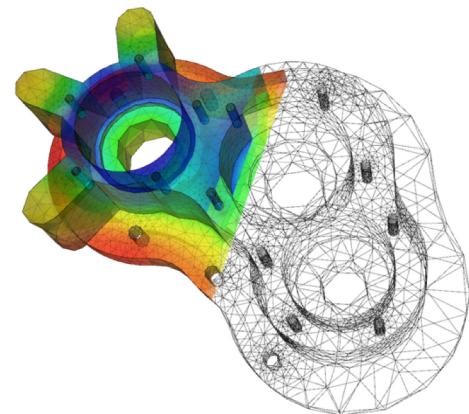


Lecture VIII: Finite Element Simulation

Motivation

- **Continuous physics:** rigid, deforming, attaching, detaching bodies.
 - “continuous” features
 - “infinite” resolution
- Computer simulation:
 - Discrete representation with limited memory.
 - Efficiency is key.
 - ...and also **consistency** and **stability**.

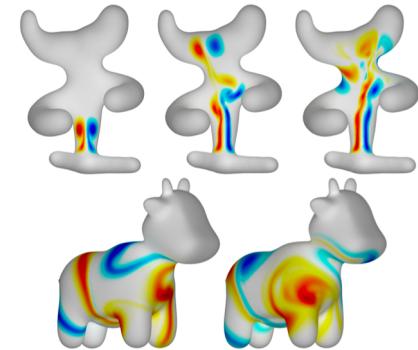


Partial Differential Equations

- Govern the connection between kinematics and dynamics.
- Examples:
 - $\vec{F} = m\vec{a} = m \frac{d\vec{v}}{dt} = -\vec{v}$ (drag).
 - $\vec{F} = -k\vec{x} = \frac{d\vec{x}}{dt}$ (spring; Hooke's law).
- We learned how to integrate in **time**.
- How to discretize in **space**?

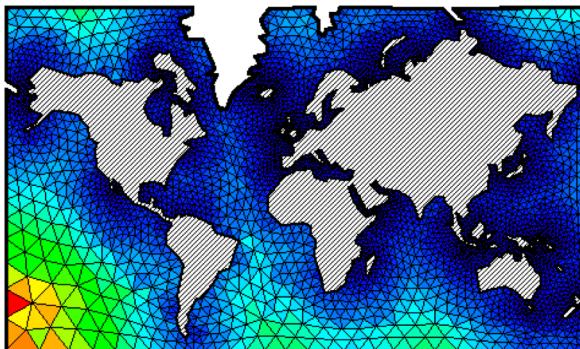
Partial Differential Equations

- Challenges:
 - Discretize (partial) derivatives
 - Obey conservation rules
 - Consistency
 - Good sampling

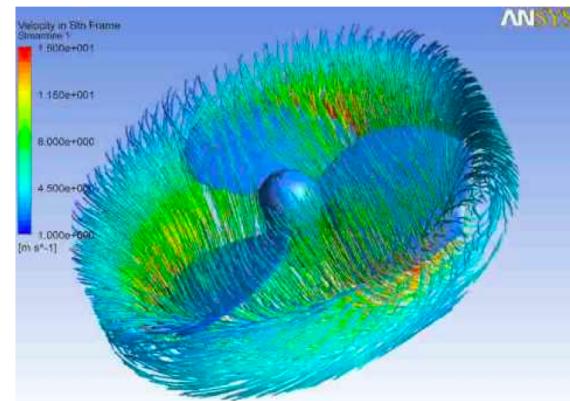


[Azencot *et al.* 2015]

- Representation choices:
 - Where do functions\vecs\tensors “live”?



<http://caewatch.com/top-5-misunderstandings-on-good-mesh/>

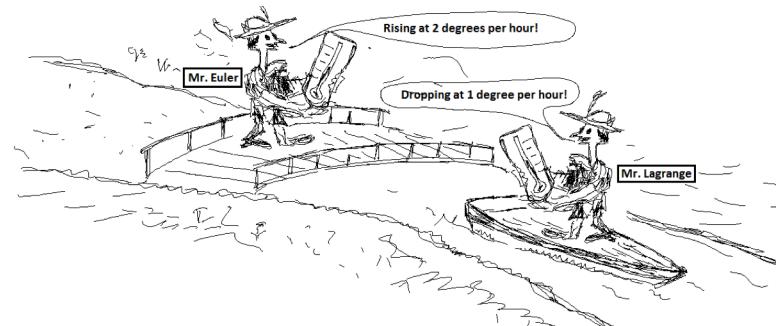


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<https://www.youtube.com/watch?v=5mNn7csNGDk>

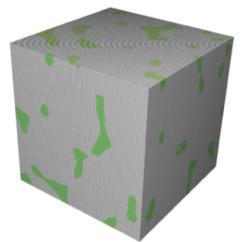
Numerical Simulation

- Lagrangian:
 - Representation: connected **mesh** or **cloud** of particles.
 - **Examples:** Finite methods, Mass-spring system, particle systems.
- Eulerian:
 - Representation: stationary point set or **grid**, where material properties change over time.
 - Boundary of object not explicitly defined.
 - Suitable for fluids.

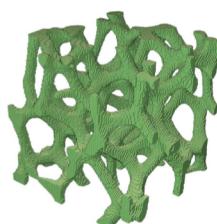


Finite Differences Method

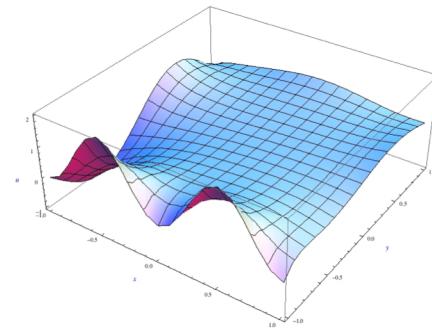
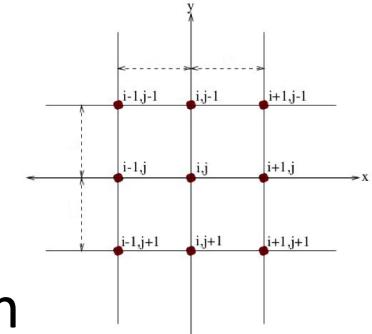
- Object sampled using **regular spatial grid**.
- PDE discretized using **finite differences**.
 - **Pro:** easier and more stable to implement than unstructured methods.
 - **Con:** difficult to approximate complex boundaries.
- Semi-implicit integration is used to move forward through time



(a) Voxelized sample cube, consisting of voxels with material index 1 (aluminum) and 0 (void).

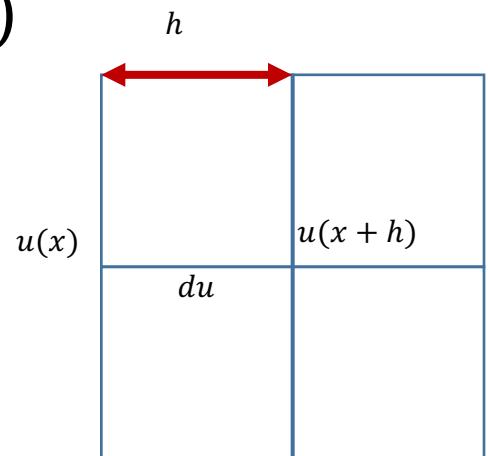


(b) Voxels with material index 1, representing the foam geometry.



Finite Differences Method

- Grid size h .
- Scalar functions: at vertices.
- Differentials: $du = u(x + h) - u(x)$
- Derivatives: $\frac{du}{dx} = \frac{u(x+h)-u(x)}{h} + O(h)$
 - "Living" on edges



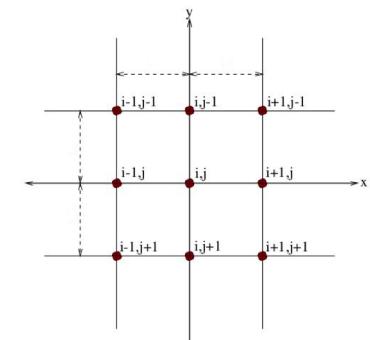
Example: Heat Equation

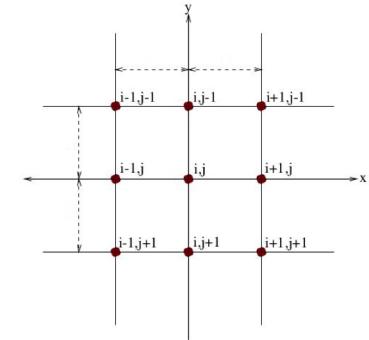
- $u_t = c\Delta u$.
 - In 2D: $u_t = c(u_{xx} + u_{yy})$
- Forward discretization in time:

$$u_t \approx \frac{u_{i,j}^{t+\Delta t} - u_{i,j}^t}{\Delta t}$$

- 2nd order derivative **central** approximation:

$$u_{xx} = (u_x)_x \approx \frac{\frac{u_{i+1}^t - u_i^t}{h} - \frac{u_i^t - u_{i-1}^t}{h}}{h} = \frac{u_{i+1}^t - 2u_i^t + u_{i-1}^t}{h^2}$$

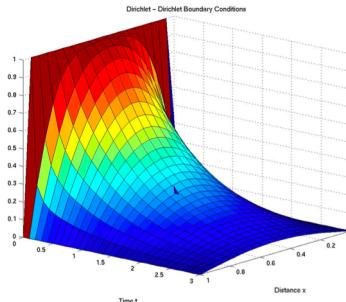




Example: Heat Equation

$$\begin{aligned}
 & u_{xx} + u_{yy} \\
 & \approx \frac{u_{i+1,j}^t - 2u_{i,j}^t + u_{i-1,j}^t}{h^2} + \frac{u_{i,j+1}^t - 2u_{i,j}^t + u_{i,j-1}^t}{h^2} \\
 & = \frac{u_{i+1,j}^t + u_{i-1,j}^t + u_{i,j+1}^t + u_{i,j-1}^t - 4u_{i,j}^t}{h^2}
 \end{aligned}$$

- Measuring difference of neighbor average from center!
- **Dirichlet Boundary conditions:** set constant values

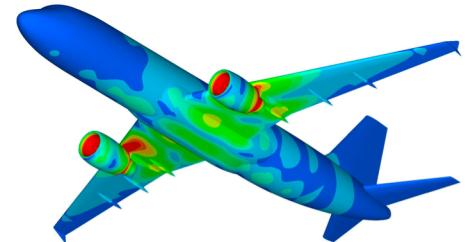


<http://www.math.oregonstate.edu/~show/images/Dd.jpg>

Boundary Element Method

- Remember Stokes theorem:

$$\int_{\partial\Omega} \omega = \int_{\Omega} d\omega$$



<http://urbana.mie.uc.edu/yliu/>

- Practical interpretation: the sum of what “happens inside” only depends on “what goes in and out”.
- Discretizing PDE only on the **boundary surface** instead of the **entire volume**.
 - Only good for **homogenous** material.
 - Topological changes more difficult to handle.

Finite Volume Methods

- Measure flux changes in small (simplicial) elements

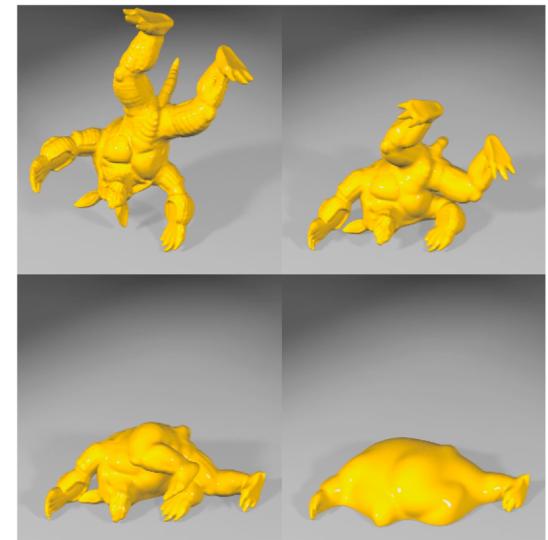
- Like a BEM on a FEM.

$$\frac{\partial}{\partial t} \iiint Q dV + \iint F dA = 0$$

- Q : variable inside the domain
 - F : flux on the boundary.

- Conserves volume by **definition!**

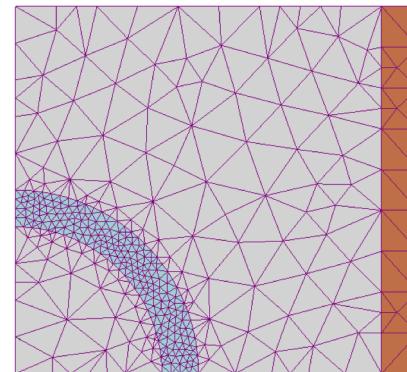
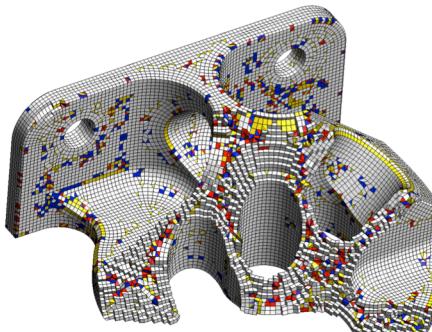
- Good for CFD.



"A simple finite volume method for adaptive viscous liquids" by [Batti *et al.* 2011]

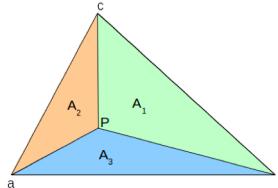
Finite Element Method (FEM)

- Tessellating the volume into a large finite number of disjoint elements (3D volumetric/surface mesh).
- Usually: simplices. Sometimes quads/hexes.
- Scalar functions: at vertices.
- Differentials on edges e_{ij} : $du = u_j - u_i$.
- Derivatives?



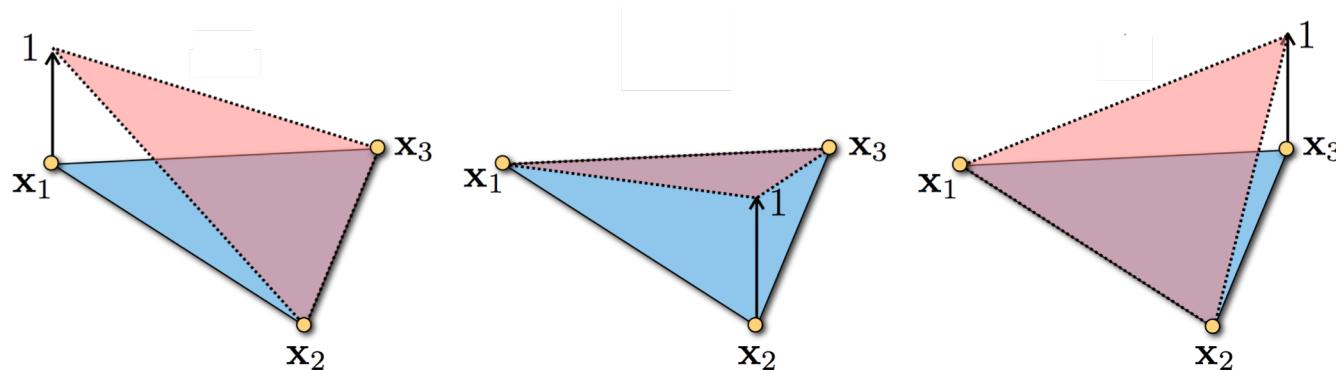
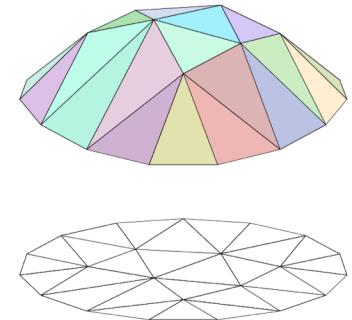
Lagrange FEM Basis

- Function values defined on vertices, but
 - Assumed to **interpolate linearly** on elements:



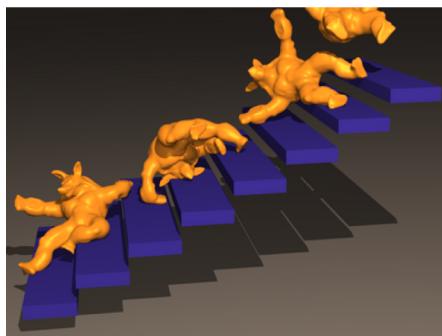
$$u(p) = \sum_{i=0}^3 B_i(p) u_i$$

- $B_i(p)$: **Barycentric coordinates** of p in element.



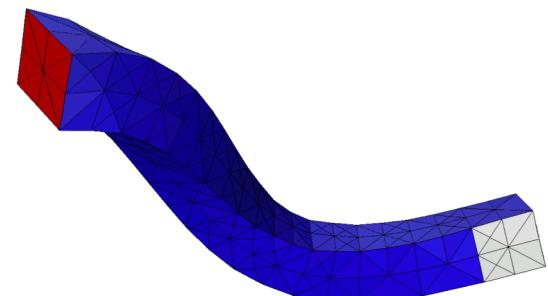
Motivation

- Discretize deformations, vectors, and tensors
- Derivative and integrals => linear operators (matrices)
 - Solving PDEs \Leftrightarrow solving linear equations
 - Either directly, or by iterations.



Irving *et al.* "Volume Conserving Finite Element Simulations of Deformable Models"

<https://www.youtube.com/watch?v=Rbq2CdUlvw4>



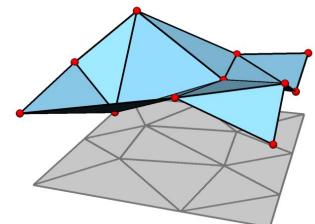
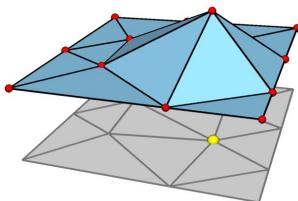
http://fetk.org/codes/mc/images/bar3ndef_wbg.gif

Lagrange FEM Basis

- Equivalent definition: function is defined by per-vertex u_v , scalar values
- Entire space is spanned by Lagrange basis functions

$$u(p) = \sum_{v \in V} u_v \varphi_v(p)$$

- $\varphi_v(p) = 1$ on v , and 0 on other vertices.
 - Piecewise linear (“hat function”).

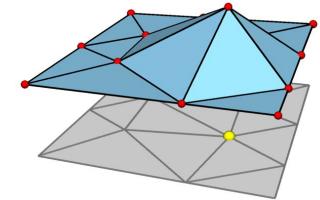


$$u(p) = \sum_{v \in V} u_i \varphi_v(p)$$

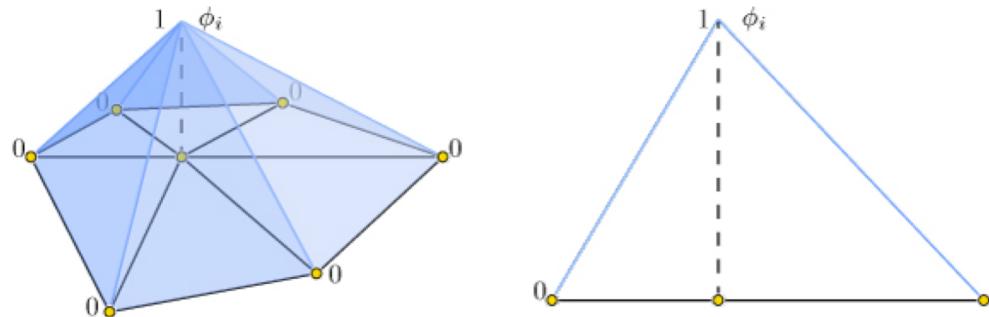
Lagrange FEM Basis

- If scalar functions are piecewise linear on vertices:
 - Gradients ∇u are piecewise constant on faces!

$$\nabla u(p) = \sum_{v \in V} u_i \nabla \varphi_v(p)$$



- Example: **position** function on vertices, **deformation** (velocity) on faces.

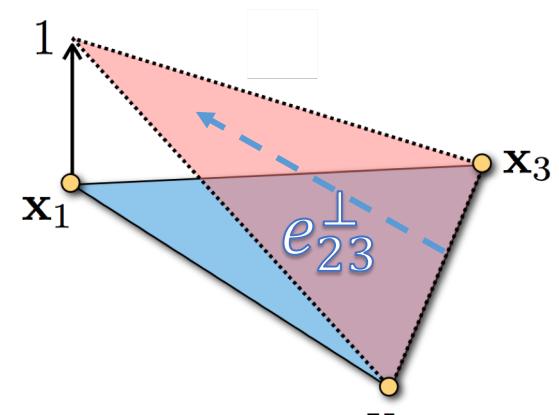


Gradients

- We want to compute $\nabla \varphi_1$ (w.l.o.g.) in face 123.
- **Insight:** in the direction of $e_{23}^\perp = \hat{n} \times e_{23}$.
- **Result:** $\nabla \varphi_1 = \frac{e_{23}^\perp}{2A_{123}}$
 - A_{123} : area of triangle 123.
- For a function u_1, u_2, u_3 :

$$\nabla u_{123} = Gu_{123} = \frac{1}{2A_{123}} (e_{23}^\perp u_1 + e_{31}^\perp u_2 + e_{12}^\perp u_3)$$

derived from geometry



Gradients

- In matrix form (\vec{e}_{23}^\perp etc. are column vectors 1x3):

$$\begin{aligned} \bullet G_{123} u &= \frac{1}{2A_{123}} (\vec{e}_{23}^\perp \quad \vec{e}_{31}^\perp \quad \vec{e}_{12}^\perp) \begin{pmatrix} u_1 \\ u_2 \\ u_3 \end{pmatrix} \\ &= \begin{pmatrix} \partial u / \partial x \\ \partial u / \partial y \\ \partial u / \partial z \end{pmatrix}_{123} \end{aligned}$$

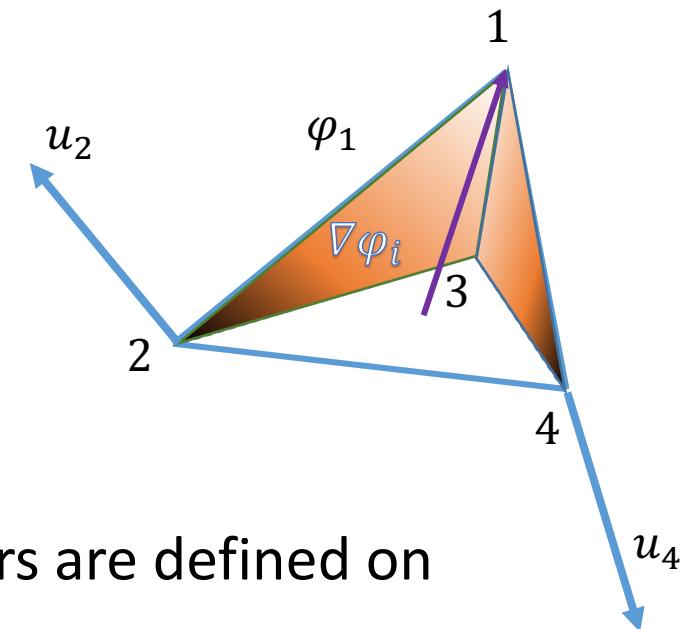
- Aggregating, we can get global matrix $G_{3|F| \times |V|}$
 - PL vertex functions -> face PC gradients.
 - $3|F|$ - one vector (xyz) in each face.

Linear Elasticity in FEM

- **Discretization:** deformation field on vertices
 - Where each vertex moves.
 - PL inside each tetrahedron\triangle.
 - Reminder: Jacobian: $Ju(p): \mathbb{R}^{3 \times 3} = \begin{pmatrix} \nabla u_x(p) \\ \nabla u_y(p) \\ \nabla u_z(p) \end{pmatrix}$
 - We simply work with three scalar functions: u_x, u_y, u_z .
 - Denoting: $\vec{u}_x = \begin{pmatrix} u_{x,1} \\ u_{y,1} \\ u_{z,1} \end{pmatrix}$
- $$\begin{pmatrix} \nabla u_x^T \\ \nabla u_y^T \\ \nabla u_z^T \end{pmatrix} = J_{123} \begin{pmatrix} \vec{u}_x \\ \vec{u}_y \\ \vec{u}_z \end{pmatrix} = \begin{pmatrix} G_{123} & & \\ & G_{123} & \\ & & G_{123} \end{pmatrix} \begin{pmatrix} \vec{u}_x \\ \vec{u}_y \\ \vec{u}_z \end{pmatrix}$$
- **Sanity check:** J_{123} is a 9×9 matrix.

Tetrahedral Meshes

- Elements $e = [1234]$.
- “straightforward” generalization
 - Scalar functions/deformation vectors are defined on vertices
 - Piecewise-linear inside **each tet volume**.
 - Gradients piecewise-constant in each tet.
 - → Gradients orthogonal to opposite face normal.
- Single tet Jacobian: J_e : 9×12 matrix
 - Input vertex deformation field $3 \times 4 = 12$.
 - Output tet jacobian: $3 \times 3 = 9$.



Linear Elasticity in FEM

- Full Lagrangian strain tensor:

$$\mathbf{E} = \frac{1}{2} (J\mathbf{u}^T J\mathbf{u} + J\mathbf{u} + J\mathbf{u}^T)$$

- Not linear inside tets!
- Linear elasticity approximation:
$$\boldsymbol{\epsilon} \approx \frac{1}{2} (J\mathbf{u} + J\mathbf{u}^T)$$
- Good for small deformations without much rotation.

Discrete Strain Tensor

- Written explicitly:

$$\boldsymbol{\epsilon} = \frac{1}{2} (Ju + Ju^T) = \frac{1}{2} \begin{pmatrix} 2\frac{\partial u_x}{\partial x} & \frac{\partial u_x}{\partial y} + \frac{\partial u_y}{\partial x} & \frac{\partial u_x}{\partial z} + \frac{\partial u_z}{\partial x} \\ \frac{\partial u_y}{\partial x} + \frac{\partial u_x}{\partial y} & 2\frac{\partial u_y}{\partial y} & \frac{\partial u_y}{\partial z} + \frac{\partial u_z}{\partial y} \\ \frac{\partial u_z}{\partial x} + \frac{\partial u_x}{\partial z} & \frac{\partial u_z}{\partial y} + \frac{\partial u_y}{\partial z} & 2\frac{\partial u_z}{\partial z} \end{pmatrix}$$

- Only 6 relevant elements (rest are symmetric):

$$\begin{pmatrix} \epsilon_{xx} \\ \epsilon_{yy} \\ \epsilon_{zz} \\ \epsilon_{xy} \\ \epsilon_{yz} \\ \epsilon_{zx} \end{pmatrix} = \begin{pmatrix} \frac{\partial}{\partial x} & & & & & \\ & \frac{\partial}{\partial y} & & & & \\ & & \frac{\partial}{\partial z} & & & \\ 0.5 \frac{\partial}{\partial y} & 0.5 \frac{\partial}{\partial x} & & & & \\ & & 0.5 \frac{\partial}{\partial z} & 0.5 \frac{\partial}{\partial y} & & \\ 0.5 \frac{\partial}{\partial z} & & & & 0.5 \frac{\partial}{\partial x} & \end{pmatrix} \begin{pmatrix} u_x(p) \\ u_y(p) \\ u_z(p) \end{pmatrix}$$

Discrete Strain Tensor

- In a **single tet**: $J_e u$ is constant $\rightarrow \boldsymbol{\varepsilon}_e$ is constant.
- We want to get $\boldsymbol{\varepsilon}_e$ as a direct function of u .
- We have:

$$\boldsymbol{\varepsilon}_e = \begin{pmatrix} \varepsilon_{xx} \\ \varepsilon_{yy} \\ \varepsilon_{zz} \\ \varepsilon_{xy} \\ \varepsilon_{yz} \\ \varepsilon_{zx} \end{pmatrix} = D J_e \begin{pmatrix} \vec{u}_x \\ \vec{u}_y \\ \vec{u}_z \end{pmatrix},$$

Where D is a permutation matrix that “chooses” the corresponding values of $J_e u$

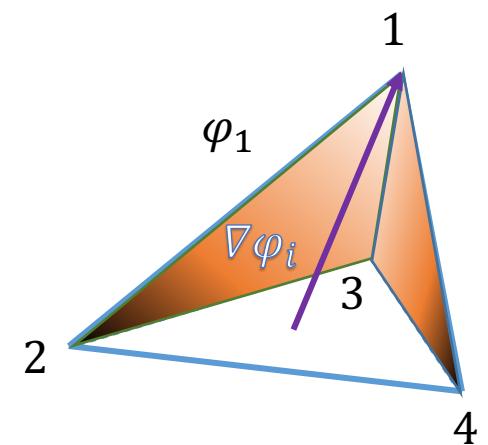
- **Sanity check**: D is of size 6×9 and **does not depend** on the shape of any tet (only contains non-zeros values “1” or “0.5”).

Discrete Strain Tensor

- Strain tensor per face:

$$\boldsymbol{\varepsilon}_e = D\mathbf{J}_e \mathbf{u}_e = \mathbf{B}_e \mathbf{u}_e$$

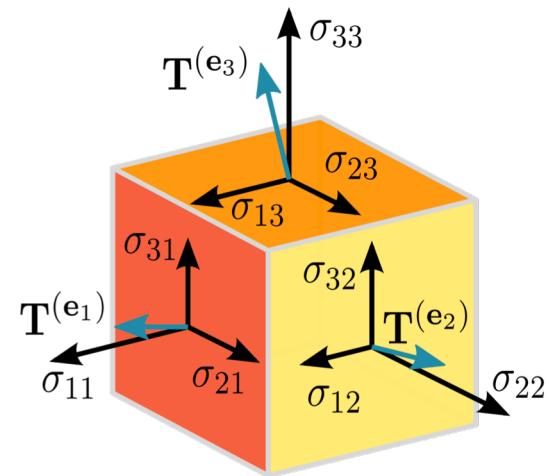
- \mathbf{u}_e : all $u_{xyz,1234}$ values (12 in total).
- **Note:** \mathbf{B}_e : $\mathbb{R}^{6 \times 12}$.
- \mathbf{B}_e contains derivatives of $\varphi_{1,2,3,4}(p)$
 - $\varphi_{1,2,3,4}$ are **linear** inside e .
 - Derivatives of φ_i are **constant** inside e .
- \mathbf{B}_e is constant inside the element!



Discrete Stress Tensor

- Stress and strain are related by Hooke's law
 - Remember $\vec{F} = -k\vec{x}$?
- Again, because of symmetry, we use the compact form:

$$\sigma_e = \begin{pmatrix} \sigma_{xx} \\ \sigma_{yy} \\ \sigma_{zz} \\ \sigma_{xy} \\ \sigma_{yz} \\ \sigma_{zx} \end{pmatrix}$$



Discrete Stress Tensor

- Stress and strain are related by discrete **stiffness tensor** $\mathbf{C}_e: \mathbb{R}^{6 \times 6}$.
$$\sigma_e = \mathbf{C}_e \varepsilon_e = \mathbf{C}_e B_e u_e$$
- Relatively easy structure in the discrete case:

$$\mathbf{C}_e = \begin{pmatrix} \lambda + 2\mu & \lambda & \lambda & & & \\ \lambda & \lambda + 2\mu & \lambda & & & \\ \lambda & \lambda & \lambda + 2\mu & & & \\ & & & \mu & & \\ & & & & \mu & \\ & & & & & \mu \end{pmatrix}$$

- Where $\mu = \frac{Y}{2(1+v)}$, $\lambda = \frac{\nu Y}{(1+\nu)(1-2\nu)}$
 - Y – Young's modulus.
 - ν – Poisson's ratio.
- μ and λ are known as Lamé coefficients.

Strain Energy

- Potential energy gained when applying strain to object:

$$U_e = \frac{1}{2} \int_T \langle \boldsymbol{\sigma}_e, \boldsymbol{\varepsilon}_e \rangle dV$$

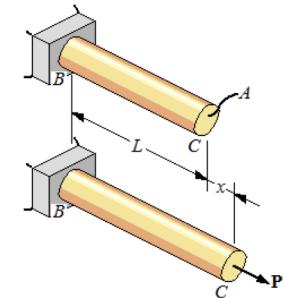
- We have that $\boldsymbol{\sigma}_e = \mathbf{C}_e B_e u_e$ and $\boldsymbol{\varepsilon}_e = B_e u_e$.
- Then:

$$\langle \boldsymbol{\sigma}_e, \boldsymbol{\varepsilon}_e \rangle = (\boldsymbol{\sigma}_e)^T \boldsymbol{\varepsilon}_e = u_e^T B_e^T C_e B_e u_e$$

- Both are constant inside volume, so:

$$U_e = \frac{1}{2} \text{Vol}(e) \times u_e^T B_e^T C_e B_e u_e = \boxed{\frac{1}{2} u_e^T K_e u_e}$$

- K_e : stiffness matrix of 12×12
 - Only depends on the original geometry and material properties!



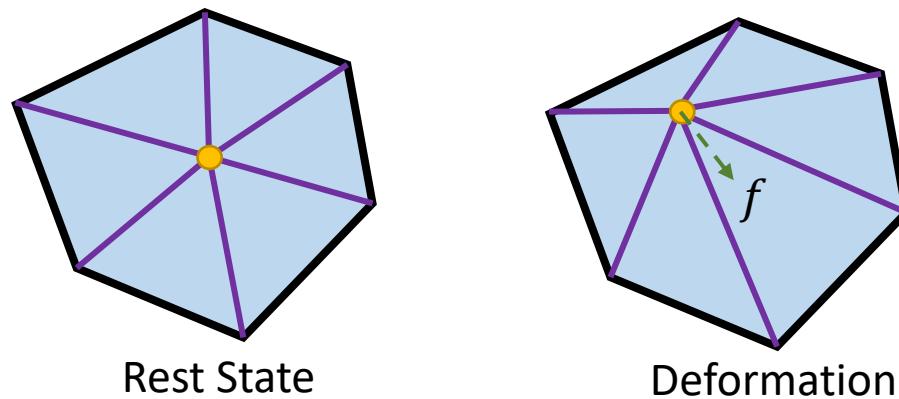
http://www.engineeringarchives.com/img/les_mom_strainenergydensity_1.png

Elastic Forces

- Derivatives of the potential energy:

$$f_e = \frac{\partial U_e}{\partial u_e} = K_e u_e$$

- Each element of f_e corresponds to an element of u_e
 - Inner force acting on that coordinate due to deformation.



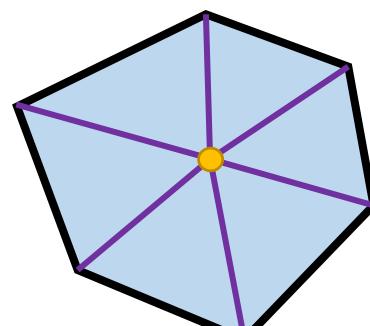
$$f_e = \frac{\partial U_e}{\partial u_e} = K_e u_e$$

Elastic Forces

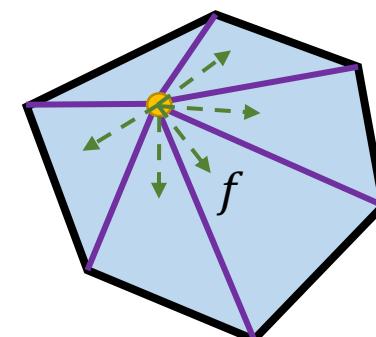
- Net force on each vertex: sum of forces from all adjacent elements.
- Aggregated into a big matrix:

$$f = Ku$$

- $u_{1 \times 3|\nu|}, f_{1 \times 3|\nu|}, K_{3|\nu| \times 3|\nu|}$
- Interpretation: a “force Laplacian”.
- Pulling vertex toward “rest-state equilibrium”.



Rest State



Deformation

Dynamic Deformation Equation

- Computing new positions $x(t)$:

$$Mx''(t) + Dx'(t) + K(x - x_0) = f_{ext}$$

- M : Mass matrix
- D : Damping matrix
- K : stiffness matrix
- f_{ext} : other forces (gravity etc.)

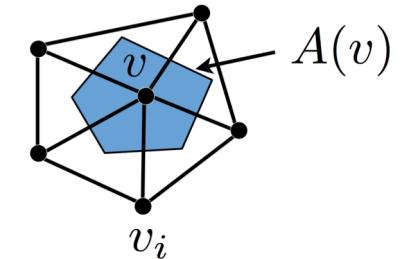


$$Mx''(t) + Dx'(t) + K(x - x_0) = f_{ext}$$

Mass Matrix

- Suppose we are given density ρ for each tetrahedron.
- We need mass **per vertex** (where acceleration is).
- Solution: use **Voronoi area**:

$$m_v = \sum_{e \in N(v)} \frac{1}{4} \rho V(e)$$



- Assuming
- Mass matrix is then

$$M = \text{diag}\{m\}$$

- Each element in the diagonal corresponds to which vertex is represented in the respective column of x .

Dampening Matrix

- We often use Rayleigh Dampening:

$$D = \alpha M + \beta K$$

- α, β are dampening parameters.
 - A whole theory of how to compute them and to what effect...
- For our purposes: keep as parameters, limit by small values $0 \leq \alpha, \beta \leq 0.02$.

$$Mx''(t) + Dx'(t) + K(x - x_0) = f_{ext}$$

Time Integration

- Best practice: a variant on implicit backward integration
- Solve for:

$$x(t + \Delta t) = x(t) + \Delta t \cdot v(t + \Delta t)$$

And similarly for $v(t + \Delta t)$ and $a(t + \Delta t)$.

- We need to solve the linear system:

Constant matrix in $[M + \Delta t D + \Delta t^2 K]v(t + \Delta t)$ variables
time!

$$= Mv(t) - \Delta t [K(x(t) - x_0) - f_{ext}]$$

Right-hand side, changes with time.

- Integrating for position with the formula above.

$$\begin{aligned}[M + \Delta t D + \Delta t^2 K]v(t + \Delta t) \\ = Mv(t) - \Delta t [K(x(t) - x_0) - f_{ext}]\end{aligned}$$

Solving the system

- We need to solve the linear system in every time step.
- Abstract form: $Ax(t) = b(t)$, A time independent.
- Sounds slow; however, we should use the fact that A does not change!
 - Reduce A to some convenient form in the beginning of time
 - In each iteration: change $b(t)$ and solve for $x(t)$.

LU Decomposition

- Every **square** matrix A can be decomposed into $A = L \cdot U$, where:
 - L **lower** triangular
 - U **upper** triangular

$$\bullet A = \begin{pmatrix} A_{11} & \cdots & A_{1n} \\ \vdots & \ddots & \vdots \\ A_{n1} & \cdots & A_{nn} \end{pmatrix} =$$
$$L \quad \boxed{\begin{pmatrix} L_{11} & \cdots & 0 \\ \vdots & \ddots & \vdots \\ L_{n1} & \cdots & L_{nn} \end{pmatrix}} \cdot \boxed{\begin{pmatrix} U_{11} & \cdots & U_{1n} \\ \vdots & \ddots & \vdots \\ 0 & \cdots & U_{nn} \end{pmatrix}} \quad U$$

Solving systems with LU Decomposition

- Decompose $Ax = LUx$.
- Solve in two steps:
 - Solve $Ly = b$ (one forward substitution)
 - Solve $Ux = y$ (one backward substitution)
- Decomposition is relatively expensive.
- But solving for a given RHS is cheap.
- **Consequence:** decomposition is very beneficial for **fixed A and varying b !**

Cholesky Decomposition

- If M is symmetric positive definite, LU reduces to:

$$M = LL^*$$

- * means “*conjugate transpose*”
 - For **real** matrices: just LL^T .
 - For **complex** matrices: every number $a + ib$ becomes $\overline{a + ib} = a - ib$.
- Cheaper, more stable decomposition.

Sparse Representation

- M, D, K have few elements w.r.t. to their sizes
 - M just diagonal.
 - K creates forces for vertices only from their neighbors.
- Efficiency and memory requirements: **use sparse matrices!**
- Direct Representation:
 - List of triplets (row, col, value).
- Example:

$$A = \begin{pmatrix} 1 & & \\ & & -5 \\ & 4 & \end{pmatrix}$$

Represented as $\{(1,1,1), (2,3, -5), (3,2,4)\}$.

- The common way to feed sparse matrices.
- Sometimes called **COOrdinate format (COO)**

Compressed Sparse Row/Column

- The common way to represent and manipulate in memory.
- For **CSR** we have 3 arrays A,IA,JA
 - A: list of non-zeros (NZ) row-by-row
 - IA: defined recursively:
 - IA[0]=0
 - IA[i]=IA[i-1]+NNZ(row(i-1))
 - JA: column value of each element in A.
- **CSC**: the same with columns instead of rows and vice versa.

$$\begin{pmatrix} 10 & 20 & 0 & 0 & 0 & 0 \\ 0 & 30 & 0 & 40 & 0 & 0 \\ 0 & 0 & 50 & 60 & 70 & 0 \\ 0 & 0 & 0 & 0 & 0 & 80 \end{pmatrix}$$

$$\begin{aligned} A &= [10 20 30 40 50 60 70 80] \\ IA &= [0 2 4 7 8] \\ JA &= [0 1 3 2 3 4 5] \end{aligned}$$

Sparse Representation

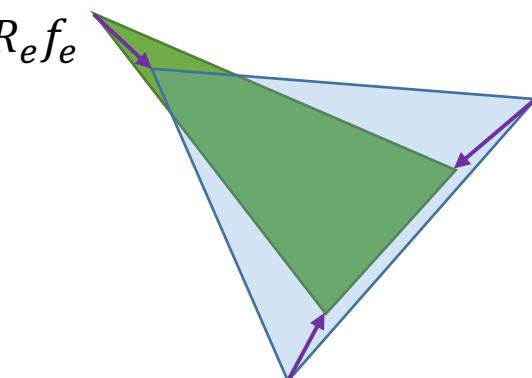
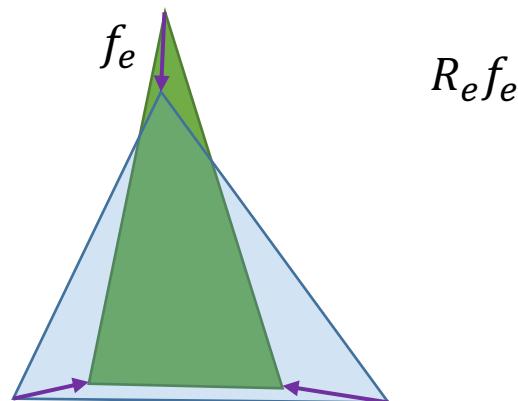
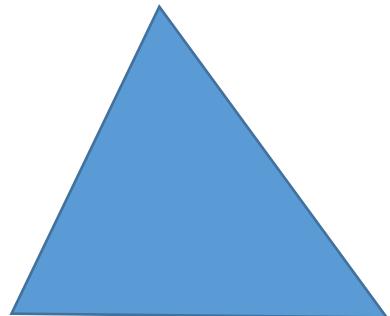
- Cheap:
 - Sparse * vector for CSR.
 - Vector * sparse for CSC.
 - Slicing (but don't use that).
- What's Relatively cheap:
 - Decompositions (sparse LU, Cholesky, etc.)
- What's Expensive:
 - Transpose (conversion from CSC to CSR)
 - Random memory access.

“Index Hell”

- Figuring out the matrices can be quite a job.
 - “which vertex, which coordinate xyz to which element multiplies with what” → index hell.
 - Easy to get wrong!
- Life-(and sanity-; and grade-) saving tip:
 - Create Complicated matrices (such as K) through the composition of easier matrices.
- Example: compute a huge K' that is just made of blocks of the much easier K_e per-element.
- Compose with a matrix M that averages numbers from elements to adjacent vertices to get $K = MK'M^T$.
 - Similar with the matrices G, J , etc.

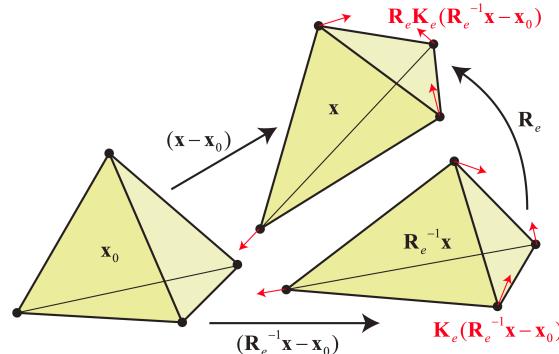
Corotational Elements

- **Insight:** rotating an element should not change the strain energy.
- **Conclusion:** Given an element in position x , with strain energy U_e , and consequent elastic forces f_e , the forces on a rotated element $R_e x$ should be $R_e f_e$!



Corotational Elements

- Method:
 - Estimate rotation R_e , and factor out rotation from the deformed object x : $R_e^{-1}x$
 - Compute elastic forces of unrotated object:
$$K_e(R_e^{-1}x - x_0)$$
 - Rotate back to deformed state to get actual forces:
$$f_e = R_e K_e(R_e^{-1}x - x_0) = K_e' u_e$$



Corotational Elements

- **Advantages:** able to work with large rotations
- **Disadvantages:** stiffness matrix not constant anymore.
 - Have to solve system using conjugate gradients...
- How to estimate rotation R_e ?

Finding Best-Fit Rotation

- Original positions x , deformed positions x' .
- *Create* stacked coordinates of edges of original points:

$$P = \begin{pmatrix} x_1 - x_2 \\ x_1 - x_d \end{pmatrix}, Q = \begin{pmatrix} x_1' - x_2' \\ x_1' - x_d' \end{pmatrix}$$

- Compute matrix: $S = P^T Q \in \mathbb{R}^{3 \times 3}$
- **Singular value decomposition** (SVD) extracts rotation from S

$$\mathbf{S} = \mathbf{U}\Sigma\mathbf{V}^T \longrightarrow \mathbf{R} = \mathbf{U}\mathbf{V}^T$$

Singular Value Decomposition

- Every linear operator (=matrix $M_{n \times m}$) can be decomposed to:
 - **Rotation** (Change of basis): $V_{m \times m}$.
 - **Stretch** in the new basis: $\Sigma_{n \times m}$
 - Note (possible) change in dimension.
 - **Rotation** (another change of basis): $U_{n \times n}$
- For vector p we get $Mp = U\Sigma V^T p$.

Examples

<https://www.youtube.com/watch?v=4WI0ksysYKM>

<https://www.youtube.com/watch?v=6f3UYHnR4zU>

https://www.youtube.com/watch?v=p5uhnSw8_Xw