

Notes

- Some example values for common materials: (VERY approximate)
 - Aluminum: $E=70$ GPa $\nu=0.34$
 - Concrete: $E=23$ GPa $\nu=0.2$
 - Diamond: $E=950$ GPa $\nu=0.2$
 - Glass: $E=50$ GPa $\nu=0.25$
 - Nylon: $E=3$ GPa $\nu=0.4$
 - Rubber: $E=1.7$ MPa $\nu=0.49\dots$
 - Steel: $E=200$ GPa $\nu=0.3$

2D FVM strain

- Triangle with corners i, j, k
- Coordinates $x_i=X(p_i)$, $x_j=X(p_j)$, $x_k=X(p_k)$
- Assume affine in this triangle: $X(p)=Ap+b$ (so $\partial X/\partial p$)
- Then $x_i-x_j = (Ap_i+b)-(Ap_j+b) = A(p_i-p_j)$
- And $x_i-x_k = A(p_i-p_k)$
- Let $\Delta x=[x_i-x_j \mid x_i-x_k]$ and $\Delta p=[p_i-p_j \mid p_i-p_k]$
- Thus $A\Delta p = \Delta x$
- Then $A = \Delta x \Delta p^{-1}$
- Can precompute and store Δp^{-1}

Finite Volume Method

- We discretize integral equation (complete with boundary integrals) over finite volumes
- Typically Voronoi regions (from Delaunay triangulation) or quads from grid
- Each finite volume has a vertex in the middle
 - Assume volume quantities are constant (equal to value at vertex)
 - Assume deformation affine in each triangle (constant strain \rightarrow constant stress)
 - [evaluate each term]
 - Exact choice of control volumes not critical - constant times normal integrates to zero

2D FVM stress calculation

- Look at single triangle i,j,k again
- Write part of path integral of σn for volume i
- Take constant σ out
- Note integral of normals around closed curve is 0
- Switch to integrating normals along triangle boundary
- So force on i due to this triangle's strain is
 - $F_i += -\sigma((x_i-x_k)^\perp + (x_j-x_i)^\perp)/2$
- Use the integral of normal=0 over triangle:
 - $F_i += \sigma(x_k-x_j)^\perp/2$

Path independence

- Note: trick involving switching path integral to triangle boundary works independent of the choice of finite volume
 - Force we compute only depends on where the finite volumes cut the triangle edges - always assume midpoint
- Choice of finite volume does affect body forces and mass calculation
- Might want to connect to centroids instead of circumcentres
 - always well defined (even if not Delaunay)
 - formulas much simpler: sum $1/3$ triangle area

Finite Element Method

- #1 most popular method for elasticity problems (and many others too)
- FEM originally began with simple idea:
 - Can solve idealized problems (e.g. that strain is constant over a triangle)
 - Call one of these problems an element
 - Can stick together elements to get better approximation
- Since then has evolved into a rigorous mathematical algorithm, a general purpose black-box method
 - Well, almost black-box...

Boundary Conditions

- Actually easier to think about numerically
 - Control volumes automatically include boundary
 - Free boundary: do nothing
 - Specified traction: integrate over decomposed boundary mesh
 - E.g. wind forces...
 - Specified displacement (position): just set x to what it's supposed to be

Modern Approach

- Galerkin framework (the most common)
- Find vector space of functions that solution (e.g. $X(p)$) lives in
 - E.g. bounded weak 1st derivative: H^1
- Say the PDE is $L[X]=0$ everywhere (“strong”)
- The “weak” statement is $\int Y(p)L[X(p)]dp=0$ for every Y in vector space
- Issue: L might involve second derivatives
 - E.g. one for strain, then one for div sigma
 - So L , and the strong form, difficult to define for H^1
- Integration by parts saves the day

Weak Momentum Equation

- Ignore time derivatives - treat acceleration as an independent quantity
 - We discretize space first, then use “method of lines”: plug in any time integrator

$$\begin{aligned}
 L[X] &= \rho \ddot{X} - f_{body} - \nabla \cdot \sigma \\
 \int_{\Omega} Y L[X] &= \int_{\Omega} Y (\rho \ddot{X} - f_{body} - \nabla \cdot \sigma) \\
 &= \int_{\Omega} Y \rho \ddot{X} - \int_{\Omega} Y f_{body} - \int_{\Omega} Y \nabla \cdot \sigma \\
 &= \int_{\Omega} Y \rho \ddot{X} - \int_{\Omega} Y f_{body} + \int_{\Omega} \sigma \nabla Y
 \end{aligned}$$

Linear Triangle Elements

- Simplest choice
- Take basis $\{\phi_i\}$ where $\phi_i(p) = 1$ at p_i and 0 at all the other p_i 's
 - It's a “hat” function
- Then $X(p) = \sum_i x_i \phi_i(p)$ is the continuous piecewise linear function that interpolates particle positions
- Similarly interpolate velocity and acceleration
- Plug this choice of X and an arbitrary $Y = \phi_j$ (for any j) into the weak form of the equation
- Get a system of equations (3 eq. for each j)

Making it finite

- The Galerkin FEM just takes the weak equation, and restricts the vector space to a finite-dimensional one
 - E.g. Continuous piecewise linear - constant gradient over each triangle in mesh, just like we used for Finite Volume Method
- This means instead of infinitely many test functions Y to consider, we only need to check a finite basis
- The method is defined by the basis
 - Very general: plug in whatever you want - polynomials, splines, wavelets, RBF's, ...

The equations

$$\begin{aligned}
 \int_{\Omega} \phi_j \sum_i \rho \ddot{x}_i \phi_i - \int_{\Omega} \phi_j f_{body} + \int_{\Omega} \sigma \nabla \phi_j &= 0 \\
 \sum_i \int_{\Omega} \rho \phi_j \phi_i \ddot{x}_i &= \int_{\Omega} \phi_j f_{body} - \int_{\Omega} \sigma \nabla \phi_j
 \end{aligned}$$

- Note that ϕ_j is zero on all but the triangles surrounding j , so integrals simplify
- Also: naturally split integration into separate triangles

Change in momentum term

- Let $m_{ij} = \int \rho \phi_i \phi_j$
- Then the first term is just $\sum_i m_{ji} \ddot{x}_i$
- Let $M=[m_{ij}]$: then first term is $M\ddot{x}$
- M is called the mass matrix
 - Obviously symmetric (actually SPD)
 - Not diagonal!
- Note that once we have the forces (the other integrals), we need to invert M to get accelerations

Stress term

- Calculate constant strain and strain rate (so constant stress) for each triangle separately
- Note $\nabla \phi_j$ is constant too
- So just take $\sigma \nabla \phi_j$ times triangle area
- [derive what $\nabla \phi_j$ is]
- Magic: exact same as FVM!
 - In fact, proof of convergence of FVM is often (in other settings too) proved by showing it's equivalent or close to some kind of FEM

Body force term

- Usually just gravity: $f_{\text{body}} = \rho g$
- Rather than do the integral with density all over again, use the fact that ϕ_i sum to 1
 - They form a “partition of unity”
 - They represent constant functions exactly - just about necessary for convergence
- Then body force term is $gM1$
- More specifically, can just add g to the accelerations; don't bother with integrals or mass matrix at all

The algorithm

- Loop over triangles
 - Loop over corners
 - Compute integral terms
 - only need to compute M once though - it's constant
 - End up with row of M and a “force”
- Solve $Ma=f$
- Plug this a into time integration scheme

Lumped Mass

- Inverting mass matrix unsatisfactory
 - For particles and FVM, each particle had a mass, so we just did a division
 - Here mass is spread out, need to do a big linear solve - even for explicit time stepping
- Idea of lumping: replace M with the “lumped mass matrix”
 - A diagonal matrix with the same row sums
 - Inverting diagonal matrix is just divisions - so diagonal entries of lumped mass matrix are the particle masses
 - Equivalent to FVM with centroid-based volumes

Locking

- Simple linear basis actually has a major problem: locking
- Notion of numerical stiffness
 - Instead of thinking of numerical method as just getting an approximate solution to a real problem,
 - Think of numerical method as exactly solving a problem that's nearby
 - For elasticity, we're exactly solving the equations for a material with slightly different (and not quite homogeneous/isotropic) stiffness
- Locking comes up when numerical stiffness is MUCH higher than real stiffness

Consistent vs. Lumped

- Original mass matrix called “consistent”
- Turns out its strongly diagonal dominant (fairly easy to solve)
- Multiplying by mass matrix = smoothing
- Inverting mass matrix = sharpening
- Rule of thumb:
 - Implicit time stepping - use consistent M (counteract over-smoothing, solving system anyways)
 - Explicit time stepping - use lumped M (avoid solving systems, don't need extra sharpening)

Locking and linear elements

- Look at nearly incompressible materials
- Can a linear triangle mesh deform incompressibly?
 - [derive problem]
- Then linear elements will resist far too much: numerical stiffness much too high
- Numerical material “locks”
- FEM isn't really a black box!
- Solutions:
 - Don't do incompressibility
 - Use other sorts of elements (quads, higher order)

Quadrature

- Formulas for linear triangle elements and constant density simple to work out
- Formulas for subdivision surfaces (or high-order polynomials, or splines, or wavelets...) and varying density are NASTY
- Instead use “quadrature”
 - I.e. numerical approximation to integrals
- Generalizations of midpoint rule
 - E.g. Gaussian quadrature (for intervals, triangles, tets) or tensor products (for quads, hexes)
- Make sure to match order of accuracy [or not]

Accuracy

- At least for SPD linear problems (e.g. linear elasticity) FEM selects function from finite space that is “closest” to solution
 - Measured in a least-squares, energy-norm sense
- Thus it’s all about how well you can approximate functions with the finite space you chose
 - Linear or bilinear elements: $O(h^2)$
 - Higher order polynomials, splines, etc.: better

Other elements

- Not so obvious ones:
 - Isoparametric elements (meshes with curved edges)
 - Radial-basis functions (mesh-free methods)
 - Mixed element meshes (triangles and quads together)
 - Embedded elements
 - Special-purpose elements (e.g. for cracks)