\|_A^2. \quad (3.1.3)$$

(3.1.3) ► simple formula for computation of energy norm of Galerkin discretization error in numerical experiments with known u .

△

Theorem 3.1.1 (Cea’s lemma). If $a(\cdot, \cdot) =$ continuous, V -elliptic, bilinear form, $V_N \subset V$ finite dimensional subspace, $u \in V / u_N \in V_N$ solve (2.6.1)/(3.1.1), then

$$\|u - u_N\|_V \leq \frac{C_A}{\gamma} \inf_{v_N \in V_N \setminus \{0\}} \|u - v_N\|_V$$

Proof. By Galerkin-orthogonality (3.1.2), for all $v_N \in V_N$

$$\begin{aligned} \gamma \|u - u_N\|_V^2 &\leq |a(u - u_N, u - u_N) + a(u - u_N, u_N - v_N)| \\ &= |a(u - u_N, u - v_N)| \leq C_A \|u - u_N\|_V \|u - v_N\|_V \\ \Rightarrow \|u - u_N\|_V &\leq \frac{C_A}{\gamma} \inf_{v_N \in V_N \setminus \{0\}} \|u - v_N\|_V, \end{aligned}$$

because v_N arbitrary.

□

Quasi-optimality of Galerkin solutions: with $C > 0$ independent of u, V_N

$$\|u - u_N\|_V \leq C \inf_{v_N \in V_N} \|u - v_N\|_V, \quad (3.1.4)$$

$$\underbrace{\|u - u_N\|_V}_{\text{(norm of) discretization error}} \leq C \underbrace{\inf_{v_N \in V_N} \|u - v_N\|_V}_{\text{best approximation error}} \quad (3.1.5)$$

► 1(b): To assess accuracy of Galerkin solution: study capability of V_N to approximate u !

► “Monotonicity” of best approximation

$$\text{Trial test spaces } V_N, V'_N \subset V: V_N \subset V'_N \Rightarrow \inf_{v_N \in V'_N} \|u - v_N\|_V \leq \inf_{v_N \in V_N} \|u - v_N\|_V.$$

Enhance accuracy by enlarging (“refining”) trial space.

Reminder:

Definition 3.1.2 (Linear operator). Let V, W be real vector spaces. A mapping $T : V \mapsto W$ is a **linear operator**, if

$$T(\alpha u + \beta v) = \alpha T(u) + \beta T(v) \quad \forall u, v \in V, \forall \alpha, \beta \in \mathbb{R}.$$

Reminder:

Definition 3.1.3 (projection). A linear operator $P : V \mapsto V$ on a vector space V is a **projection**, if $P^2 = P$.

Definition 3.1.4 (Galerkin projection). Under the assumptions of Cea’s lemma Thm. 3.1.1 the **Galerkin projection** $P_N : V \mapsto V_N \subset V$ is defined by

$$a(P_N u, v_N) = a(u, v_N) \quad \forall v_N \in V_N.$$

[Lax-Milgram Lemma Thm 2.6.7 $\Rightarrow P_N$ well defined and continuous]

3.1.2 The (linear) algebraic setting

[Now we tackle issue 2. (conversion of (3.1.1) into system of equations)]

I.

Introduce (ordered) **basis** \mathfrak{B}_N of V_N :



$$\mathfrak{B}_N := \{b_N^1, \dots, b_N^N\} \subset V_N, \quad V_N = \text{Span}\{\mathfrak{B}_N\}, \quad N := \dim(V_N).$$

II. Basis representations

$$\begin{aligned} u_N &= \mu_1 b_N^1 + \dots + \mu_N b_N^N, \quad \mu_i \in \mathbb{R} \\ v_N &= \nu_1 b_N^1 + \dots + \nu_N b_N^N, \quad \nu_i \in \mathbb{R} \end{aligned} \quad \text{in (3.1.1).}$$

$$(3.1.1): \quad a(u_N, v_N) = f(v_N) \quad \forall v_N \in V_N$$

$$a(\mu_1 b_N^1 + \dots + \mu_N b_N^N, \nu_1 b_N^1 + \dots + \nu_N b_N^N) = f(\nu_1 b_N^1 + \dots + \nu_N b_N^N) \quad \forall \nu_1, \dots, \nu_N \in \mathbb{R},$$



$$\sum_{k=1}^N \sum_{j=1}^N \mu_k \nu_j a(b_N^k, b_N^j) = \sum_{j=1}^N \nu_j f(b_N^j) \quad \forall \nu_1, \dots, \nu_N \in \mathbb{R},$$



$$\sum_{j=1}^N \nu_j \left(\sum_{k=1}^N \mu_k a(b_N^k, b_N^j) - f(b_N^j) \right) = 0 \quad \forall \nu_1, \dots, \nu_N \in \mathbb{R},$$



$$\sum_{k=1}^N \mu_k a(b_N^k, b_N^j) = f(b_N^j) \quad \text{for } j = 1, \dots, N.$$



$$\boxed{\mathbf{A} \vec{\mu} = \vec{\varphi}}, \quad \mathbf{A} = \left(a(b_N^k, b_N^j) \right)_{j,k=1}^N \in \mathbb{R}^{N,N}, \quad \vec{\varphi} = \left(f(b_N^j) \right)_{j=1}^N, \\ \vec{\mu} = (\mu_1, \dots, \mu_N)^T \in \mathbb{R}^N$$

Discrete variational problem
 $u_N \in V_N: a(u_N, v_N) = f(v_N) \quad \forall v_N \in V_N$

Choosing basis \mathfrak{B}_N

Linear system
of equations
 $\mathbf{A} \vec{\mu} = \vec{\varphi}$

Stiffness matrix: $\mathbf{A} = \left(a(b_N^k, b_N^j) \right)_{j,k=1}^N \in \mathbb{R}^{N,N},$

Load vector: $\vec{\varphi} = \left(f(b_N^j) \right)_{j=1}^N \in \mathbb{R}^N,$

Coefficient vector: $\vec{\mu} = (\mu_1, \dots, \mu_N)^T \in \mathbb{R}^N,$

Recovery of solution: $u_N = \sum_{k=1}^N \mu_k b_N^k.$

Corollary 3.1.5.

(3.1.1) has unique solution $\Leftrightarrow \mathbf{A}$ regular

Impact of choice of basis ?

Choice of \mathfrak{B}_N does **not** affect $u_N \Rightarrow$ No impact on discretization error !

Properties of matrix \mathbf{A} crucially depend on basis \mathfrak{B}_N !

Lemma 3.1.6. Consider (3.1.1) and two bases of V_N ,

$$\mathfrak{B}_N := \{b_N^1, \dots, b_N^N\}, \quad \underline{\mathfrak{B}}_N := \{\underline{b}_N^1, \dots, \underline{b}_N^N\},$$

related by

$$\underline{b}_N^j = \sum_{k=1}^N s_{jk} b_N^k \quad \text{with } \mathbf{S} = (s_{jk})_{j,k=1}^N \in \mathbb{R}^{N,N} \text{ regular.}$$

► Stiffness matrices $\mathbf{A}, \underline{\mathbf{A}} \in \mathbb{R}^{N,N}$, load vectors $\vec{\varphi}, \underline{\vec{\varphi}} \in \mathbb{R}^N$, and coefficient vectors $\vec{\mu}, \underline{\vec{\mu}} \in \mathbb{R}^N$, respectively, satisfy

$$\underline{\mathbf{A}} = \mathbf{S} \mathbf{A} \mathbf{S}^T, \quad \underline{\vec{\varphi}} = \mathbf{S} \vec{\varphi}, \quad \underline{\vec{\mu}} = \mathbf{S}^{-T} \vec{\mu}. \quad (3.1.6)$$

Proof.

$$\underline{\mathbf{A}}_{lm} = a(\underline{b}_N^m, \underline{b}_N^l) = \sum_{k=1}^N \sum_{j=1}^N s_{mk} a(b_N^k, b_N^j) s_{lj} = \sum_{k=1}^N \left(\sum_{j=1}^N s_{lj} \mathbf{A}_{jk} \right) s_{mk} = (\mathbf{S} \mathbf{A} \mathbf{S}^T)_{lm},$$

$(\mathbf{S} \mathbf{A})_{lk}$

Reminder of linear algebra:

Definition 3.1.7 (Congruent matrices). Two matrices $\mathbf{A} \in \mathbb{R}^{N,N}$, $\mathbf{B} \in \mathbb{R}^{N,N}$, $N \in \mathbb{N}$, are called **congruent**, if there is a regular matrix $\mathbf{S} \in \mathbb{R}^{N,N}$ such that $\mathbf{B} = \mathbf{S} \mathbf{A} \mathbf{S}^T$.

► Equivalence relation on square matrices

Lemma 3.1.8. Matrix property invariant under congruence \Leftrightarrow Property of stiffness matrix invariant under change of basis \mathfrak{B}_N

Matrix properties invariant under congruence =

- regularity
- symmetry
- positive definiteness

Reminder:

Definition 3.1.9 (Positive definite matrix). Matrix $\mathbf{B} \in \mathbb{R}^{N,N}$, $N \in \mathbb{N}$, is **positive definite** $\Leftrightarrow \vec{\xi}^T \mathbf{B} \vec{\xi} > 0$ for all $\vec{\xi} \in \mathbb{R}^N \setminus \{0\}$.

3.1.3 Principles of FEM

1D case → Sect. 1.2.3.1, now higher dimensional, “complicated” domain Ω :

$\Omega \subset \mathbb{R}^d$, $d = 2, 3$, bounded **computational domain**: assumed polygonal $d = 2$, polyhedral $d = 3$

First main ingredient: **triangulation/mesh** of Ω , cf. Sect. 1.2.1

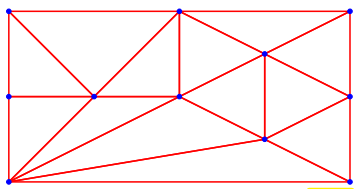
Definition 3.1.10. A **mesh** (or **triangulation**) of $\Omega \subset \mathbb{R}^d$ is a finite collection $\{K_i\}_{i=1}^M$, $M \in \mathbb{N}$, of open non-degenerate polygons ($d = 2$)/polyhedra ($d = 3$) such that

- (A) $\bar{\Omega} = \bigcup \{\bar{K}_i, i = 1, \dots, M\}$,
- (B) $K_i \cap K_j = \emptyset \Leftrightarrow i \neq j$,
- (C) for all $i, j \in \{1, \dots, M\}$, $i \neq j$, the intersection $\bar{K}_i \cap \bar{K}_j$ is a vertex, edge, or face of both K_i and K_j .

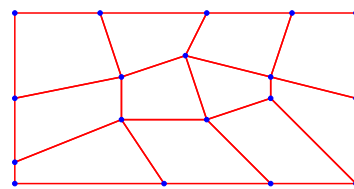
► “vertex”, “edge”, “face” of polygon/polyhedron: → geometric intuition

Terminology: Given mesh $\mathcal{M} := \{K_i\}_{i=1}^M$: K_i called **cell** or **element**.
Vertices of a mesh → **nodes** (set $\mathcal{N}(\mathcal{M})$)

Types of meshes:



Triangular mesh in 2D



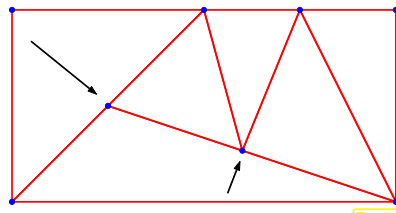
Quadrilateral mesh in 2D

If (C) does not hold

► Triangular **non-conforming** mesh
(with **hanging nodes**)

$\bar{K}_i \cap \bar{K}_j$ is only part of an edge/face for at most one of the adjacent cells.

(However, conforming if degenerate quadrilaterals admitted)



Simplicial mesh = triangular mesh in 2D
tetrahedral mesh in 3D

Second main ingredient: space of **piecewise polynomial functions**

$$V_N := \{v \in V : v|_K \in \mathcal{P}_p(K) \forall K \in \mathcal{M}\},$$

$$\mathcal{P}_p(K) = \text{polynomials of degree } \leq p \text{ on cell } K.$$

Note:

$v \in V \rightarrow$ **conformity conditions** at interelement boundaries

Lemma 3.1.11 (Conformity condition for H^1). Let $\mathcal{M} := \{K_i\}_{i=1}^M$ be a triangulation (→ Def. 3.1.10) of $\Omega \subset \mathbb{R}^d$ and assume that $v : \Omega \mapsto \mathbb{R}$ satisfies that $v|_K$ can be extended to a function in $C^\infty(\bar{K})$ for any $K \in \mathcal{M}$. Then

$$v \in H^1(\Omega) \Leftrightarrow v \in C^0(\bar{\Omega}).$$

► Conformity condition for H^1 = global continuity (C^0 , **not** C^1 ! → Ex. 28)
(recall physical constraints on temperature distributions!)

Thanks to notion of weak derivative, Sect. 2.4 !

Definition 3.1.12 (Conformity). Let V be a function space. A \mathcal{M} -piecewise polynomial space V_N is called **V-conforming**, if $V_N \subset V$.

Third main ingredient: **Locally supported basis functions**

Basis functions b_N^1, \dots, b_N^N for a finite element trial/test space V_N built on a mesh \mathcal{M} satisfy:

- each b_N^i **associated** with a single cell/edge/face/vertex of \mathcal{M} ,
- $\text{supp}(b_N^i) = \bigcup \{\bar{K} : K \in \mathcal{M}, \mathbf{p} \subset \bar{K}\}$, if b_N^i associated with cell/edge/face/vertex \mathbf{p} .

Finite element terminology: b_N^i = **global shape/basis functions**

Example 33 (Supports of global shape functions in 1D). → Sect. 1.2.3.1

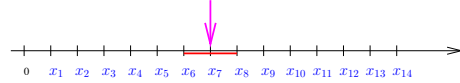
• $\Omega =]a, b[\hat{=}$ interval

• Equidistant mesh

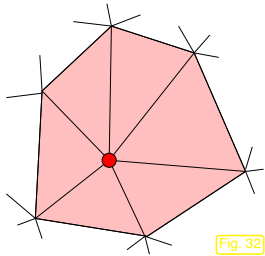
$$\mathcal{M} := \{x_{j-1}, x_j[, j = 1, \dots, N\},$$

$$x_j := a + hj, h := (b - a)/N, N \in \mathbb{N}.$$

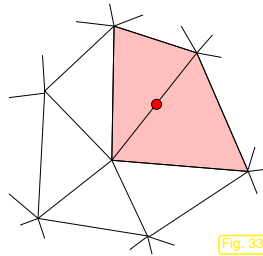
Support of global shape function associated with x_7



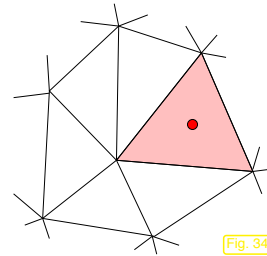
Example 34 (Supports of global shape functions on triangular mesh).



Support of node-associated basis function



Support of edge-associated basis function



Support of cell-associated basis function

Rationale for small supports ?

Recall bilinear form $\leftrightarrow -\Delta$:

$$a(u, v) := \int_{\Omega} \text{grad } u \cdot \text{grad } v \, dx$$

Use triangular mesh \mathcal{M} , test/trial space $V_N \subset H^1(\Omega)$ with basis $\mathfrak{B}_N := \{b_N^1, \dots, b_N^N\}$ (→ Sect. 3.1.2)

► Stiffness matrix $\mathbf{A} \in \mathbb{R}^{N,N}$ with $a_{ij} := a(b_N^j, b_N^i), i, j = 1, \dots, N$

b_N^i, b_N^j associated with nodes
not linked by an edge

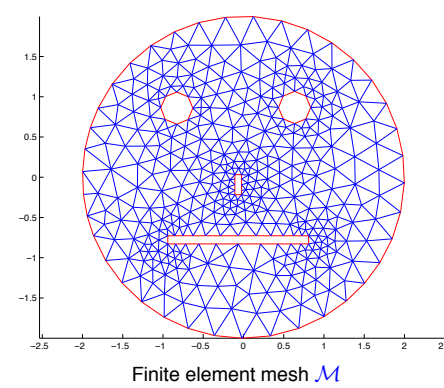
$a_{ij} = 0$
(because
 $\text{vol}(\text{supp}(b_N^i) \cap \text{supp}(b_N^j)) = 0$)

Finite element stiffness matrices are **sparse**

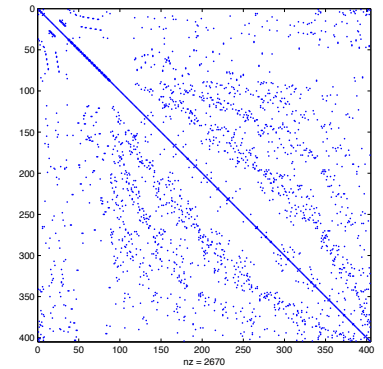
Definition 3.1.13 (Sparse matrix). A matrix $\mathbf{A} \in \mathbb{R}^{N,N}$ is called **sparse**, if $\text{nnz}(\mathbf{A}) := \#\{(i, j) : a_{ij} \neq 0\} \ll N^2$.

Example 35 (Sparse stiffness matrices).

V_N : one basis function associated with each vertex



Finite element mesh \mathcal{M}



Resulting **sparsity pattern** of stiffness matrix

Visualization of sparsity pattern: MATLAB-spy () -Funktion

Remark 36 (Storing sparse matrices).

- Special (efficient) storage formats for sparse matrices, e.g., **CRS-format**
- Special MATLAB commands: `sparse`, `spones`, `speye`, `spdiags`
(→ use **mandatory** !)

3.1.4 Linear H^1 -conforming finite elements

\mathcal{M} = simplicial mesh of polygonal/polyhedral computational domain $\Omega \subset \mathbb{R}^d, d = 2, 3$

Linear H^1 -conforming finite elements

= Simplest $H^1(\Omega)$ -conforming finite element space

= Simplest finite element scheme for scalar second order elliptic BVP on Ω

$$\mathcal{S}_1^0(\mathcal{M}) := \{v \in C^0(\bar{\Omega}) : v|_K \in \mathcal{P}_1(K) \forall K \in \mathcal{M}\} \subset H^1(\Omega)$$

Representation: $\mathcal{P}_1(K) := \{\mathbf{x} \mapsto \alpha + \beta \cdot \mathbf{x}, \mathbf{x} \in K, \alpha \in \mathbb{R}, \beta \in \mathbb{R}^d\}$.
(space of d -variate polynomials of total degree ≤ 1)
 $\dim \mathcal{P}_1(K) = d + 1$

Notation: $\mathcal{S}_1^0(\mathcal{M})$ — continuous functions, cf. $C^0(\Omega)$
— locally 1st degree polynomials, cf. \mathcal{P}_1

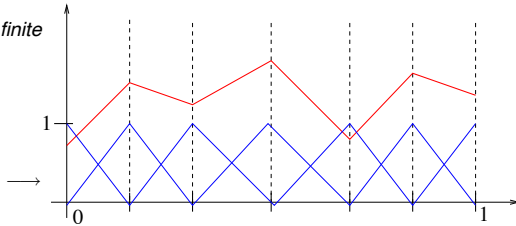
Example 37. ($H^1(\Omega)$ -conforming linear finite element space in 1D)

$d = 1, \Omega =]0, 1[$,

mesh \mathcal{M} = partition of $]0, 1[$ into intervals

red: function $\in \mathcal{S}_1^0(\mathcal{M})$

blue: hat function basis of $\mathcal{S}_1^0(\mathcal{M})$



Locally supported basis functions in 2D ?

On a triangle T with vertices $\mathbf{a}^1, \mathbf{a}^2, \mathbf{a}^3$: $q \in \mathcal{P}_1(T)$ uniquely determined by values $q(\mathbf{a}^i)$.

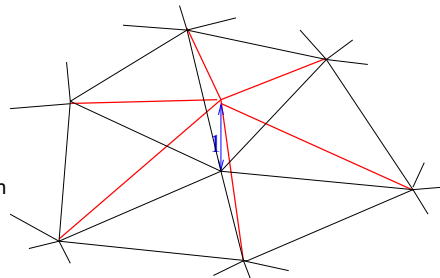
$v_N \in \mathcal{S}_1^0(\mathcal{M})$ uniquely determined by $\{v_N(\mathbf{x}), \mathbf{x} \text{ node of } \mathcal{M}\}$!

$\dim \mathcal{S}_1^0(\mathcal{M}) = \#\mathcal{N}(\mathcal{M})$ ($\mathcal{N}(\mathcal{M})$ = set of vertices of \mathcal{M})

If $\mathcal{N}(\mathcal{M}) = \{\mathbf{x}^1, \dots, \mathbf{x}^N\}$, nodal basis $\mathfrak{B}_N := \{b_N^1, \dots, b_N^N\}$ of $\mathcal{S}_1^0(\mathcal{M})$ defined by $b_N^i(\mathbf{x}^j) = \delta_{ij}$.

Piecewise linear nodal basis function
("hat function")
(= global shape function for $\mathcal{S}_1^0(\mathcal{M})$)

coefficient μ_j = "nodal value" of u_N at j -th node of \mathcal{M}



Global shape functions

Restriction to element

local shape functions

(3.1.7)

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Example 38 (Local shape functions for $\mathcal{S}_1^0(\mathcal{M})$).

Triangle K with vertices $\mathbf{a}^1 = \begin{pmatrix} 0 \\ 0 \end{pmatrix}, \mathbf{a}^2 = \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \mathbf{a}^3 = \begin{pmatrix} 0 \\ 1 \end{pmatrix}$:

Local shape functions: $b_K^1(\mathbf{x}) = 1 - x_1 - x_2, b_K^2(\mathbf{x}) = x_1, b_K^3(\mathbf{x}) = x_2$.

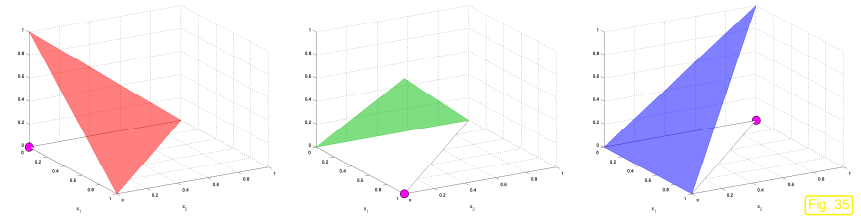


Fig. 35

Local shape functions for $\mathcal{S}_1^0(\mathcal{M})$ on triangle/tetrahedron = barycentric coordinate functions

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Definition 3.1.14 (Barycentric coordinates). Given $d+1$ points $\mathbf{a}^1, \dots, \mathbf{a}^{d+1} \in \mathbb{R}^d$ that do not lie in a hyperplane the **barycentric coordinates** $\lambda_1 = \lambda_1(\mathbf{x}), \dots, \lambda_{d+1} = \lambda_{d+1}(\mathbf{x}) \in \mathbb{R}$ of $\mathbf{x} \in \mathbb{R}^d$ are uniquely defined by

$$\lambda_1(\mathbf{x}) + \dots + \lambda_{d+1}(\mathbf{x}) = 1, \quad \lambda_1(\mathbf{x}) \mathbf{a}^1 + \dots + \lambda_{d+1}(\mathbf{x}) \mathbf{a}^{d+1} = \mathbf{x} \quad \forall \mathbf{x} \in \mathbb{R}^d.$$

Barycentric coordinates obtained by solving

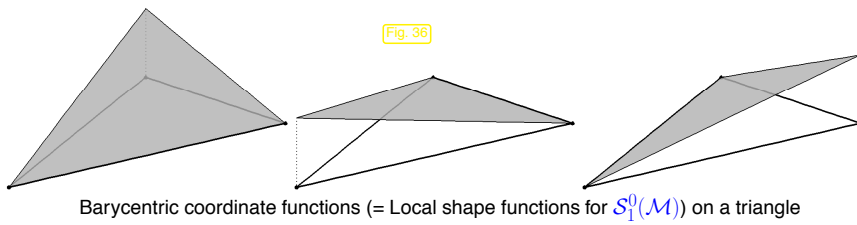
$$\begin{pmatrix} a_1^1 & \dots & a_1^{d+1} \\ \vdots & & \vdots \\ a_d^1 & \dots & a_d^{d+1} \\ 1 & \dots & 1 \end{pmatrix} \begin{pmatrix} \lambda_1(\mathbf{x}) \\ \vdots \\ \lambda_d(\mathbf{x}) \\ \lambda_{d+1}(\mathbf{x}) \end{pmatrix} = \begin{pmatrix} x_1 \\ \vdots \\ x_d \\ 1 \end{pmatrix}. \quad (3.1.8)$$

Corollary 3.1.15. Given $d+1$ points $\mathbf{a}^1, \dots, \mathbf{a}^{d+1} \in \mathbb{R}^d$ as in Def. 3.1.14, the barycentric coordinates are affine linear functions on \mathbb{R}^d , which satisfy

$$\lambda_j(\mathbf{a}^i) = \delta_{ij} := \begin{cases} 1, & \text{if } i = j, \\ 0, & \text{else,} \end{cases} \quad 1 \leq i, j \leq d+1.$$

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How to get $H_0^1(\Omega)$ -conforming finite element space $\mathcal{S}_{1,0}^0(\mathcal{M}) := \mathcal{S}_1^0(\mathcal{M}) \cap H_0^1(\Omega)$?

► Discard nodal basis functions associated with vertices on $\partial\Omega$!

Remark 39. Piecewise linear finite element subspace of $H_*^1(\Omega)$?

► There exist no locally supported piecewise linear basis functions.

3.1.5 Simplicial Lagrangian finite elements

\mathcal{M} = simplicial mesh of polygonal/polyhedral computational domain $\Omega \subset \mathbb{R}^d$, $d = 2, 3$

Idea: Use higher degree polynomials \rightarrow "better accuracy" (cf. interpolation)

Higher degree polynomials $\mathcal{P}_p(\mathbb{R}^d) := \{\mathbf{x} \in \mathbb{R}^d \mapsto \sum_{\alpha \in \mathbb{N}_0^d, |\alpha| \leq p} \kappa_\alpha \mathbf{x}^\alpha, \kappa_\alpha \in \mathbb{R}\}$.

Notation: α = "multiindex" $(\alpha_1, \dots, \alpha_d)$, $|\alpha| = \alpha_1 + \dots + \alpha_d$, $\mathbf{x}^\alpha := x_1^{\alpha_1} \dots x_d^{\alpha_d}$.

Example: $\mathcal{P}_2(\mathbb{R}^2) = \text{Span} \{1, x_1, x_2, x_1^2, x_2^2, x_1 x_2\}$

Lemma 3.1.16.

$$\dim \mathcal{P}_p(\mathbb{R}^d) = \binom{d+p}{p} \text{ for all } p \in \mathbb{N}, d \in \mathbb{N}$$

Definition 3.1.17 (Higher order Lagrangian finite element spaces). Space of p -th degree Lagrangian finite element functions on mesh \mathcal{M}

$$\mathcal{S}_p^0(\mathcal{M}) := \{v \in C^0(\bar{\Omega}) : v|_K \in \mathcal{P}_p(K) \quad \forall K \in \mathcal{M}\}.$$

Notation: $\mathcal{S}_p^0(\mathcal{M})$ continuous functions, cf. $C^0(\Omega)$
locally polynomials of degree p , cf. $\mathcal{P}_p(\mathbb{R}^d)$

Construction: Local shape functions $\xrightarrow{\text{"Glueing"}}$ Global FE space
(Glueing must ensure global continuity $\leftrightarrow H^1(\Omega)$ -conformity)

! Design of local shape functions must make glueing possible

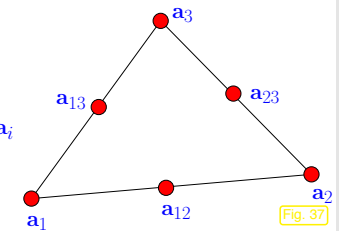
Example 40 (Quadratic Lagrangian finite elements).

Local shape functions for $\mathcal{P}_2(K)$, K triangle:

$$\begin{aligned} b_1^K &= -\lambda_1(1 - 2\lambda_1), & b_{12}^K &= 4\lambda_1\lambda_2, \\ b_2^K &= -\lambda_2(1 - 2\lambda_2), & b_{13}^K &= 4\lambda_1\lambda_3, \\ b_3^K &= -\lambda_3(1 - 2\lambda_3), & b_{23}^K &= 4\lambda_2\lambda_3. \end{aligned}$$

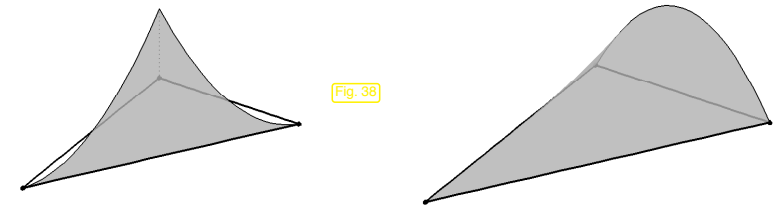
λ_i = barycentric coordinate function (\rightarrow Def. 3.1.14) for vertex \mathbf{a}_i

$$b_j^K(\mathbf{a}_i) = \delta_{ij}, \quad i, j \in \{1, 2, 3, (12), (23), (13)\}.$$



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► Local shape functions = Lagrangian (interpolatory) polynomials for local nodes in K

► Specifying local interpolation nodes \Leftrightarrow specifying local shape functions

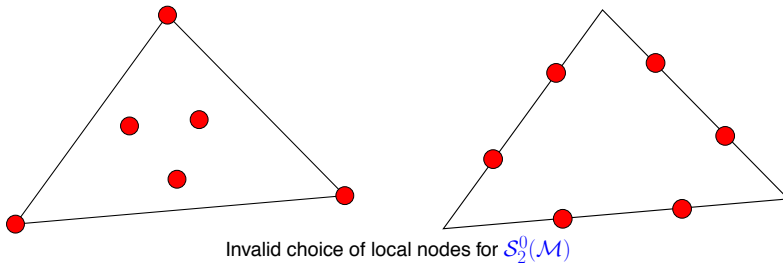
When are local nodes \mathbf{q}_i , $i = 1, \dots, Q$, (for $\mathcal{P}_p(K)$) suitable for "glueing" ?

• Unisolvence: $v \in \mathcal{P}_p(K): v(\mathbf{q}_i) = 0 \quad \forall i \Leftrightarrow v \equiv 0$

• Fixing traces: locally unisolvent interpolation on each vertex/edge/face

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• **Interelement matching:**

corresponding nodes on joint edges/faces

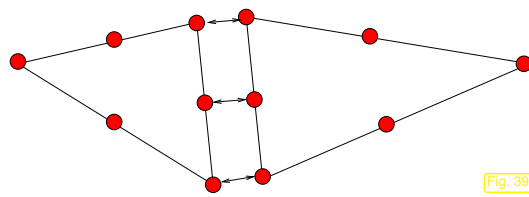


Fig. 39

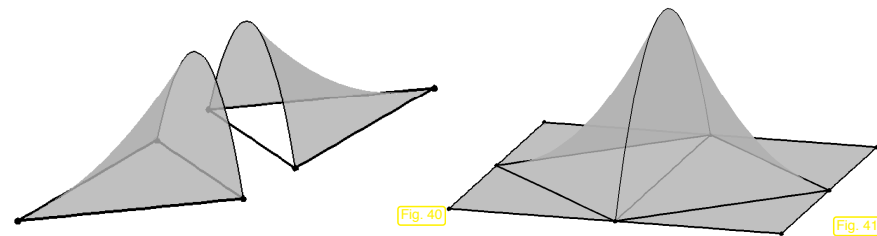
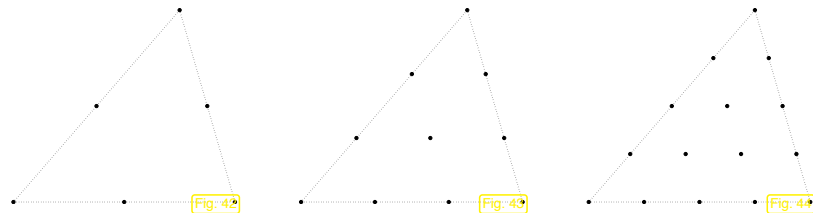
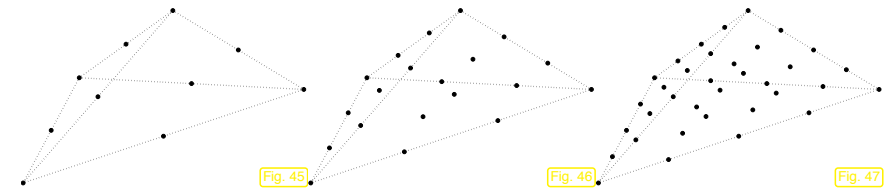


Fig. 40

Fig. 41



Location of local (interpolation) nodes for triangular Lagrangian finite elements of degree 2 (left), degree 3 (middle), and degree 4 (right)



Local nodes for tetrahedral Lagrangian finite elements (left: $p = 2$, middle: $p = 3$, right: $p = 4$)

Can we find other locally supported bases for $S_2^0(\mathcal{M})$?

YES!



Alternative **p-hierarchical** local shape functions:

$$\begin{aligned} b_1^K &= \lambda_1, & b_{12}^K &= 4\lambda_1\lambda_2, \\ b_2^K &= \lambda_2, & b_{13}^K &= 4\lambda_1\lambda_3, \\ b_3^K &= \lambda_3, & b_{23}^K &= 4\lambda_2\lambda_3. \end{aligned}$$

Set comprises local shape functions for $p = 1$.



Glueing can easily be accomplished

No “canonical” local shape functions/global basis functions for higher order Lagrangian finite elements.



Selection of \mathfrak{B}_N to get “good” matrix properties.

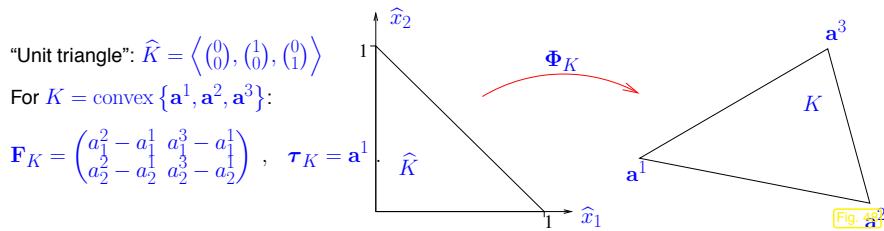
3.1.6 Parametric finite elements

Definition 3.1.18 (Affine transformation). Mapping $\Phi : \mathbb{R}^d \mapsto \mathbb{R}^d$ **affine**, if $\Phi(\mathbf{x}) = \mathbf{F}\mathbf{x} + \boldsymbol{\tau}$ with $\mathbf{F} \in \mathbb{R}^{d,d}$, $\boldsymbol{\tau} \in \mathbb{R}^d$.

Usually:

All elements of a mesh = **affine** images of **reference element(s)** \hat{K}

$$\exists \hat{K} \quad \forall K \in \mathcal{M}: \quad \exists \mathbf{F}_K \in \mathbb{R}^{d,d} \text{ regular, } \boldsymbol{\tau}_K \in \mathbb{R}^d: \quad K = \Phi_K(\hat{K}) \quad \text{with} \quad \Phi_K(\hat{\mathbf{x}}) := \mathbf{F}_K \hat{\mathbf{x}} + \boldsymbol{\tau}_K.$$



Transformations of elements \Rightarrow transformation of functions:

Definition 3.1.19 (Pullback). Given domains $\Omega, \hat{\Omega}$ and a bijective mapping $\Phi : \hat{\Omega} \mapsto \Omega$, the pullback $\Phi^*u : \hat{\Omega} \mapsto \mathbb{R}$ of a function $u : \Omega \mapsto \mathbb{R}$ is defined by

$$(\Phi^*u)(\hat{\mathbf{x}}) := u(\Phi(\hat{\mathbf{x}})), \quad \hat{\mathbf{x}} \in \hat{\Omega}.$$

Notation: If unambiguous: $\hat{u} := \Phi^*u$

Note: Consider $\mathcal{S}_1^0(\mathcal{M})$, triangle $K \in \mathcal{M}$, unit triangle \hat{K} , affine mapping $\Phi_K : \hat{K} \mapsto K$

- b_1^K, b_2^K, b_3^K (standard) local shape functions on K
- $\hat{b}_1, \hat{b}_2, \hat{b}_3$ (standard) local shape functions on \hat{K}

Observation: $\hat{b}_i = \Phi_K^* b_i^K$

Terminology: **affine equivalent finite elements**

► All families of Lagrangian finite elements (Sect. 3.1.5) (equipped with "natural" local shape functions) are affine equivalent.



STEP 1: **define** local shape functions on reference element \hat{K}

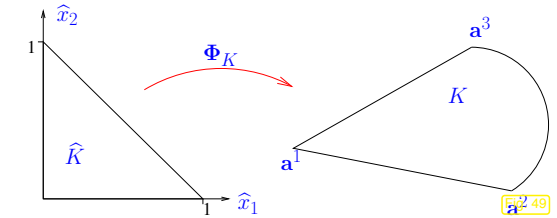
STEP 2: local shape functions on $K \in \mathcal{M}$ via pullback $(\Phi_K^{-1})^*$ (\rightarrow Def. 3.1.19)

Parametric finite elements

Generalization: **curvilinear** meshes with curved edges/faces (\rightarrow Sect. 3.2.8)

Now: Φ_K diffeomorphism!
(i.e. Φ_K and Φ_K^{-1} are differentiable)

$$b_i^K := (\Phi_K^{-1})^* \hat{b}_i.$$



► Application: approximation of curved interfaces/boundaries (\rightarrow Sect 3.2.8)

3.1.7 Lagrangian finite elements on quadrilaterals/hexahedra

Parametric construction: start from reference element $\hat{K} =]0, 1[^d$ (unit cube)

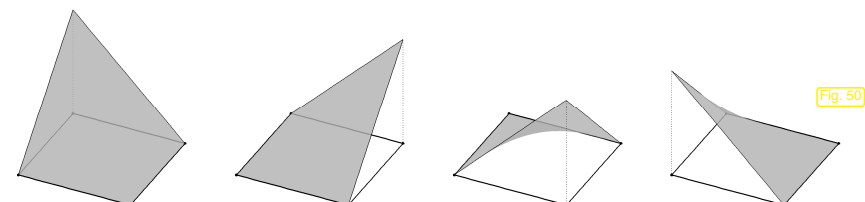
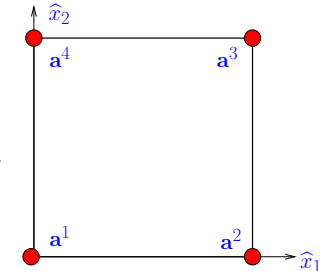
Lowest polynomial degree $p = 1$, 2D: piecewise **bilinear** finite elements

Local shape functions on $\hat{K} =]0, 1[^2$

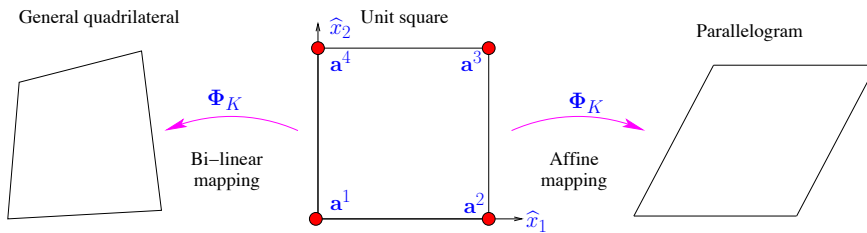
$$\begin{aligned} \hat{b}_1 &= (1 - \hat{x}_1)(1 - \hat{x}_2), & \hat{b}_2 &= \hat{x}_1(1 - \hat{x}_2), \\ \hat{b}_3 &= \hat{x}_1\hat{x}_2, & \hat{b}_4 &= (1 - \hat{x}_1)\hat{x}_2. \end{aligned}$$

Bilinear Lagrangian interpolation polynomials w.r.t. corner points \mathbf{a}_i of \hat{K}

Note: \hat{b}_i linear on edges



Bilinear local shape functions on unit square \hat{K}



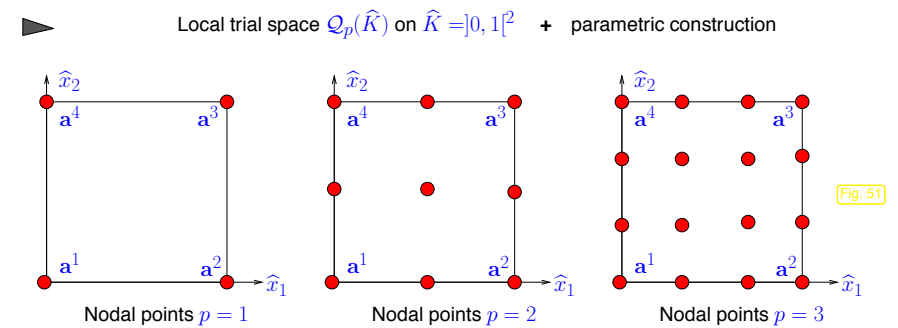
Affine mapping = linear transformation + translation

Bilinear mapping to general quadrilateral:

$$\Phi_K(\hat{\mathbf{x}}) = \begin{pmatrix} \alpha_1 + \beta_1 \hat{x}_1 + \gamma_1 \hat{x}_2 + \delta_1 \hat{x}_1 \hat{x}_2 \\ \alpha_2 + \beta_2 \hat{x}_1 + \gamma_2 \hat{x}_2 + \delta_2 \hat{x}_1 \hat{x}_2 \end{pmatrix}, \quad \alpha_i, \beta_i, \gamma_i, \delta_i \in \mathbb{R}.$$

$\hat{b}_i, i = 1, 2, 3, 4$, bilinear local shape functions on $]0, 1[^2$, $\Rightarrow b_i^K = (\Phi_K^{-1})^* \hat{b}_i$ linear on edges of K
 $\Phi_K : \hat{K} \mapsto K$ bilinear

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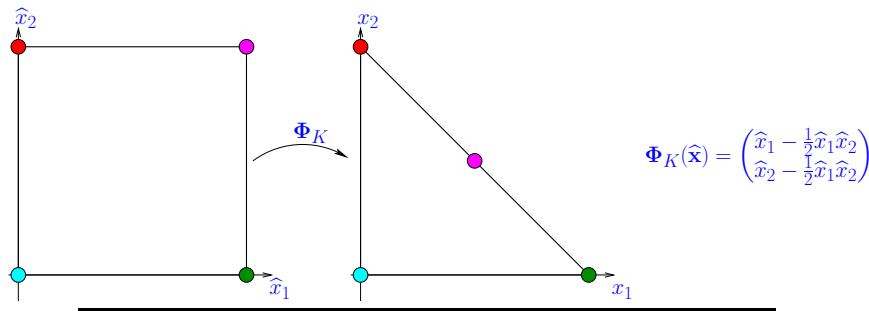


3.1.8 Degrees of freedom*

Recall: Lagrangian local shape functions b_i^K fixed by $b_i^K(\mathbf{q}_j) = \delta_{ij}$ for nodes $\mathbf{q}_i, i, j = 1, \dots, Q$, $Q \in \mathbb{N}$ (\rightarrow Sects. 3.1.5, 3.1.7).

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Example 41 (Triangle as degenerate quadrilateral).



Higher order quadrilateral Lagrangian finite elements:

Definition 3.1.20 (Tensor product polynomials). Space of *tensor product polynomials* of degree $p \in \mathbb{N}$ in each coordinate direction

$$\mathcal{Q}_p(\mathbb{R}^d) := \{\mathbf{x} \mapsto p_1(x_1) \cdots p_d(x_d), p_i \in \mathcal{P}_p(\mathbb{R}), i = 1, \dots, d\}.$$

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Definition 3.1.21 (Dual basis). Given a vector space V with basis $\mathfrak{B} := \{b_1, \dots, b_Q\}$, $Q \in \mathbb{N}$, the corresponding *dual basis* is a set l_1, \dots, l_Q of linear forms on V such that

$$l_j(b_i) = \delta_{ij}, \quad i, j \in \{1, \dots, Q\}.$$

For Lagrangian finite elements $\mathcal{S}_p^0(\mathcal{M})$, on element K :

Nodal evaluation functionals $v \mapsto v(\mathbf{q}_j)$, $j = 1, \dots, Q$, form dual basis w.r.t. basis $\{b_1^K, \dots, b_Q^K\}$ of local trial space $\mathcal{S}_p(K)$.

Definition 3.1.22 (Local degrees of freedom). A dual basis of the local trial space corresponding to the local shape functions provides *local degrees of freedom* (d.o.f.).

Role reversal: degrees of freedom \Rightarrow local shape functions

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Example 42 (Cubic Hermitian Finite Elements on triangular mesh).

- Local trial space $V_K := \mathcal{P}_3(K)$ for each $K \in \mathcal{M}$,

- Local degrees of freedom, K triangle with vertices $\mathbf{a}^1, \mathbf{a}^2, \mathbf{a}^3 \in \mathbb{R}^2$

$$\begin{aligned} l_1(v) &= v(\mathbf{a}^1), & l_2(v) &= v(\mathbf{a}^2), & l_3(v) &= v(\mathbf{a}^3), \\ l_4(v) &= \mathbf{grad} v(\mathbf{a}^1) \cdot (\mathbf{a}^2 - \mathbf{a}^1), & l_5(v) &= \mathbf{grad} v(\mathbf{a}^2) \cdot (\mathbf{a}^3 - \mathbf{a}^2), & l_6(v) &= \mathbf{grad} v(\mathbf{a}^3) \cdot (\mathbf{a}^1 - \mathbf{a}^3) \\ l_7(v) &= \mathbf{grad} v(\mathbf{a}^1) \cdot (\mathbf{a}^3 - \mathbf{a}^1), & l_8(v) &= \mathbf{grad} v(\mathbf{a}^2) \cdot (\mathbf{a}^1 - \mathbf{a}^2), & l_9(v) &= \mathbf{grad} v(\mathbf{a}^3) \cdot (\mathbf{a}^2 - \mathbf{a}^3) \\ l_{10}(v) &= v(\tfrac{1}{3}(\mathbf{a}^1 + \mathbf{a}^2 + \mathbf{a}^3)). \end{aligned}$$

Dual basis (\rightarrow Def. 3.1.21) ?

- # functionals = $\dim V_K = \dim \mathcal{P}_3(\mathbb{R}^2) = \binom{3+2}{2} = 10$,

- If $l_j(v) = 0$ for all $j = 1, \dots, 10$, then $\Rightarrow v(\mathbf{a}^i) = 0$ and $\mathbf{grad} v(\mathbf{a}^i) = 0, i = 1, 2, 3$,
 $\Rightarrow v \in \mathcal{P}_3(K) \equiv 0$ on any edge,
 $\Rightarrow (= 0 \text{ at center of gravity}) v \equiv 0$ for any $v \in V_K$.

► **Unisolvence** of local d.o.f.

Suitable for glueing ?

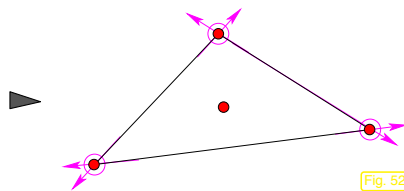
YES, because $v|_{\text{edge}}$ uniquely determined by d.o.f. associated with the edge.

A local degree of freedom l is regarded as **associated with an edge E** , if $l(v)$ only depends on $v|_E$, $\mathbf{grad} v|_E, \dots$

Symbolic notation for local d.o.f. for cubic

Hermitian elements:

(filled circle = nodal values, circle = first derivatives, arrows = directional derivatives)



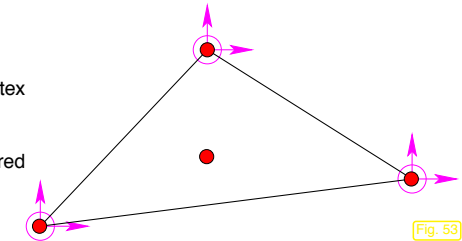
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HOWEVER, alternative choice of local degrees of freedom possible
(on triangle K with vertices $\mathbf{a}^1, \mathbf{a}^2, \mathbf{a}^3 \in \mathbb{R}^2$)

$$\begin{aligned} l_1(v) &= v(\mathbf{a}^1), & l_2(v) &= v(\mathbf{a}^2), & l_3(v) &= v(\mathbf{a}^3), \\ l_4(v) &= \frac{\partial v}{\partial x_1}(\mathbf{a}^1), & l_5(v) &= \frac{\partial v}{\partial x_1}(\mathbf{a}^2), & l_6(v) &= \frac{\partial v}{\partial x_1}(\mathbf{a}^3), \\ l_7(v) &= \frac{\partial v}{\partial x_2}(\mathbf{a}^1), & l_8(v) &= \frac{\partial v}{\partial x_2}(\mathbf{a}^2), & l_9(v) &= \frac{\partial v}{\partial x_2}(\mathbf{a}^3), \\ l_{10}(v) &= v(\mathbf{a}^{123}). \end{aligned}$$

Three d.o.f. associated with each vertex

Fewer global shape functions compared to previous choice!

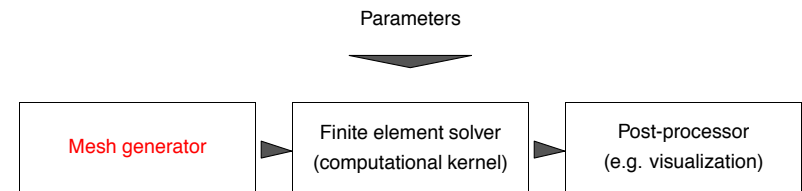


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3.2 Implementation

3.2.1 Mesh file format

Data flow in (most) finite element software packages:



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Example 43 (Mesh file format (triangular mesh of polygonal domain)).

```
# Two-dimensional simplicial mesh
1   $\xi_1$    $\eta_1$           # Coordinates of first node
2   $\xi_2$    $\eta_2$           # Coordinates of second node
:
N   $\xi_N$    $\eta_N$           # Coordinates of N-th node
1   $n_1^1$   $n_2^1$   $n_3^1$   $X_1$   # Indices of nodes of first triangle
2   $n_1^2$   $n_2^2$   $n_3^2$   $X_2$   # Indices of nodes of second triangle
:
M   $n_1^M$   $n_2^M$   $n_3^M$   $X_M$  # Indices of nodes of M-th triangle
 $X_i, i = 1, \dots, M \rightarrow$  extra information (e.g. material properties in triangle #i).
```

(3.2.1)

Optional: additional information about edges (on $\partial\Omega$):

```
 $K \in \mathbb{N}$           # Number of edges on  $\partial\Omega$ 
 $n_1^1$   $n_2^1$   $Y_1$     # Indices of endpoints of first edge
 $n_1^2$   $n_2^2$   $Y_2$     # Indices of endpoints of second edge
:
 $n_1^K$   $n_2^K$   $Y_K$     # Indices of endpoints of K-th edge
 $Y_k, k = 1, \dots, K \rightarrow$  extra information
```

(3.2.2)

Example 44 (Mesh file format for MATLAB code).

Vertex coordinate file:	Cell information file:
% List of vertices	% List of elements
1 +0.000000e+00 -1.000000e+00	1 1 2 5
2 +1.000000e+00 +0.000000e+00	2 2 3 5
3 +0.000000e+00 +1.000000e+00	3 3 4 5
4 -1.000000e+00 +0.000000e+00	4 4 1 5
5 +0.000000e+00 +0.000000e+00	

```
Loading a mesh
m = load_Mesh('Coord_Circ.dat',...
              'Elem_Circ.dat');
plot_Mesh(m, 'apts');
```

Option flags:

'a': with axes
'p': vertex labels on
't': cell labels on
's': caption/title on

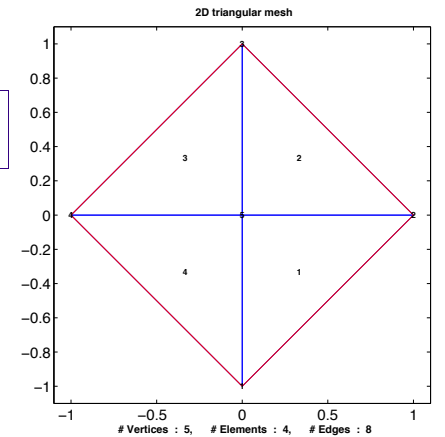


Fig. 54

How to create a mesh ?

→ Mesh generation (beyond scope of this course)

→ <http://www.andrew.cmu.edu/user/sowen/mesh.html>

Free software:

- NETGEN (<http://www.hpfem.jku.at/netgen/>)
- Triangle (<http://www.cs.cmu.edu/~quake/triangle.html>)
- TETGEN (<http://tetgen.berlios.de>)

Example 45 (Mesh generation in MATLAB code).

Algorithm & details → [?]

```
MATLAB-CODE: mesh generation for circular domain
BBOX = [-1 -1; 1 1];
H0 = 0.1;
DHD = @(x) sqrt(x(:,1).^2+x(:,2).^2)-1;
HHANDLE = @(x) ones(size(x,1),1);
Mesh = init_Mesh(BBOX,H0,DHD,...
                 HHANDLE,[],1);
save_Mesh(Mesh,'Coordinates.dat',...
           'Elements.dat');
```

Bounding box

Largest reasonable edge length

Signed distance function $\varphi(\mathbf{x})$:
(distance from $\partial\Omega$, $\varphi(\mathbf{x}) < 0 \Leftrightarrow \mathbf{x} \in \Omega$)

Element size function
(determines local edge length)

3.2.2 Assembly

→ term used for computing entries of stiffness matrix/load vector.

Discrete variational problem (V_N = FE space, $\dim V_N = N \in \mathbb{N}$, see Sect. 3.1.1)

$$u_N \in V_N: \quad a(u_N, v_N) = f(v_N) \quad \forall v_N \in V_N. \quad (3.1.1)$$

To be computed:

- Global stiffness matrix: $\mathbf{A} = \left(a(b_N^j, b_N^i) \right)_{i,j=1}^N \in \mathbb{R}^{N,N}$

- Global load vector: $\vec{\varphi} := \left(f(b_N^i) \right)_{i=1}^N \in \mathbb{R}^N$

both can be written in terms of **local cell contributions**:

$$a(u, v) = \sum_{K \in \mathcal{M}} a_K(u|_K, v|_K) \quad , \quad f(v) = \sum_{K \in \mathcal{M}} f_K(v|_K). \quad (3.2.3)$$

Example: bilinear form/linear form arising from 2nd-order elliptic BVPs (→ Sect. 2.5)

$$a(u, v) := \int_{\Omega} D \mathbf{grad} u \cdot \mathbf{grad} v \, dx = \sum_{K \in \mathcal{M}} \underbrace{\int_K D \mathbf{grad} u \cdot \mathbf{grad} v \, dx}_{=: a_K(u|_K, v|_K)},$$

$$f(v) := \int_{\Omega} f v \, dx = \sum_{K \in \mathcal{M}} \underbrace{\int_K f v \, dx}_{=: f_K(v|_K)}.$$

Recall (3.1.7): Restrictions of global shape functions to cells = local shape functions

Definition 3.2.1. Given local shape functions $\{b_1^K, \dots, b_Q^K\}$, we call

element stiffness matrix $\mathbf{A}_K := \left(a_K(b_j^K, b_i^K) \right)_{i,j=1}^Q \in \mathbb{R}^{Q,Q},$

element load vector $\vec{\varphi}_K := \left(f_K(b_i^K) \right)_{i=1}^Q \in \mathbb{R}^Q.$

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Theorem 3.2.2. The stiffness matrix and load vector can be obtained from their cell counterparts by

$$\mathbf{A} = \sum_K \mathbf{T}_K^T \mathbf{A}_K \mathbf{T}_K \quad , \quad \vec{\varphi} = \sum_K \mathbf{T}_K^T \vec{\varphi}_K, \quad (3.2.4)$$

with the **index mapping matrices** ("T-matrices") $\mathbf{T}_K \in \mathbb{R}^{Q,N}$, defined by

$$(\mathbf{T}_K)_{ij} := \begin{cases} 1 & , \text{ if } (b_N^j)|_K = b_i^K, \\ 0 & , \text{ otherwise.} \end{cases} \quad 1 \leq i \leq Q, 1 \leq j \leq N.$$

Proof.

$$(\mathbf{A})_{ij} = a(b_N^j, b_N^i) = \sum_{K \in \mathcal{M}} a_K(b_N^j|_K, b_N^i|_K) = \sum_{\substack{K \in \mathcal{M}, \text{supp}(b_N^j) \cap K \neq \emptyset, \\ \text{supp}(b_N^i) \cap K \neq \emptyset}} a_K(b_{l(j)}^K, b_{l(i)}^K) = \sum_{\substack{K \in \mathcal{M}, \text{supp}(b_N^j) \cap K \neq \emptyset, \\ \text{supp}(b_N^i) \cap K \neq \emptyset}} (\mathbf{A}_K)_{l(i), l(j)}$$

$l(i) \in \{1, \dots, k_K\}$, $1 \leq i \leq N \triangleq$ index of the local shape function corresponding to the global shape function b_N^i on K .

$$\Rightarrow (\mathbf{A})_{ij} = \sum_{\substack{K \in \mathcal{M}, \text{supp}(b_N^j) \cap K \neq \emptyset, \\ \text{supp}(b_N^i) \cap K \neq \emptyset}} \sum_{l=1}^Q \sum_{n=1}^Q (\mathbf{T}_K)_{li} (\mathbf{A}_K)_{ln} (\mathbf{T}_K)_{nj}. \quad \square$$

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Example 46 (Assembly for linear Lagrangian finite elements on triangular mesh).

Using the local/global numbering indicated beside

$$\rightarrow \mathbf{T}_{K^*} = \begin{pmatrix} 0 & 1 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 & 0 & 0 & 1 & 0 & 0 \end{pmatrix}$$

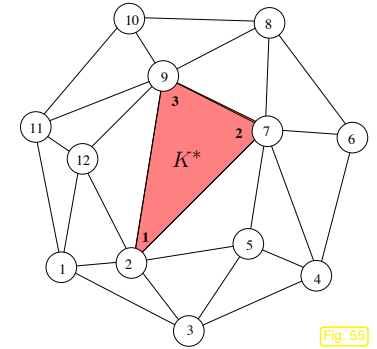


Fig. 55

Definition 3.2.3 (Abstract assembly operator). We denote the assembly operator in (3.2.4) symbolically by

$$\mathbf{A} = \mathcal{A}_{K \in \mathcal{M}} \mathbf{A}_K, \quad \vec{\varphi} = \mathcal{A}_{K \in \mathcal{M}} \vec{\varphi}_K.$$

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3.2.3 Mesh data structures

Issue: internal representation of mesh (\rightarrow Def. 3.1.10) in computer code

mesh data structure must provide:

1. offer unique identification of cells/(faces)/edges/vertices
2. represent **mesh topology** (= incidence relationships of cells/faces/edges/vertices)
3. describe **mesh geometry** (= location/shape of cells/faces/edges/vertices)
4. allow sequential access to edges/faces of a cell
(\rightarrow traversal of local shape functions/degrees of freedom)
5. make possible traversal of cells of the mesh (\rightarrow **global numbering**)

Focus: **array oriented data layout** (\rightarrow MATLAB, FORTRAN)

Notation:

\mathcal{M} = mesh (set of elements), $\mathcal{N}(\mathcal{M})$ = set of nodes (vertices) in \mathcal{M} , $\mathcal{E}(\mathcal{M})$ = set of edges in \mathcal{M}

Case: d -dimensional simplicial triangulation \mathcal{M} , *minimal* data structure (cf. Sect. 3.2.1)

\rightarrow Coordinates of vertices $\mathcal{N}(\mathcal{M})$: $\# \mathcal{N}(\mathcal{M}) \times d$ -array Coordinates of reals

\rightarrow Vertex indices for cells: $\# \mathcal{M} \times (d+1)$ -array Elements of integers.

► Already offers complete description of the mesh topology and geometry !

Optional extra information:

\rightarrow Edge connecting vertices: $\# \mathcal{N}(\mathcal{M}) \times \# \mathcal{N}(\mathcal{M})$ symmetric sparse integer matrix $I_{\mathcal{E}}$

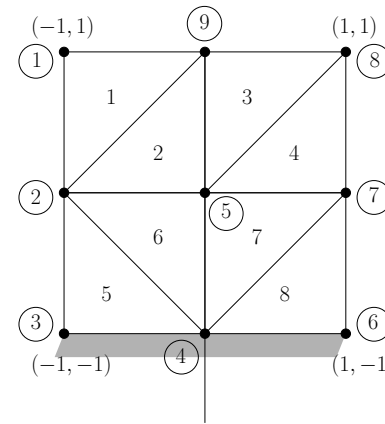
$$(I_{\mathcal{E}})_{ij} := \begin{cases} 0 & , \text{ if vertex } \#i \text{ not linked to } \#j \\ e_{ij} & , \text{ if edge connecting } \#i \text{ and } \#j \end{cases}$$

here e_{ij} is the unique edge number $\in \{1, 2, \dots, \# \mathcal{E}(\mathcal{M})\}$

\rightarrow End points of the edges: $\# \mathcal{E}(\mathcal{M}) \times 2$ array of integer (= vertex indices of end points).

\rightarrow Cell adjacent to edges: $\# \mathcal{E}(\mathcal{M}) \times 2$ array of integers (=cell indices)
(one cell index =0 if edge is on $\partial \Omega$)

Note: Global shape functions associated with edges/faces \rightarrow extra information required !



(Fig. 56)

Global shape functions associated with edges/faces \rightarrow extra information required !

Example 48 (Extended MATLAB mesh data structure).

```
mesh = add_Edge2Elem(add_Edges(init_Mesh(BBOX,H0,DHD,HHANDLE,[ ],1)))
```

(init_Mesh \rightarrow Ex. 45)

mesh =
Coordinates: [5x2 double] vertex coordinates, see Ex. 44
Elements: [4x3 double] vertex indices of triangles, see Ex. 44
Edges: [8x2 double] indices of endpoints in Coordinates array
Vert2Edge: [5x5 double] $\# \mathcal{N}(\mathcal{M}) \times \# \mathcal{N}(\mathcal{M})$ sparse integer matrix:
Edge2Elem: [8x2 double] entry (i, j) = edge index, if $\neq 0$
EdgeLoc: [8x2 double] $\# \mathcal{E}(\mathcal{M}) \times 2$ integer array:
indices of adjacent cells in Elements array
 $\# \mathcal{E}(\mathcal{M}) \times 2$ integer array: **local** indices of edges w.r.t. adjacent cells

Notation: $\mathcal{E}(\mathcal{M}) \doteq$ edges of 2D mesh

How to number \leftrightarrow order local shape functions
global shape functions ?

Elements, Edges arrays \rightarrow ordering of vertices of cells/endpoints of edges

Arrays (of vertices, cells, edges) \rightarrow array indices \rightarrow numbering of global shape functions

3.2.4 Algorithms

Cell oriented assembly \leftrightarrow (3.2.4) $\leftrightarrow \mathbf{A} = \mathbf{A}_{K \in \mathcal{M}} \mathbf{A}_K$

$$\mathbf{A} = \mathbf{A}_{K \in \mathcal{M}} \mathbf{A}_K := \left\{ \begin{array}{l} \text{foreach } K \in \mathcal{M} \text{ do} \\ \quad \text{local operations on } K \rightarrow \mathbf{A}_K \text{ and } \mathbf{A} = \mathbf{A} + \mathbf{T}_K^T \mathbf{A}_K \mathbf{T}_K \\ \text{enddo} \end{array} \right\}$$

Notion: **local operations** $\hat{=}$ \bullet required only data from fixed “neighbourhood” of K
 \bullet computational effort “ $O(1)$ ”: independent of $\#\mathcal{M}$

Cell oriented assembly in MATLAB

```
function A = assemble(Mesh)
for k = Mesh.Elements'
    idx = ❶
    Aloc = ❷
    A(idx,idx) = A(idx,idx)+Aloc;
end
```

❶ row vector of index numbers of global shape functions $b_{i_1}, \dots, b_{i_Q} \in V_N$ corresponding to local shape functions b_1^K, \dots, b_Q^K :
 \blacktriangleright $\text{idx} = (i_1, \dots, i_Q)$ (encodes index mapping matrix \mathbf{T}_K)
❷ $Q \times Q$ element stiffness matrix

For Lagrangian FEM (\rightarrow Sect. 3.1.5):

the total computational effort is of the order $O(\#\mathcal{M}) = O(N)$, $N := \dim V_N$.

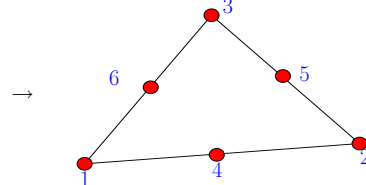
Example 49 (Assembly for quadratic Lagrangian FE in MATLAB code).

Setting: FE space $\mathcal{S}_2^0(\mathcal{M})$ on triangular mesh \mathcal{M} of polygon $\Omega \subset \mathbb{R}^2$

Recall: 6 local shape functions: 3 vertex-associated, 3 edge-associated \rightarrow Ex. 40, Sect. 3.1.5

Convention: vertex-associated global shape functions $\rightarrow b_1, \dots, b_{\#\mathcal{M}}$
edge-associated global shape functions $\rightarrow b_{\#\mathcal{M}+1}, \dots, b_{\#\mathcal{M}+\#\mathcal{E}(\mathcal{M})}$

Local numbering



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MATLAB-CODE: assembly for quadratic Lagrangian FE

```
function A = assemMat_QFE(Mesh,EHandle,varargin)

nV = size(Mesh.Coordinates,1);
nE = size(Mesh.Elements,1)

I = zeros(36*nE,1); J = I; a = I; offset = 0;
for k = 1:nE
    vidx = Mesh.Elements(k,:);
    idx = [vidx,...
           Mesh.Vert2Edge(vidx(1),vidx(2))+nV,...
           Mesh.Vert2Edge(vidx(2),vidx(3))+nV,...
           Mesh.Vert2Edge(vidx(3),vidx(1))+nV];
    Aloc = transpose(EHandle(Mesh.Coordinates(vidx,:),...
                             Mesh.ElemFlag(k),varargin{:}));

    Qsq = prod(size(Aloc)); range = offset + 1:Qsq;
    t = idx(ones(length(idx),1),:); I(range) = t(:);
    t = idx(ones(1,length(idx)),:); J(range) = t(:);
    a(range) = Aloc(:);
    offset = offset + Qsq;
end
A = sparse(I,J,a);
```

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❶: EHandle (function handle) \rightarrow provides element stiffness matrix $\mathbf{A}_K \in \mathbb{R}^{6,6}$

❷: $\mathbf{I}, \mathbf{J}, \mathbf{a} \hat{=}$ linear arrays storing (i, j, a_{ij}) for stiffness matrix \mathbf{A} .
 Initialized with 0 for the sake of efficiency \rightarrow Ex. 50

❸: $\text{idx} \hat{=}$ index mapping vector, see ❶ above

❹: $\mathbf{Aloc} = \mathbf{A}_K \in \mathbb{R}^{6,6}$ (element stiffness matrix)

❺: Mesh.ElemFlag(k) marks groups of elements (e.g. to select local heat conductivity D in (2.5.4))

❻: Build *sparse* MATLAB-matrix (\rightarrow Def. 3.1.13) from index-entry arrays

◇

Example 50 (Efficient implementation of assembly).

tic-toc-timing (min of 4v runs), MATLAB V7, Intel Pentium 4 Mobile CPU 1.80GHz, Linux
 Computation of element stiffness matrices skipped !

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- **Sparse assembly:**

```

A(idx,idx) = A(idx,idx) + Aloc;
• Array assembly I: "growing arrays"
I = []; J = []; a = [];
...
t = idx(: , ones(length(idx),1))';
I = [I;t(:)];
t = idx(: , ones(1,length(idx)));
J = [J;t(:)];
a = [a; Aloc(:)];

```

- **Array assembly III**

→ see code fragment above

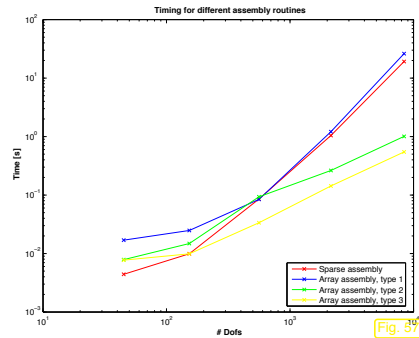
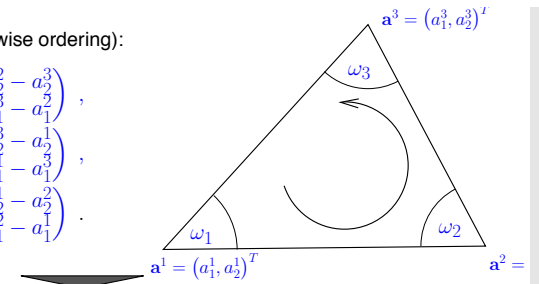


Fig. 3.9

If $\mathbf{a}^1, \mathbf{a}^2, \mathbf{a}^3$ vertices of K (counterclockwise ordering):

$$\begin{aligned}\lambda_1(\mathbf{x}) &= \frac{1}{2|K|} \left(\mathbf{x} - \begin{pmatrix} a_1^2 \\ a_2^2 \end{pmatrix} \right) \cdot \begin{pmatrix} a_2^2 - a_2^3 \\ a_1^3 - a_1^2 \end{pmatrix}, \\ \lambda_2(\mathbf{x}) &= \frac{1}{2|K|} \left(\mathbf{x} - \begin{pmatrix} a_1^3 \\ a_2^3 \end{pmatrix} \right) \cdot \begin{pmatrix} a_2^3 - a_2^1 \\ a_1^1 - a_1^3 \end{pmatrix}, \\ \lambda_3(\mathbf{x}) &= \frac{1}{2|K|} \left(\mathbf{x} - \begin{pmatrix} a_1^1 \\ a_2^1 \end{pmatrix} \right) \cdot \begin{pmatrix} a_2^1 - a_2^2 \\ a_1^2 - a_1^1 \end{pmatrix}.\end{aligned}$$



$$\mathbf{grad} \lambda_1 = \frac{1}{2|K|} \begin{pmatrix} a_2^2 - a_2^3 \\ a_1^3 - a_1^2 \end{pmatrix}, \quad \mathbf{grad} \lambda_2 = \frac{1}{2|K|} \begin{pmatrix} a_2^3 - a_2^1 \\ a_1^1 - a_1^3 \end{pmatrix}, \quad \mathbf{grad} \lambda_3 = \frac{1}{2|K|} \begin{pmatrix} a_2^1 - a_2^2 \\ a_1^2 - a_1^1 \end{pmatrix}. \quad (3.2.8)$$

3.2.5 Local computations

First option:

analytic evaluations

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We discuss bilinear form related to $-\Delta$, triangular Lagrangian finite elements of degree p :

K triangle: $a_K(u, v) := \int_K \mathbf{grad} u \cdot \mathbf{grad} v \, d\mathbf{x}$ ► element stiffness matrix.

Use **barycentric coordinate representations** of local shape functions

$$b_i^K = \sum_{\alpha \in \mathbb{N}_0^3, |\alpha| \leq p} \kappa_\alpha \lambda_1^{\alpha_1} \lambda_2^{\alpha_2} \lambda_3^{\alpha_3}, \quad \kappa_\alpha \in \mathbb{R}, \quad (3.2.5)$$

$$\Rightarrow \mathbf{grad} b_i^K = \sum_{\alpha \in \mathbb{N}_0^3, |\alpha| \leq p} \kappa_\alpha \left(\alpha_1 \lambda_1^{\alpha_1-1} \lambda_2^{\alpha_2} \lambda_3^{\alpha_3} \mathbf{grad} \lambda_1 + \alpha_2 \lambda_1^{\alpha_1} \lambda_2^{\alpha_2-1} \lambda_3^{\alpha_3} \mathbf{grad} \lambda_2 + \alpha_3 \lambda_1^{\alpha_1} \lambda_2^{\alpha_2} \lambda_3^{\alpha_3-1} \mathbf{grad} \lambda_3 \right). \quad (3.2.6)$$

$$\text{to evaluate } \int_K \lambda_1^{\beta_1} \lambda_2^{\beta_2} \lambda_3^{\beta_3} \mathbf{grad} \lambda_i \cdot \mathbf{grad} \lambda_j \, d\mathbf{x}, \quad i, j \in \{1, 2, 3\}, \beta_k \in \mathbb{N}. \quad (3.2.7)$$

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$$\begin{aligned} \left(\int_K \mathbf{grad} \lambda_i \cdot \mathbf{grad} \lambda_j \, d\mathbf{x} \right)_{i,j=1}^3 &= \\ &= \frac{1}{2} \begin{pmatrix} \cot \omega_3 + \cot \omega_2 & -\cot \omega_3 & -\cot \omega_2 \\ -\cot \omega_3 & \cot \omega_3 + \cot \omega_1 & -\cot \omega_1 \\ -\cot \omega_2 & -\cot \omega_1 & \cot \omega_2 + \cot \omega_1 \end{pmatrix}. \quad (3.2.9) \end{aligned}$$

→ Exercise.

Lemma 3.2.4 (Integration of powers of barycentric coordinate functions). *For any non-degenerate d -simplex K and $\alpha_j \in \mathbb{N}$, $j = 1, \dots, d+1$,*

$$\int_K \lambda_1^{\alpha_1} \dots \lambda_{d+1}^{\alpha_{d+1}} \, d\mathbf{x} = d!|K| \frac{\alpha_1! \alpha_2! \dots \alpha_{d+1}!}{(\alpha_1 + \alpha_2 + \dots + \alpha_{d+1} + d)!} \quad \forall \alpha \in \mathbb{N}_0^{d+1}. \quad (3.2.10)$$

Remark. Alternative: **symbolic computing** (MAPLE, Mathematica) for local computations

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3.2.6 Numerical quadrature

Second option (for local evaluations): **Numerical quadrature**

$$\int_{\Omega} f(\mathbf{x}) d\mathbf{x} \approx \sum_{K \in \mathcal{M}} |K| \sum_{l=1}^{P_K} \omega_l^K f(\pi_l^K), \quad \pi_l^K \in K, \omega_l^K \in \mathbb{R}. \quad (3.2.11)$$

Terminology:

$$\omega_l^K \rightarrow \text{weights}, \quad \pi_l^K \rightarrow \text{quadrature nodes} \\ (3.2.11) = \text{local quadrature rule}$$

- Mandatory • for computation of load vector (f complicated/only available in procedural form)
• for computation of stiffness matrix, if $D = D(\mathbf{x})$ does not permit analytic integration.

Guideline: only quadrature rules with positive weights are numerically stable.

For affine equivalent finite elements (\rightarrow Sect. 3.1.6):

► Parametric definition of local quadrature rules on reference cell \hat{K} :

$$\int_{\hat{K}} f(\hat{\mathbf{x}}) d\hat{\mathbf{x}} \approx |\hat{K}| \sum_{l=1}^P \hat{\omega}_l f(\hat{\pi}_l) \quad \blacktriangleright \quad \int_{\Omega} f(\mathbf{x}) d\mathbf{x} \approx \sum_{K \in \mathcal{M}} |K| \sum_{l=1}^P \omega_l^K f(\pi_l^K) \\ \text{with } \omega_l^K = \hat{\omega}_l, \pi_l^K = \Phi_K(\hat{\pi}_l).$$

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How to gauge the quality of parametric local quadrature rules ?

Quality of a parametric local quadrature rule on $K \sim$ largest space of polynomials on \hat{K} integrated exactly by the corresponding quadrature rule on \hat{K} .

Parlance: Quadrature rule exact for $\mathcal{P}_p(\hat{K}) \Rightarrow$ quadrature rule of order $p+1$
degree of exactness p

Example 51 (Local quadrature rules on triangles).

If K triangle $\Rightarrow \hat{K} := \text{convex} \left\{ \begin{pmatrix} 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \begin{pmatrix} 0 \\ 1 \end{pmatrix} \right\}$.

Quadrature rules described by pairs $(\hat{\omega}_1, \hat{\pi}_1), \dots, (\hat{\omega}_P, \hat{\pi}_P)$, $P \in \mathbb{N}$.

- Quadrature rule of order 2 (exact for $\mathcal{P}_1(\hat{K})$)

$$\left\{ \left(\frac{1}{3}, \begin{pmatrix} 0 \\ 0 \end{pmatrix} \right), \left(\frac{1}{3}, \begin{pmatrix} 0 \\ 1 \end{pmatrix} \right), \left(\frac{1}{3}, \begin{pmatrix} 1 \\ 0 \end{pmatrix} \right) \right\}. \quad (3.2.12)$$

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- Quadrature rule of order 3 (exact for $\mathcal{P}_2(\hat{K})$)

$$\left\{ \left(\frac{1}{3}, \begin{pmatrix} 1/2 \\ 0 \end{pmatrix} \right), \left(\frac{1}{3}, \begin{pmatrix} 0 \\ 1/2 \end{pmatrix} \right), \left(\frac{1}{3}, \begin{pmatrix} 1/2 \\ 1/2 \end{pmatrix} \right) \right\}. \quad (3.2.13)$$

- One-point quadrature rule of order 2 (exact for $\mathcal{P}_1(\hat{K})$)

$$\left\{ \left(1, \begin{pmatrix} 1/3 \\ 1/3 \end{pmatrix} \right) \right\}. \quad (3.2.14)$$

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- Quadrature rule of order 6 (exact for $\mathcal{P}_5(\hat{K})$)

$$\left\{ \left(\frac{9}{40}, \begin{pmatrix} 1/3 \\ 1/3 \end{pmatrix} \right), \left(\frac{155 + \sqrt{15}}{1200}, \begin{pmatrix} 6 + \sqrt{15}/21 \\ 6 + \sqrt{15}/21 \end{pmatrix} \right), \left(\frac{155 + \sqrt{15}}{1200}, \begin{pmatrix} 9 - 2\sqrt{15}/21 \\ 6 + \sqrt{15}/21 \end{pmatrix} \right), \right. \\ \left. \left(\frac{155 + \sqrt{15}}{1200}, \begin{pmatrix} 6 + \sqrt{15}/21 \\ 9 - 2\sqrt{15}/21 \end{pmatrix} \right), \left(\frac{155 - \sqrt{15}}{1200}, \begin{pmatrix} 6 - \sqrt{15}/21 \\ 9 + 2\sqrt{15}/21 \end{pmatrix} \right), \right. \\ \left. \left(\frac{155 - \sqrt{15}}{1200}, \begin{pmatrix} 9 + 2\sqrt{15}/21 \\ 6 - \sqrt{15}/21 \end{pmatrix} \right), \left(\frac{155 - \sqrt{15}}{1200}, \begin{pmatrix} 6 - \sqrt{15}/21 \\ 6 - \sqrt{15}/21 \end{pmatrix} \right) \right\} \quad (3.2.15)$$

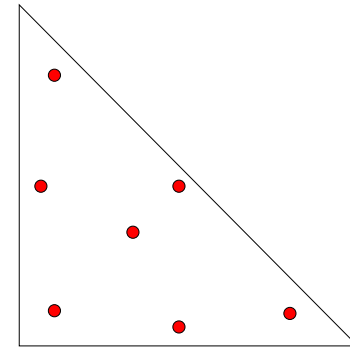


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In [?]: quadrature rules up to order $p = 21$ with $P \leq 1/6p(p+1) + 5$

Example 52 (Local quadrature rules on quadrilaterals).

If K quadrilateral $\Rightarrow \hat{K} := \text{convex} \left\{ \begin{pmatrix} 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 1 \\ 0 \end{pmatrix}, \begin{pmatrix} 0 \\ 1 \end{pmatrix}, \begin{pmatrix} 1 \\ 1 \end{pmatrix} \right\}$ (unit square).

On \hat{K} : **tensor product construction:**

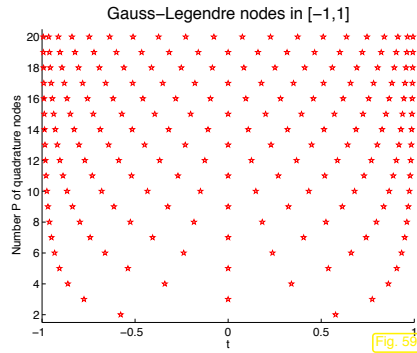
If $\{(\omega_1, \pi_1), \dots, (\omega_P, \pi_P)\}$, $P \in \mathbb{N}$, quadrature rule on the interval $]0, 1[$, exact for $\mathcal{P}_P]0, 1[$, then

$$\left\{ \begin{array}{ccc} (\omega_1^2, \begin{pmatrix} \pi_1 \\ \pi_1 \end{pmatrix}) & \cdots & (\omega_1 \omega_P, \begin{pmatrix} \pi_1 \\ \pi_P \end{pmatrix}) \\ \vdots & & \vdots \\ (\omega_1 \omega_P, \begin{pmatrix} \pi_P \\ \pi_1 \end{pmatrix}) & \cdots & (\omega_P^2, \begin{pmatrix} \pi_P \\ \pi_P \end{pmatrix}) \end{array} \right\}$$

quadrature rule on \hat{K} , exact for $\mathcal{Q}_P(\hat{K})$.

Quadrature rules on $]0, 1[$ (\rightarrow basic numerics):

- classical **Newton-Cotes formulas** (equidistant quadrature nodes).
- **Gauss-Legendre quadrature rules**, exact for $\mathcal{P}_{2P}]0, 1[$ using only P nodes.
- **Gauss-Lobatto quadrature rules**: P nodes including $\{0, 1\}$, exact for $\mathcal{P}_{2P-1}]0, 1[$.



3.2.7 Treatment of essential boundary conditions

Remember Sect. 2.7: extension $g \rightarrow \tilde{g}$ of Dirichlet data into Ω yielded linear variational problem.

Adaptation to finite element setting:

V_N = finite element space **without** constraints on $\partial\Omega$.

FIRST STEP:

Interpolation/projection of boundary data

FE-space $V_N \Rightarrow W_N := V_N|_{\partial\Omega}$ (FE trace space)

Example: if $V_N = \mathcal{S}_1^0(\mathcal{M})$, then W_N = set of piecewise linear, continuous functions on **boundary mesh** $\mathcal{M}_{|\partial\Omega}$.

BUT,

not necessarily $g \in W_N$!

► Replace g by (interpolant, least squares fit, etc.) $g_N \in W_N$

Example: if $V_N = \mathcal{S}_1^0(\mathcal{M})$ and $g \in C^0(\partial\Omega)$, then choose g_N as p.w. linear interpolant.

SECOND STEP:

Trivial extension of $g_N \rightarrow \tilde{g}_N \in V_N$

► Only nodal basis functions associated with node/edge/face $\subset \partial\Omega$ contribute to \tilde{g}_N !

Example: if $V_N = \mathcal{S}_1^0(\mathcal{M})$, g_N p.w. linear continuous on $\mathcal{M}_{|\partial\Omega}$

$$\tilde{g}_N = \sum_{\mathbf{p} \in \mathcal{N}(\mathcal{M}_{|\partial\Omega})} g_N(\mathbf{p}) b_N^{\mathbf{p}}, \quad \text{where } b_N^{\mathbf{p}} = \text{"hat function" for node } \mathbf{p}.$$

$$u_N \in V_{N,0}: a(u_N + \tilde{g}_N, v_N) = f(v_N) \quad \forall v_N \in V_{N,0}. \quad (3.2.16)$$

$V_{N,0} := \{v_N \in V_N: v_N = 0 \text{ on } \partial\Omega\}$ = span of "interior" basis functions.

$$\text{Elimination} \quad \blacktriangleright \quad \mathbf{A}_{\Omega\Omega} \tilde{\boldsymbol{\mu}}_\Omega = \tilde{\boldsymbol{\varphi}}_\Omega - \mathbf{A}_{\Gamma\Omega} \tilde{\boldsymbol{\mu}}_\Gamma. \quad (\text{cf. Sect. 2.7}) \quad (3.2.17)$$

Remark 53. Alternative: elimination on element level \Rightarrow modified $\tilde{\boldsymbol{\varphi}}_K$

$$\mathbf{A}_K = \begin{pmatrix} \mathbf{A}_{ii} & \mathbf{A}_{bi} \\ \mathbf{A}_{ib} & \mathbf{A}_{bb} \end{pmatrix}, \quad \tilde{\boldsymbol{\varphi}}_K = \begin{pmatrix} \tilde{\varphi}_i \\ \tilde{\varphi}_b \end{pmatrix} \quad \blacktriangleright \quad \tilde{\mathbf{A}}_K = \mathbf{A}_{ii}, \quad \tilde{\tilde{\boldsymbol{\varphi}}}_K = \tilde{\varphi}_i - \mathbf{A}_{bi} \tilde{\boldsymbol{\mu}}_{\Gamma,K}.$$

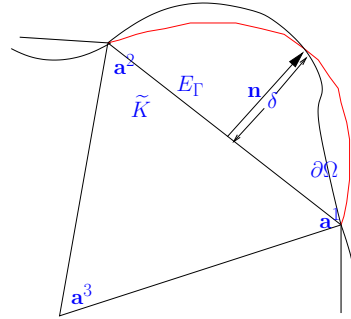
Then do assembly based on $\tilde{\mathbf{A}}_K$ and $\tilde{\tilde{\boldsymbol{\varphi}}}_K$.

3.2.8 Boundary approximation

Sect 3.1.6 → approximate treatment of curved $\partial\Omega$ by parametric FE:

Idea: **Piecewise polynomial approximation** of boundary (boundary fitting)
($\partial\Omega$ locally considered as function over straight edge of an element)

Example: Piecewise quadratic boundary approximation
(Part of $\partial\Omega$ between \mathbf{a}^1 and \mathbf{a}^2 approximated by parabola)



Mapping $\tilde{K} \rightarrow$ “curved element” K :

$$\Phi(\tilde{\mathbf{x}}) := \tilde{\mathbf{x}} + 4\delta \lambda_1(\tilde{\mathbf{x}})\lambda_2(\tilde{\mathbf{x}}) \mathbf{n}.$$

(λ_i barycentric coordinate functions on \tilde{K} , \mathbf{n} normal to E_Γ)

Note: Essential: Φ diffeomorphism $\leftrightarrow \delta$ sufficiently small

Transformation formula for gradients: for $u : K \mapsto \mathbb{R}$, diffeomorphism $\Phi : \tilde{K} \mapsto K$

$$(\mathbf{grad}_{\tilde{\mathbf{x}}}(\Phi^*u))(\tilde{\mathbf{x}}) = (D\Phi(\tilde{\mathbf{x}}))^T (\mathbf{grad}_{\mathbf{x}}u)(\Phi(\tilde{\mathbf{x}})) \quad \forall \tilde{\mathbf{x}} \in \tilde{K}. \quad (3.2.18)$$

Proof: **chain rule**:

$$\frac{\partial u}{\partial x_i}(\mathbf{x}) = \frac{\partial}{\partial x_i} \Phi^*u(\Phi^{-1}(\mathbf{x})) = \sum_{j=1}^d \frac{\partial \Phi^*u}{\partial \tilde{x}_j}(\Phi^{-1}(\mathbf{x})) \frac{\partial \Phi_j^{-1}}{\partial x_i}(\mathbf{x}).$$

$$\begin{aligned} \blacktriangleright \quad \mathbf{grad} u(\mathbf{x}) &= (D\Phi^{-1}(\mathbf{x}))^T \mathbf{grad}_{\tilde{\mathbf{x}}}(\Phi^*u)(\Phi^{-1}(\mathbf{x})) \\ &= D\Phi(\Phi^{-1}(\mathbf{x}))^{-T} (\mathbf{grad} \Phi^*u)(\Phi^{-1}(\mathbf{x})). \end{aligned}$$

Parametric construction:

$$b_i^{\tilde{K}} = \Phi^*b_i^K, \quad i = 1, \dots, Q$$

Local shape functions on \tilde{K} Local shape functions on K

Local computations use (3.2.18) & **transformation formula** (for multidimensional integrals):

$$\int_K f(\Phi(\mathbf{x})) d\mathbf{x} = \int_{\tilde{K}} f(\tilde{\mathbf{x}}) |\det D\Phi(\tilde{\mathbf{x}})| d\tilde{\mathbf{x}} \quad \text{for } f : K \mapsto \mathbb{R},$$

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$$\begin{aligned} \int_K \mathbf{grad} u \cdot \mathbf{grad} v d\mathbf{x} &= \int_{\tilde{K}} (\mathbf{grad} u)(\Phi(\tilde{\mathbf{x}})) \cdot (\mathbf{grad} v)(\Phi(\tilde{\mathbf{x}})) |\det D\Phi(\tilde{\mathbf{x}})| d\tilde{\mathbf{x}} \\ &= \int_{\tilde{K}} D\Phi^{-T}(\tilde{\mathbf{x}}) \mathbf{grad}_{\tilde{\mathbf{x}}}(\Phi^*u) \cdot D\Phi^{-T}(\tilde{\mathbf{x}}) \mathbf{grad}_{\tilde{\mathbf{x}}}(\Phi^*v) |\det D\Phi(\tilde{\mathbf{x}})| d\tilde{\mathbf{x}}. \end{aligned}$$

$$\int_K \mathbf{grad} b_i^K \cdot \mathbf{grad} b_j^K d\mathbf{x} = \int_{\tilde{K}} \left\{ D\Phi(\tilde{\mathbf{x}})^T D\Phi(\tilde{\mathbf{x}}) \right\}^{-1} \mathbf{grad} b_i^{\tilde{K}} \cdot \mathbf{grad} b_j^{\tilde{K}} |\det D\Phi(\tilde{\mathbf{x}})| d\tilde{\mathbf{x}}.$$

Note: local shape functions $b_i^{\tilde{K}}$ simple polynomials !

For parabolic boundary fitting:

$$D\Phi = Id + 4\delta \mathbf{n} \cdot \mathbf{grad}(\lambda_1\lambda_2)^T \in \mathbb{R}^{2,2}, \quad \det(D\Phi) = 1 + 4\delta \mathbf{n} \cdot \mathbf{grad}(\lambda_1\lambda_2).$$

Next: numerical quadrature (→ Sect. 3.2.6) on \tilde{K}

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3.2.9 Static condensation

interior basis functions = global basis functions supported inside a cell

(occur for $\mathcal{S}_3^0(\mathcal{M})$ on triangular mesh \mathcal{M} in 2D)

Sorting of global basis functions: coefficients for interior basis functions last

► Block structure of resulting linear system $\mathbf{A}\vec{\mu} = \vec{\varphi}$

$$\mathbf{A}\vec{\mu} = \begin{pmatrix} \mathbf{A}_{oo} & \mathbf{A}_{oi} \\ \mathbf{A}_{io} & \mathbf{A}_{ii} \end{pmatrix} \begin{pmatrix} \vec{\mu}_o \\ \vec{\mu}_i \end{pmatrix} = \begin{pmatrix} \vec{\varphi}_o \\ \vec{\varphi}_i \end{pmatrix} = \vec{\varphi}. \quad (3.2.19)$$

\mathbf{A}_{ii} ← coupling among interior basis functions

\mathbf{A}_{oi} ← coupling between interior b.f. & basis functions on nodes/edges

Note: \mathbf{A}_{ii} is **block-diagonal** with small blocks ► “easy to invert”

[Elimination of $\vec{\mu}_i$ (**Static condensation**)]

$$\text{Schur complement system: } (\mathbf{A}_{oo} - \mathbf{A}_{oi}\mathbf{A}_{ii}^{-1}\mathbf{A}_{io}) \vec{\mu}_o = \vec{\varphi}_o - \mathbf{A}_{oi}\mathbf{A}_{ii}^{-1}\vec{\varphi}_i.$$

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