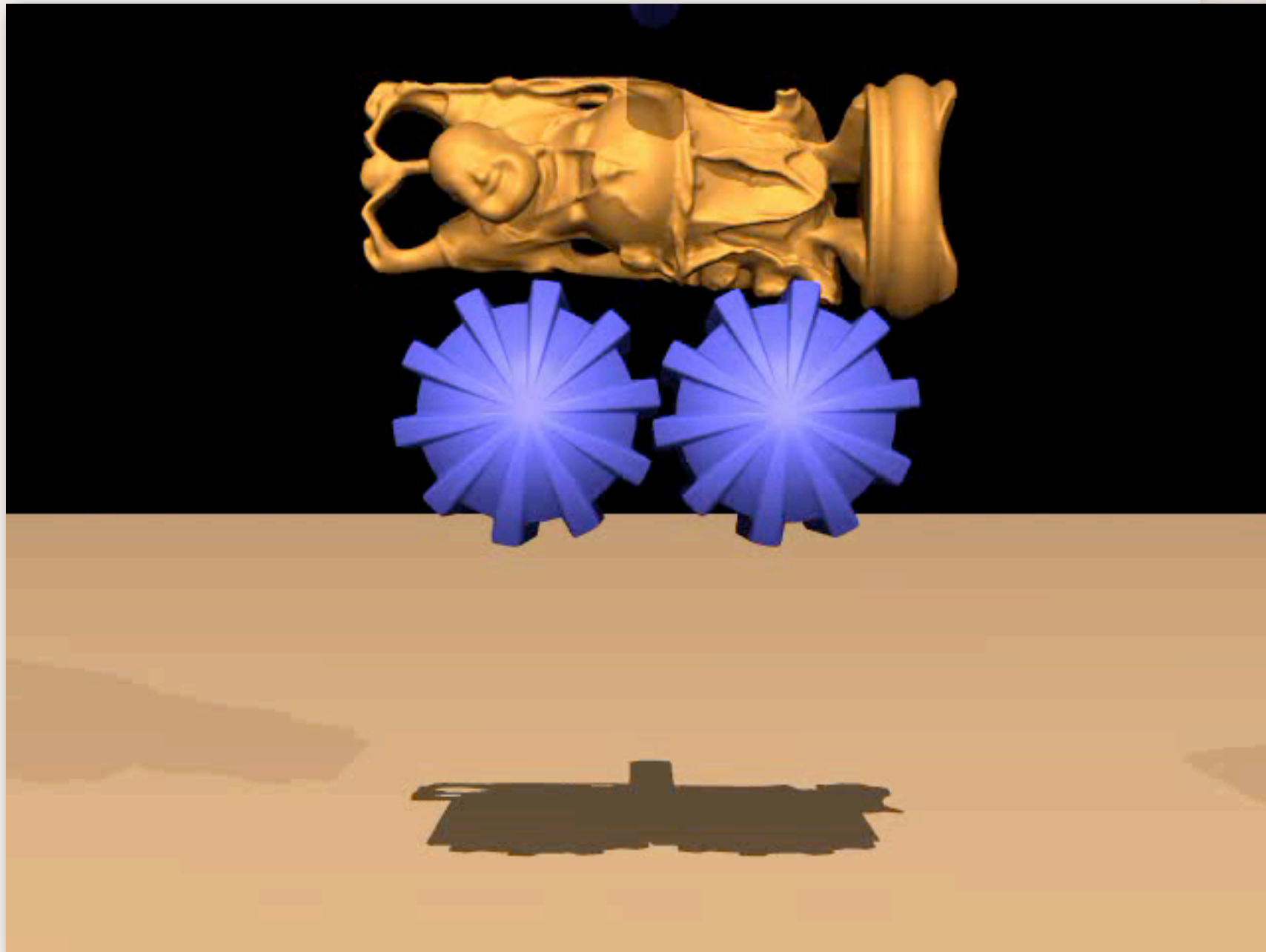


Deformable Materials

Adrien Treuille

Deformable Materials



Taking a Hard Look at Soft Things

Spring with rest length 1:



Deforms by Δx :



$$\text{energy} = E \frac{1}{2} (\Delta x)^2$$

$$\text{force} = - \frac{d \text{ energy}}{dx}$$

$$\text{force} = -E \Delta x$$

Deformations

Spring deformed by Δx :



$$\text{force} = -E \Delta x$$

stress: σ Young's modulus strain: ϵ

Hooke's Law:

$$\sigma = -E \epsilon$$

Steel: $E=10^{11} \text{ N/m}^2$

Rubber: $E=10^7 \sim 10^8 \text{ N/m}^2$



Hooke's Law

$$\sigma = -E\epsilon$$

- **Want to generalize in two ways:**
 - **Continuum Deformations**
 - **3D**



Hooke's Law

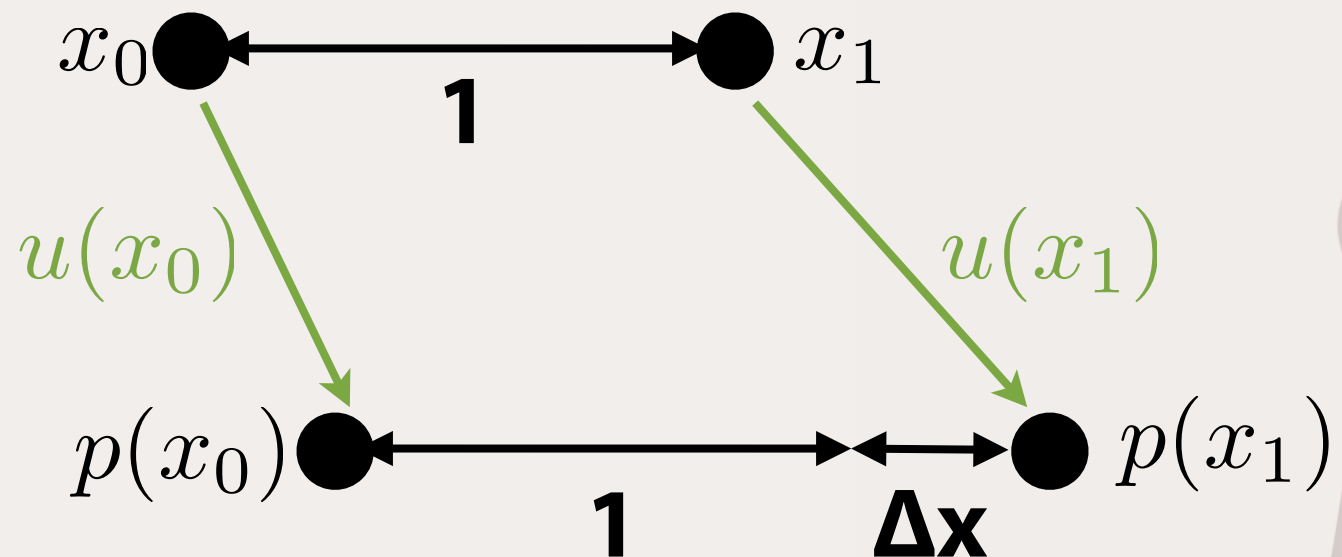
$$\sigma = -E\epsilon$$

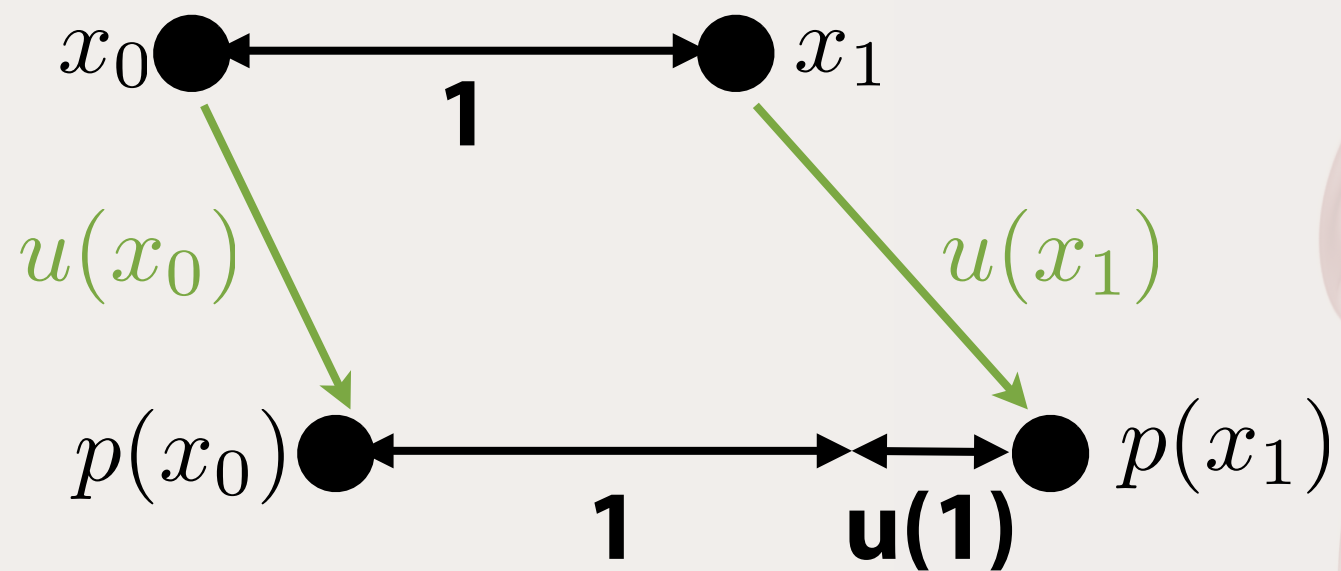
- Want to generalize in two ways:
 - **Continuum Deformations**
 - 3D



Continuum Deformations

- Given a displacement field $p(x)$:
- Which defines a deformation field:
 - $u(x) = p(x) - x$
 - (Like velocities in a fluid.)





- **Suppose: $x_0 = 0$ and $x_1 = 1$**
- **$p(0) = u(0)$**
- **$p(1) = 1 + u(1)$**
- **energy = $\frac{1}{2} E (p(1) - p(0) - 1)^2$**
- **$p(1) \approx 1 + u(0) + \nabla u(0)$**
- **energy $\approx \frac{1}{2} E (1 + u(0) + \nabla u(0) - u(0) - 1)^2$**
- **energy $\approx \frac{1}{2} E \nabla u^2$**
- **force = $-E \nabla u$**

Therefore...

- **force** = $-E \nabla u$

$$\sigma = -E \epsilon$$

In 1D, $\epsilon = \nabla u$.

(Would like to generalize to 3D.)



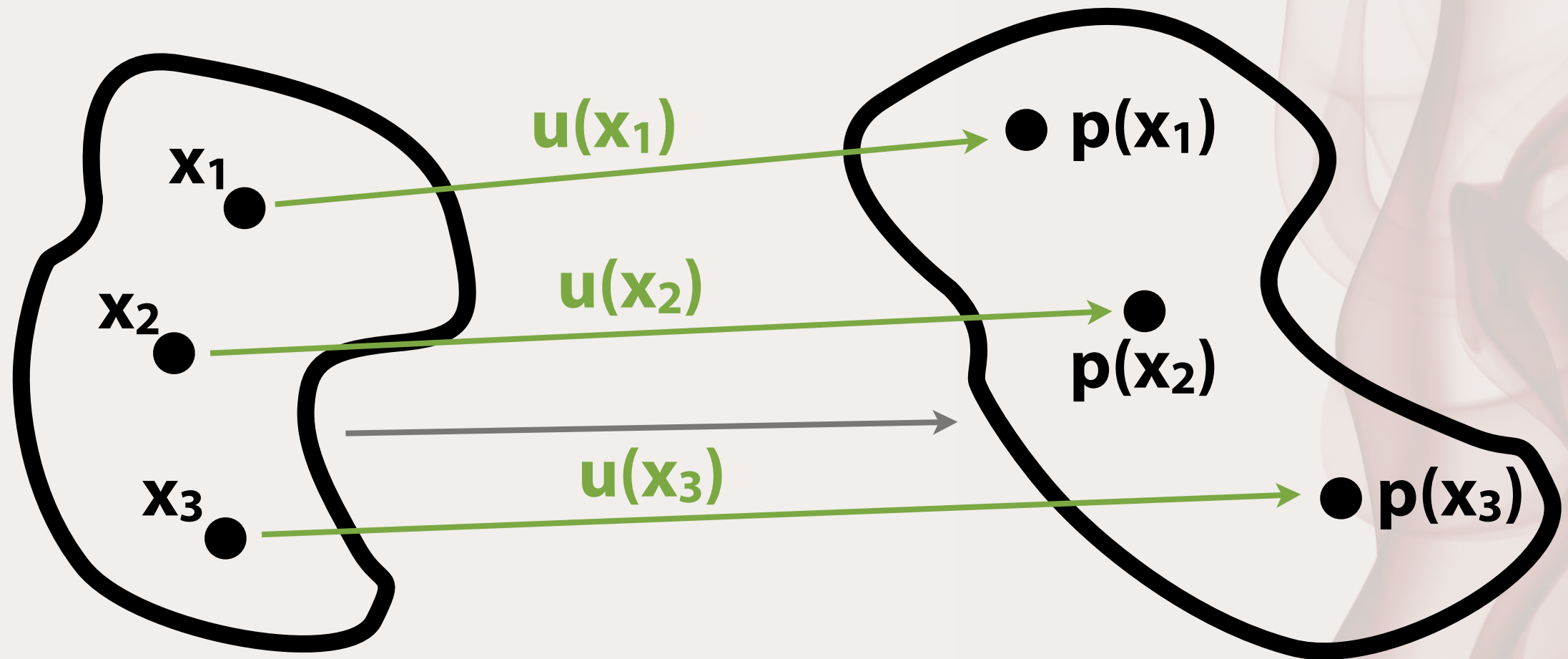
Hooke's Law

$$\sigma = -E\epsilon$$

- Want to generalize in two ways:
 - Continuum Deformations
 - **3D (things will get a little silly)**

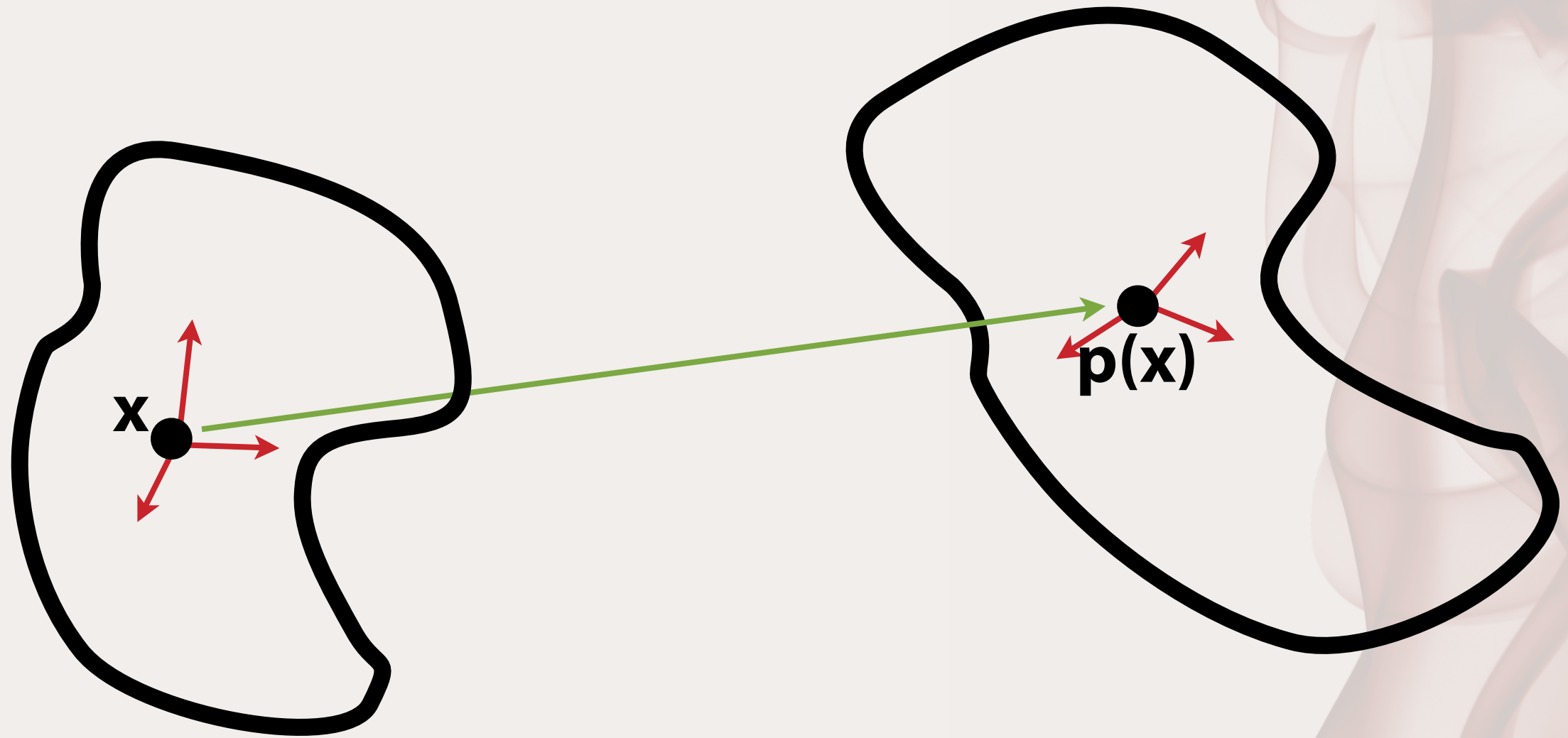


In 3D...



$$u(x) = p(x) - x$$

Coord Sys Transform



point: x

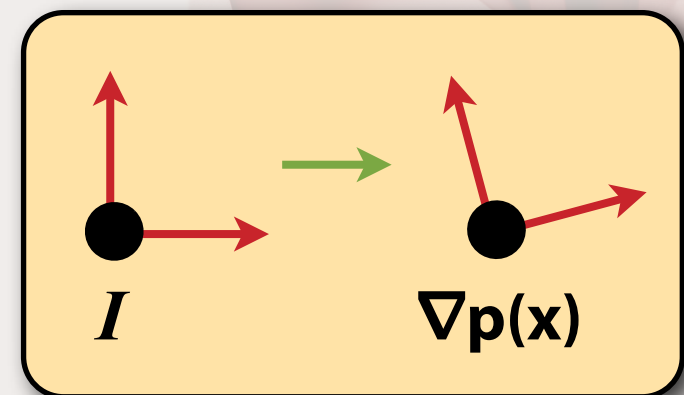
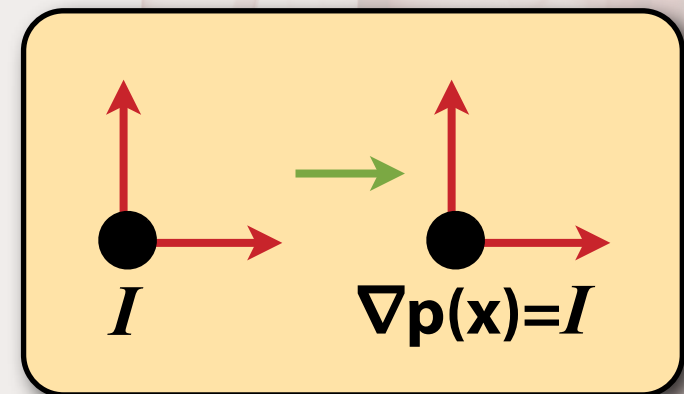
local coordinates: I

point: $p(x)$

local coordinates: ∇p

Defining Strain

- Strain is **invariant to translation**.
 - Ignore $p(x)$
 - Define in terms of local coordinate system transform: $\nabla p(x)$.
- Strain is **invariant to rotation**.
 - If $[\nabla p(x)]^T \nabla p(x) = I$,
 - Then $\epsilon = 0$
- Natural to define strain as:
 - $\epsilon = \frac{1}{2}([\nabla p(x)]^T \nabla p(x) - I)$
 - 6 DOFs



$$\epsilon = \begin{bmatrix} \epsilon_{xx} & \epsilon_{xy} & \epsilon_{xz} \\ \epsilon_{xy} & \epsilon_{yy} & \epsilon_{yz} \\ \epsilon_{xz} & \epsilon_{yz} & \epsilon_{zz} \end{bmatrix}$$

Defining Strain

$$\epsilon = \frac{1}{2} [\nabla p(x)]^T \nabla p(x) - I$$

$$u(x) = p(x) - x$$

$$\nabla u(x) = \nabla p(x) - I$$

$$\epsilon = \frac{1}{2} [\nabla u(x) + I]^T [\nabla u(x) + I] - I$$

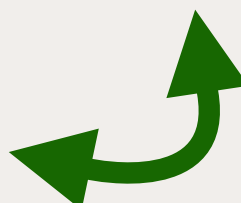
Green's Strain:

$$\epsilon_G = \frac{1}{2} \left(\nabla u + [\nabla u]^T + [\nabla u]^T \nabla u \right)$$

Cauchy's Strain:

$$\epsilon_C = \frac{1}{2} \left(\nabla u + [\nabla u]^T \right) \quad \text{(no rotation)}$$

1D Strain:

$$\epsilon_{1D} = \nabla u$$


Defining Strain

$$\epsilon = \frac{1}{2} [\nabla p(x)]^T \nabla p(x) - I$$

$$u(x) = p(x) - x$$

$$\nabla u(x) = \nabla p(x) - I$$

$$\epsilon = \frac{1}{2} [\nabla u(x) + I]^T [\nabla u(x) + I] - I$$

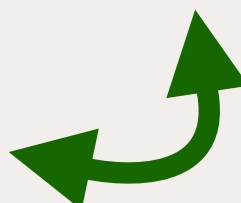
Green's Strain:

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1D Strain:

$$\epsilon_{1D} = \nabla u$$


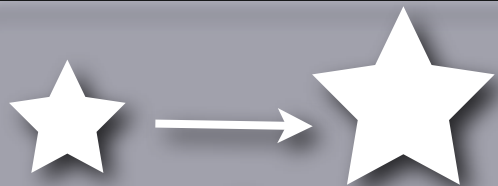
Cauchy's vs. Green's Strain

$$\epsilon_C = \frac{1}{2} \left(\nabla u + [\nabla u]^T \right)$$

Cauchy's Strain

$$\epsilon_G = \frac{1}{2} \left(\nabla u + [\nabla u]^T + [\nabla u]^T \nabla u \right)$$

Green's Strain

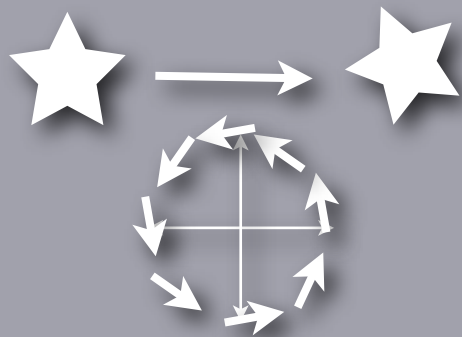


$$u(x) = x$$

$$\nabla u(x) = I$$

$$\begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

$$\frac{3}{2} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$



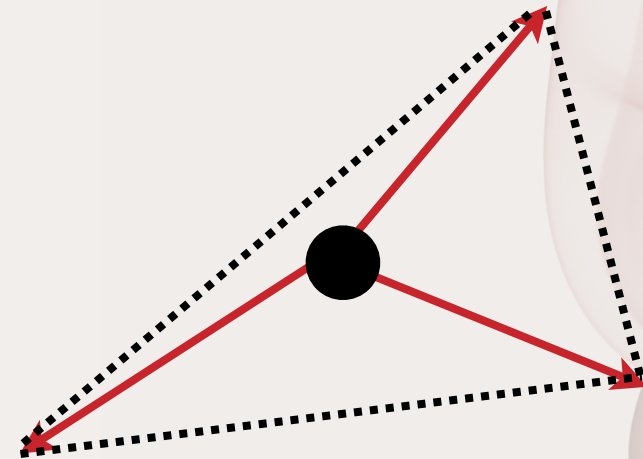
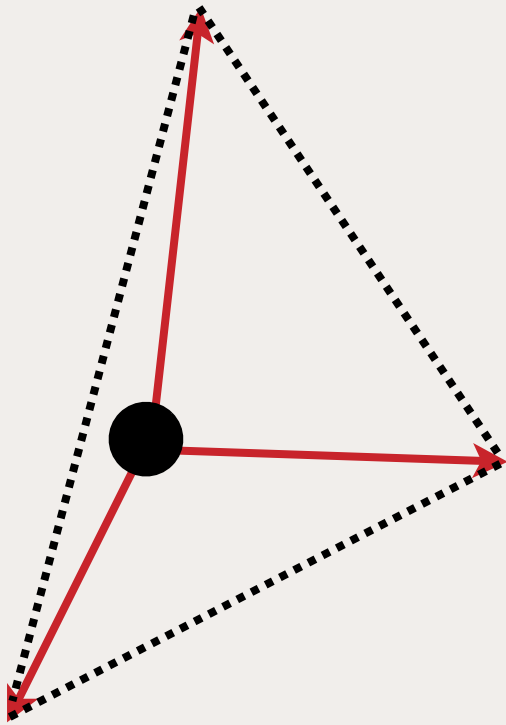
$$\nabla u = \begin{bmatrix} -1 & -1 & 0 \\ 1 & -1 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$

$$\begin{bmatrix} -1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$



$$\begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}$$

Question



Question

- **How could we reduce the cost of simulation for a very finely discretized surface?**
- **Are there cheap ways of getting volumetric behavior without a full tetrahedralization?**
- **How can collision constraints be integrated?**
- **How to simulate plasticity?**



Student Answers

- **How could we reduce the cost of simulation for a very finely discretized surface?**

- make a more coarse tetrahedral mesh, and then track the barycentric coordinates of the original mesh points
 - then, just simulate the coarser mesh!
 - how do we get the coarser mesh:
 - for example, in planar (or low curvature) regions, use larger elements
 - how can we "coalesce" tetrahedra into larger elements
 - what other principles are there?
- if we can precompute (or if we just know) how some object reacts to forces
 - we can use this `_proxy_` instead of the real thing!

- **Are there cheap ways of getting volumetric behavior without a full tetrahedralization?**

- use springs inside the mesh instead of doing the whole stress/strain thing
- forget the interior/ only simulate on the surface!!!
 - ...and maybe we can even simulate the surface too
 - maybe we can use angular springs
- build a "tree" of the structure : the mesh refines itself in areas of deformation

- **How can collision constraints be integrated?**

- is there a local way (looking at deformation gradients) to detect self collisions?
- apply forces to points on the mesh (rather than the whole object itself)
-

- **How to simulate plasticity?**

- non monotonically increasing stress-strain relationship!
 - implementation: if it stresses (gradient of the displacement) too much, stop applying forces!