

# *CONSTRAINT-BASED SIMULATION*

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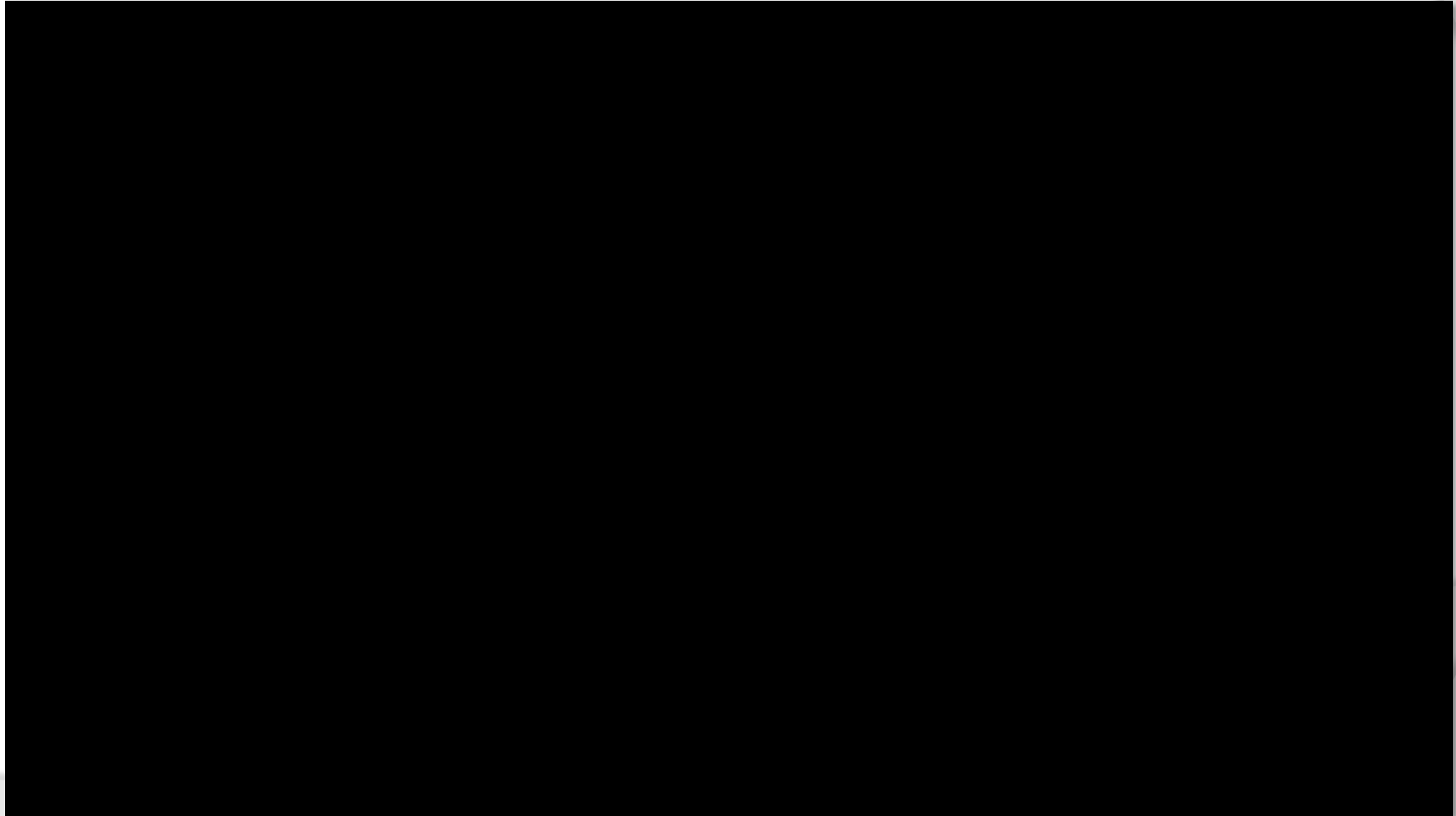
# Content

- ⦿ Motivation
- ⦿ Position-Based Dynamics (PBD)
- ⦿ eXtended Position-Based Dynamics (XPBD)
- ⦿ Deformable Solids
- ⦿ Rigid Bodies
- ⦿ Fluids
- ⦿ Conclusion

# Motivation

- ⦿ Real-time physically-based simulation in visual computing:
  - virtual reality
  - training simulators
  - robotics
  - computer games
  - medical simulation
  - ...
- ⦿ Requirements:
  - high performance
  - stability
  - controllability
  - realistic results (however, high accuracy is often secondary)

# Motivation



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# Constraints

- The motion of a body can be restricted by constraints of the form

$$\mathbf{C}(\mathbf{x}) = 0,$$

where  $\mathbf{x}$  contains all particle positions.

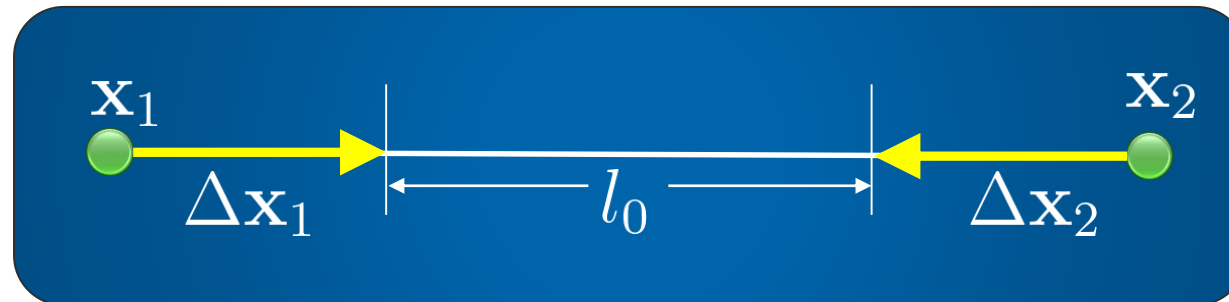
- PBD enforces constraints by position changes.

# Constraints - Example

- Distance constraint:

$$\mathbf{C}(\mathbf{x}) = \|\mathbf{x}_1 - \mathbf{x}_2\| - l_0 = 0$$

- Position changes:

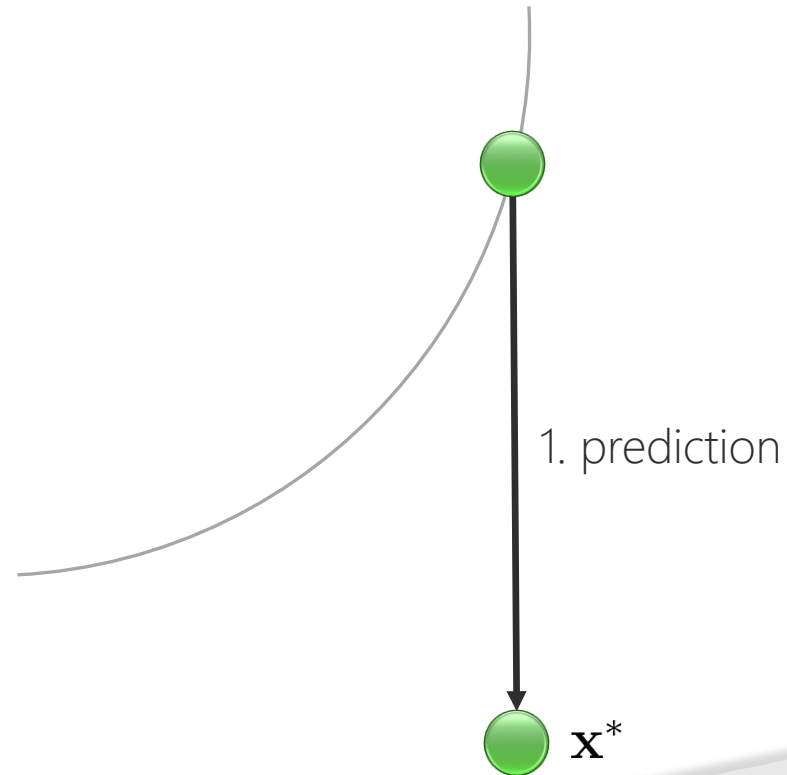


# PBD-Algorithm

1. Prediction step:
  - Time integration only considering external forces to get candidate positions  $\mathbf{x}^*$ .
2. Constraint projection:
  - Check if constraints are fulfilled, otherwise modify candidate positions:
$$\mathbf{x}(t + \Delta t) = \text{constraintProjection}(C(\mathbf{x}), \mathbf{x}^*)$$
3. Velocity update:
  - Update velocities according to position changes to get a consistent system.

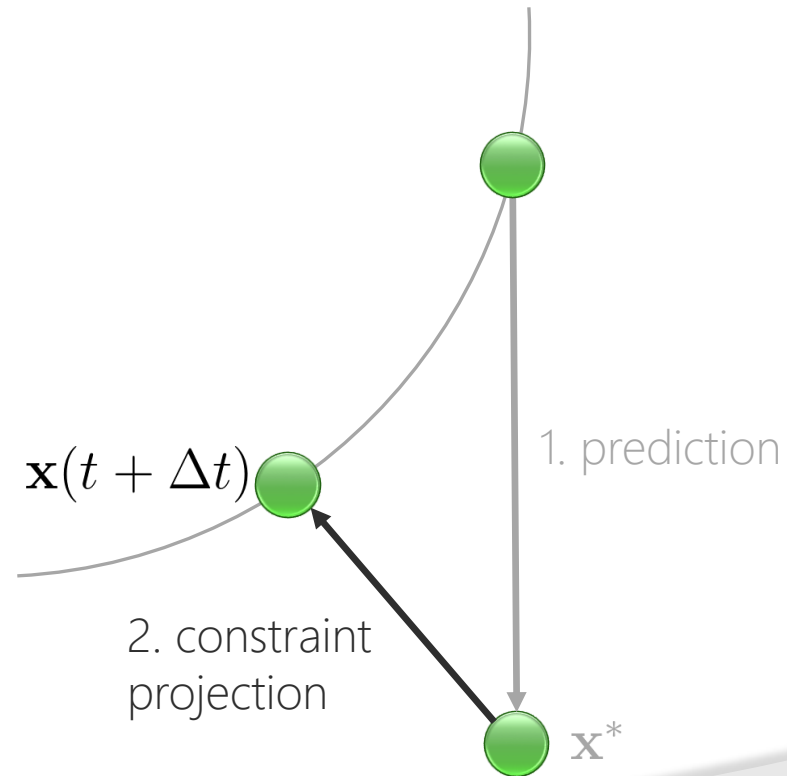
# Example

- Position correction for a particle on a circle



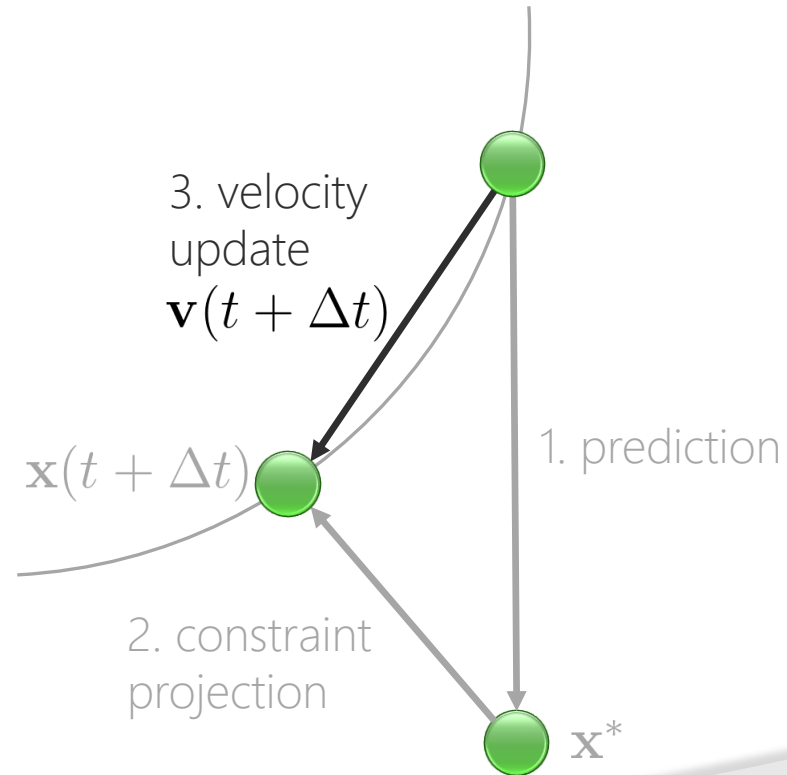
# Example

- Position correction for a particle on a circle



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- Position correction for a particle on a circle

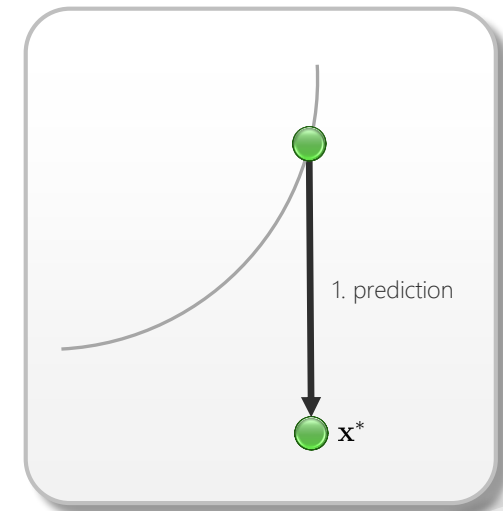


# 1. Prediction Step

- Symplectic Euler integration to get candidate positions:

$$\mathbf{v}^* = \mathbf{v}^n + \frac{\Delta t}{m} \mathbf{F}^{\text{ext}}$$

$$\mathbf{x}^* = \mathbf{x}^n + \Delta t \mathbf{v}^*$$

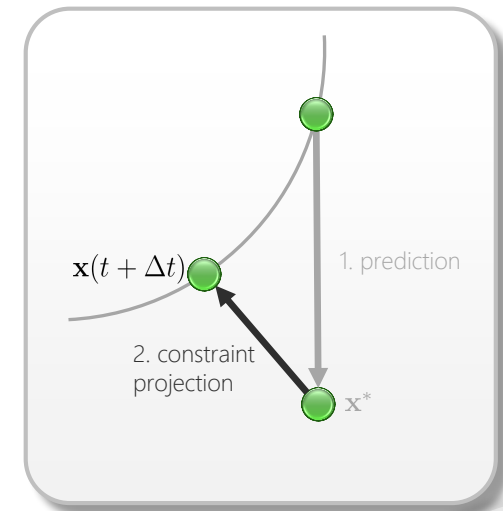


## 2. Constraint Projection

- Linearize the constraint by a first-order Taylor approximation:

$$\mathbf{C}(\mathbf{x}^* + \Delta\mathbf{x}) \approx \mathbf{C}(\mathbf{x}^*) + \mathbf{J}\Delta\mathbf{x} = 0$$
$$\mathbf{J} = \nabla\mathbf{C}^T$$

- However, the resulting linear system is underdetermined.



# Constraint Projection

- ⦿ Solution: we restrict position correction to be in the direction of the gradient (constraint space):

$$\Delta \mathbf{x} = \mathbf{J}^T \boldsymbol{\lambda}$$

- ⦿ This guarantees momentum conservation.
- ⦿  $\boldsymbol{\lambda}$  is a Lagrange multiplier which defines the position correction in constraint space.

# Constraint Projection

- Finally, we weight the corrections with the inverse particle masses:

$$\Delta \mathbf{x} = \mathbf{M}^{-1} \mathbf{J}^T \boldsymbol{\lambda}$$

# Constraint Projection

- Substituting this position correction

$$\Delta \mathbf{x} = \mathbf{M}^{-1} \mathbf{J}^T \boldsymbol{\lambda}$$

in

$$\mathbf{C}(\mathbf{x}^*) + \mathbf{J} \Delta \mathbf{x} = 0$$

yields

$$\mathbf{J} \mathbf{M}^{-1} \mathbf{J}^T \boldsymbol{\lambda} = -\mathbf{C}(\mathbf{x}^*)$$

which is the final system of linear equations that has to be solved.

# Interpretation

$$\mathbf{J}\mathbf{M}^{-1}\mathbf{J}^T\boldsymbol{\lambda} = -\mathbf{C}(\mathbf{x}^*)$$

- ⦿ Mass-weighted position correction (CS):  $\boldsymbol{\lambda}$
- ⦿ Mass-weighted position correction (WS):  $\mathbf{J}^T\boldsymbol{\lambda}$
- ⦿ Position correction (WS):  $\mathbf{M}^{-1}\mathbf{J}^T\boldsymbol{\lambda}$
- ⦿ Position correction (CS):  $\mathbf{J}\mathbf{M}^{-1}\mathbf{J}^T\boldsymbol{\lambda}$
- ⦿ Predicted constraint error (CS):  $-\mathbf{C}(\mathbf{x}^*)$

# Example: Distance Constraint

- Constraint function:

$$C(\mathbf{x}_1, \mathbf{x}_2) = \|\mathbf{x}_1 - \mathbf{x}_2\| - l_0$$

- Gradients:

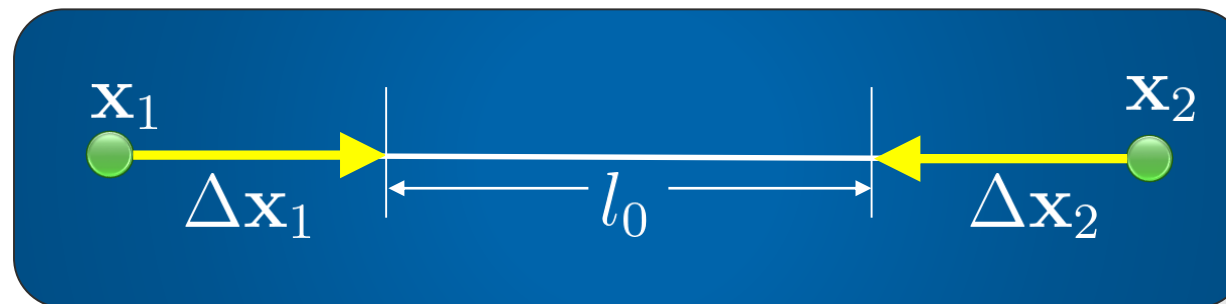
$$\frac{\partial}{\partial \mathbf{x}_1} C(\mathbf{x}_1, \mathbf{x}_2) = \frac{\mathbf{x}_1 - \mathbf{x}_2}{\|\mathbf{x}_1 - \mathbf{x}_2\|}$$
$$\frac{\partial}{\partial \mathbf{x}_2} C(\mathbf{x}_1, \mathbf{x}_2) = -\frac{\mathbf{x}_1 - \mathbf{x}_2}{\|\mathbf{x}_1 - \mathbf{x}_2\|}$$

# Example: Distance Constraint

- This yields the following position corrections:

$$\Delta \mathbf{x}_1 = -\frac{\frac{1}{m_1}}{\frac{1}{m_1} + \frac{1}{m_2}} (\|\mathbf{x}_1 - \mathbf{x}_2\| - l_0) \frac{\mathbf{x}_1 - \mathbf{x}_2}{\|\mathbf{x}_1 - \mathbf{x}_2\|}$$

$$\Delta \mathbf{x}_2 = +\frac{\frac{1}{m_2}}{\frac{1}{m_1} + \frac{1}{m_2}} (\|\mathbf{x}_1 - \mathbf{x}_2\| - l_0) \frac{\mathbf{x}_1 - \mathbf{x}_2}{\|\mathbf{x}_1 - \mathbf{x}_2\|}$$



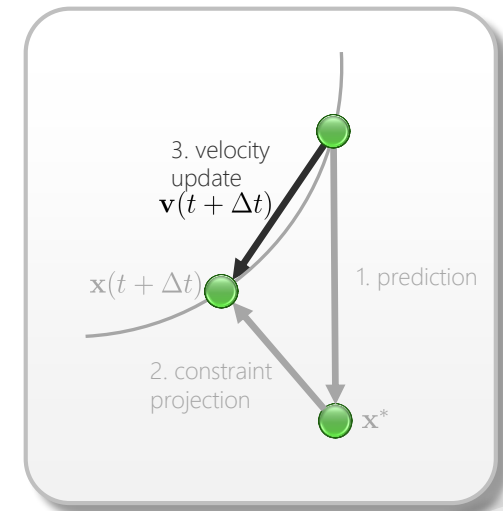
# Constraint Solver

- ⦿ Modified Jacobi iteration
  - In each iteration solve all constraints in parallel
  - Accumulate position corrections
  - Average corrections after each iteration
- ⦿ Gauss-Seidel iteration
  - In each iteration solve constraints one by one
  - Better convergence
  - Atomic operations required for parallelization

# 3. Velocity Update

- Update velocities to get a consistent system as:

$$\mathbf{v}(t + \Delta t) = \frac{1}{\Delta t}(\mathbf{x}(t + \Delta t) - \mathbf{x}(t))$$



# Discussion

- ⦿ Advantages:
  - Fast
  - Stable
  - Simple to implement
- ⦿ Disadvantages
  - Material stiffness depends on number of iterations and time step size

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# eXtended Position-Based Dynamics

- ⦿ XPBD solves the stiffness problem by using a compliant constraint formulation to minimize the energy:

$$E(\mathbf{x}) = \frac{1}{2} \mathbf{C}(\mathbf{x})^T \boldsymbol{\alpha}^{-1} \mathbf{C}(\mathbf{x})$$

where compliance matrix  $\boldsymbol{\alpha}$  corresponds to the inverse stiffness.

# PBD - XPBD

- Position Based Dynamics

$$\mathbf{J}\mathbf{M}^{-1}\mathbf{J}^T\boldsymbol{\lambda} = -C(\mathbf{x}^n)$$

$$\Delta\mathbf{x} = \mathbf{M}^{-1}\mathbf{J}^T\boldsymbol{\lambda}$$

- eXtended Position-Based Dynamics (XPBD)

$$\left(\mathbf{J}\mathbf{M}^{-1}\mathbf{J}^T + \frac{\alpha}{\Delta t^2}\right)\Delta\boldsymbol{\lambda} = -C(\mathbf{x}^n) - \frac{\alpha}{\Delta t^2}\boldsymbol{\lambda}^n$$

$$\Delta\mathbf{x} = \mathbf{M}^{-1}\mathbf{J}^T\Delta\boldsymbol{\lambda}$$

$$\boldsymbol{\lambda}^{n+1} = \boldsymbol{\lambda}^n + \Delta\boldsymbol{\lambda}$$

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# PBD Energy Minimization

- ⦿ To minimize an energy, we define the constraint:

$$C(\mathbf{x}) = E(\mathbf{x})$$

- ⦿ For a continuous model of a deformable solid, we use the strain energy.
- ⦿ In the following we will derive the constraint function and its gradient.

# Deformation

- Deformation

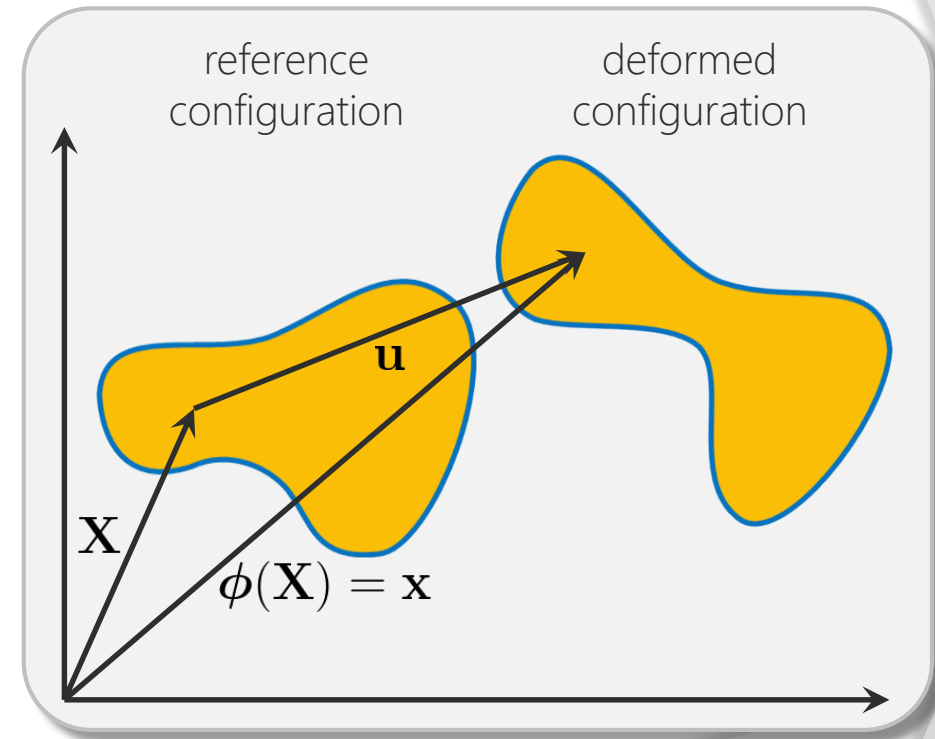
$$\phi(\mathbf{X}) = \mathbf{X} + \mathbf{u} = \mathbf{x}$$

- Deformation gradient

$$\mathbf{F} = \partial\phi(\mathbf{X})/\partial\mathbf{X}$$

- Non-linear Green strain tensor

$$\boldsymbol{\epsilon} = \frac{1}{2} (\mathbf{F}^T \mathbf{F} - \mathbb{1})$$



# Strain Energy

- Energy:

$$E = \int_{\Omega} \Psi d\mathbf{X}$$

- Possible constitutive models of materials:

- Hooke's generalized law

$$\Psi = \frac{1}{2} \boldsymbol{\epsilon} : \mathbf{C} \boldsymbol{\epsilon}$$

- St. Venant-Kirchhoff model

$$\Psi = \frac{\lambda}{2} [\text{tr}(\boldsymbol{\epsilon})]^2 + \mu \text{tr}(\boldsymbol{\epsilon} : \boldsymbol{\epsilon})$$

- Neohookean elasticity

$$\Psi = \frac{\mu}{2} (\mathbf{I}_1 - 3) - \mu \log(J) + \frac{\lambda}{2} \log^2(J)$$

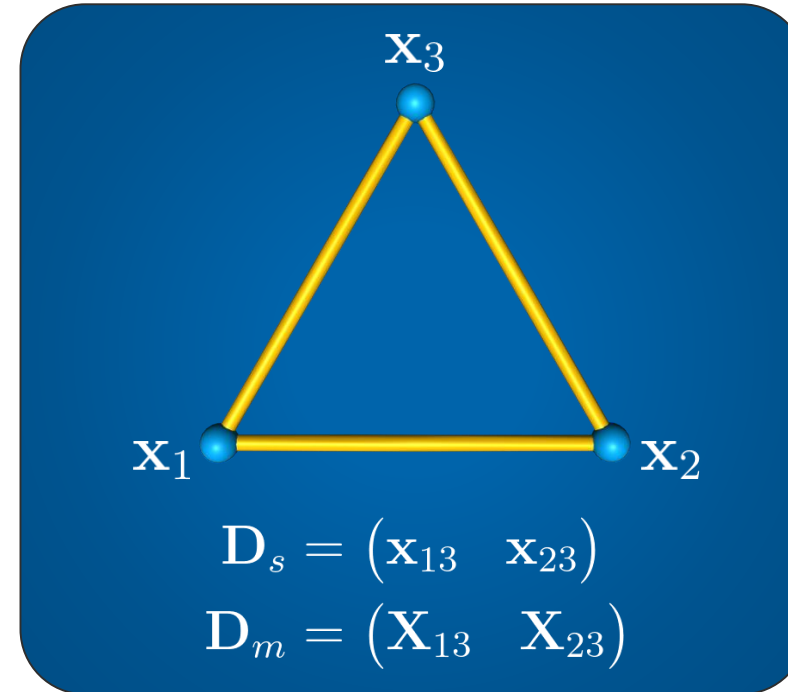
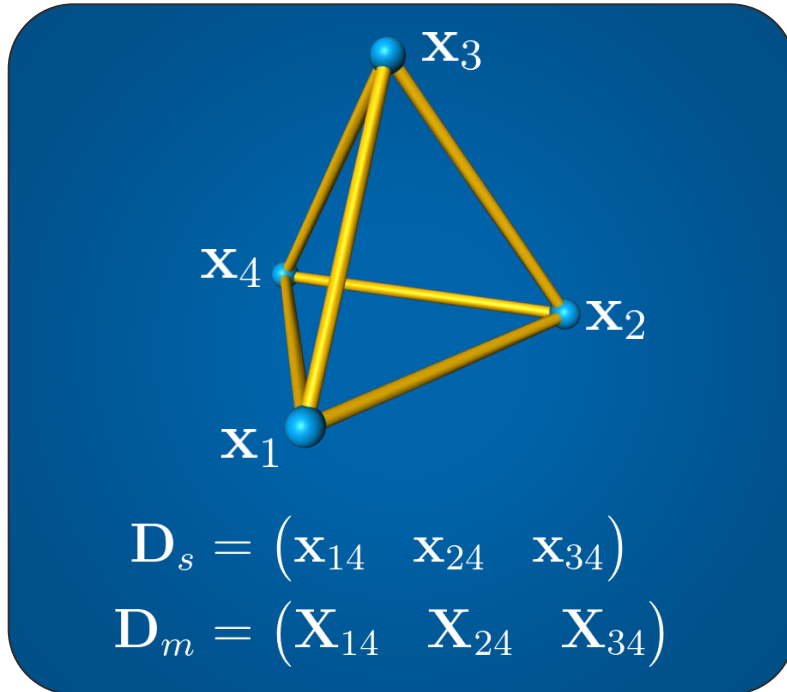
# Strain Energy Gradient

- Strain energy gradient

$$\frac{\partial E}{\partial \mathbf{x}} = \int_{\Omega} \frac{\partial \Psi}{\partial \mathbf{x}} d\mathbf{X} = \int_{\Omega} \mathbf{P}(\mathbf{F}) \frac{\partial \mathbf{F}}{\partial \mathbf{x}} d\mathbf{X}$$

where  $\mathbf{P}(\mathbf{F})$  is the first Piola-Kirchhoff stress tensor.

# Discretization – Linear Elements

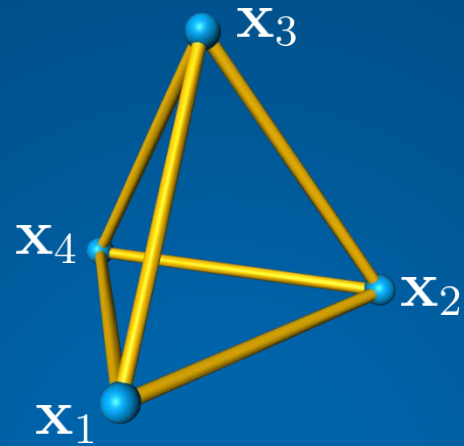


Deformation gradient:

$$\mathbf{F} = \mathbf{D}_s \mathbf{D}_m^{-1}$$

$$\mathbf{x}_{ij} = \mathbf{x}_i - \mathbf{x}_j, \quad \mathbf{X}_{ij} = \mathbf{X}_i - \mathbf{X}_j$$

# Elastic Energy

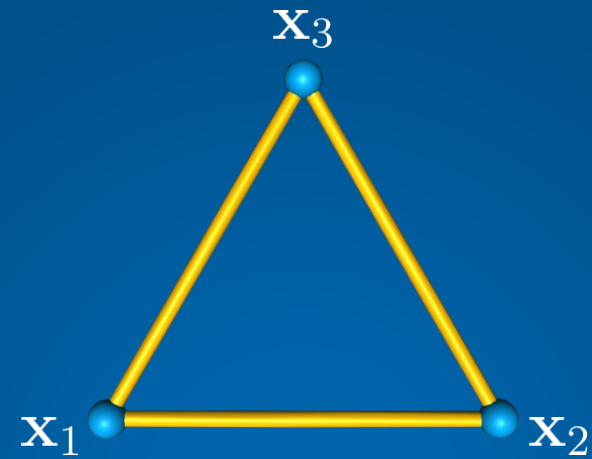


Energy:  $E = V\Psi(\mathbf{F})$

Gradient:

$$\begin{bmatrix} \frac{\partial E}{\partial \mathbf{x}_1} & \frac{\partial E}{\partial \mathbf{x}_2} & \frac{\partial E}{\partial \mathbf{x}_3} \end{bmatrix} = V\mathbf{P}(\mathbf{F})\mathbf{D}_m^{-T}$$

$$\frac{\partial E}{\partial \mathbf{x}_4} = - \sum_{i=1}^3 \frac{\partial E}{\partial \mathbf{x}_i}$$



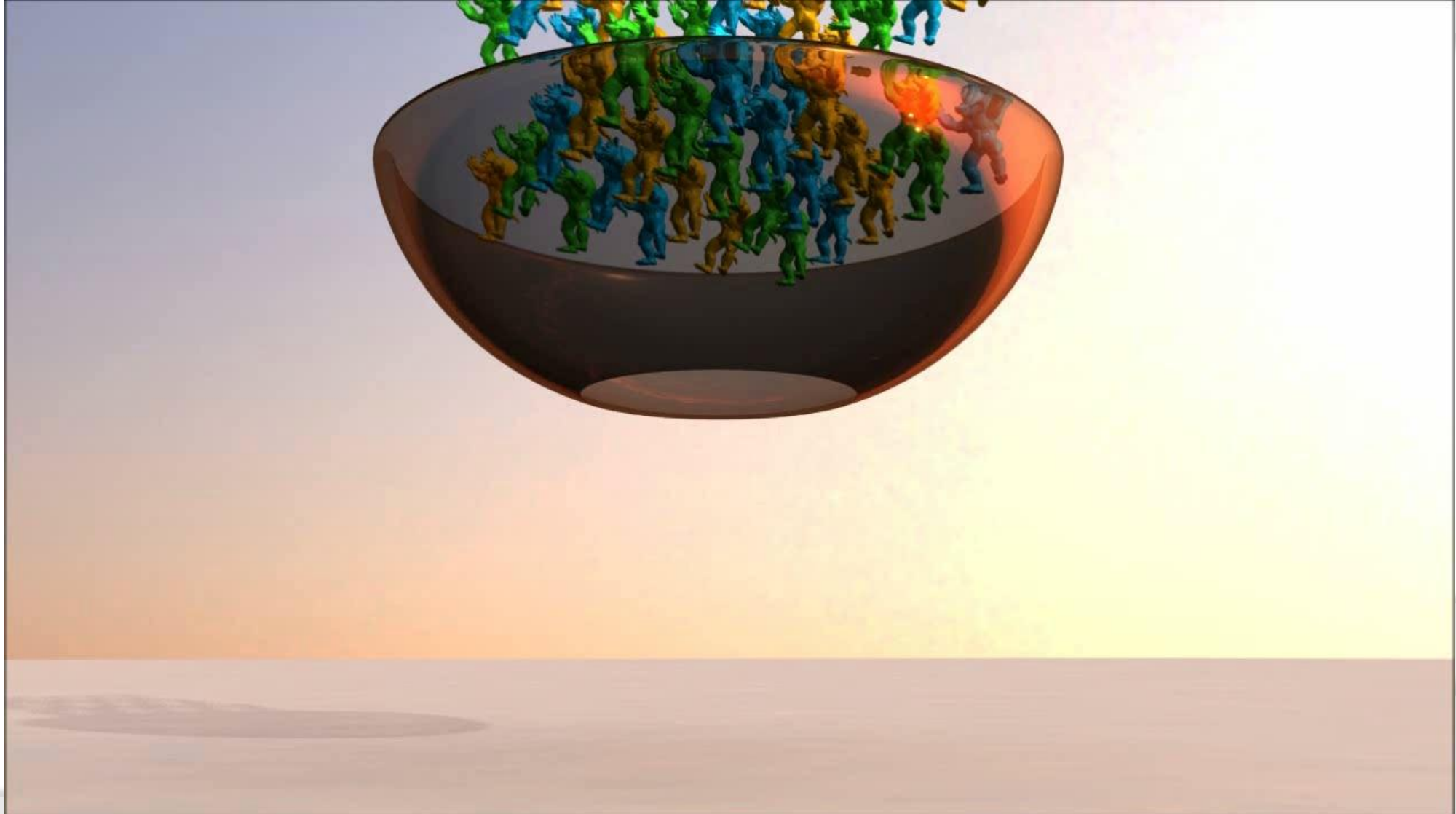
Energy:  $E = A\Psi(\mathbf{F})$

Gradient:

$$\begin{bmatrix} \frac{\partial E}{\partial \mathbf{x}_1} & \frac{\partial E}{\partial \mathbf{x}_2} \end{bmatrix} = A\mathbf{P}(\mathbf{F})\mathbf{D}_m^{-T}$$

$$\frac{\partial E}{\partial \mathbf{x}_3} = - \sum_{i=1}^2 \frac{\partial E}{\partial \mathbf{x}_i}$$

# Results



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# Prediction Step

- Translational DOFs:

$$\mathbf{v}^* = \mathbf{v}^n + \Delta t \frac{\mathbf{F}_{\text{ext}}^n}{m}$$

$$\mathbf{x}^* = \mathbf{x}^n + \Delta t \mathbf{v}^*$$

- Rotational DOFs:

$$\boldsymbol{\omega}^* = \boldsymbol{\omega}^n + \Delta t \mathbf{I}^{-1} (\boldsymbol{\tau}_{\text{ext}}^n - (\boldsymbol{\omega}^n \times (\mathbf{I} \boldsymbol{\omega}^n)))$$

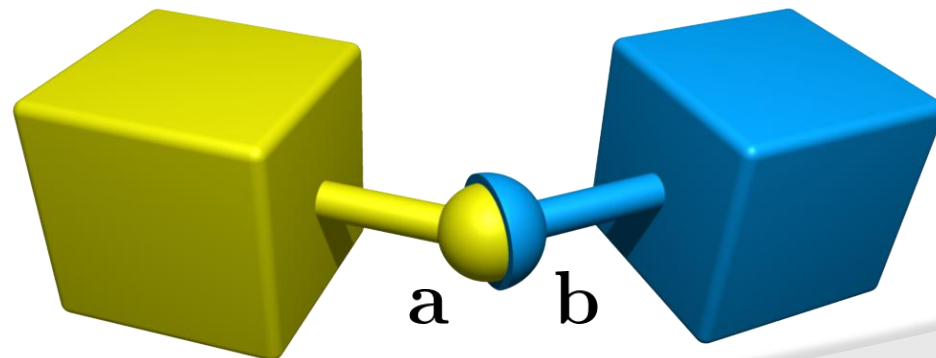
$$\mathbf{q}' = \mathbf{q}^n + \Delta t \frac{1}{2} [0, \omega_x^*, \omega_y^*, \omega_z^*] \mathbf{q}^n$$

$$\mathbf{q}^* = \frac{\mathbf{q}'}{\|\mathbf{q}'\|}$$

# Position Correction

- We split up the constraint in its magnitude  $c$  and its direction  $\mathbf{n}$ .
- Example: ball joint

$$\mathbf{C} = \mathbf{a} - \mathbf{b} = \underbrace{\|\mathbf{a} - \mathbf{b}\|}_c \underbrace{\frac{\mathbf{a} - \mathbf{b}}{\|\mathbf{a} - \mathbf{b}\|}}_{\mathbf{n}} = \mathbf{0}$$

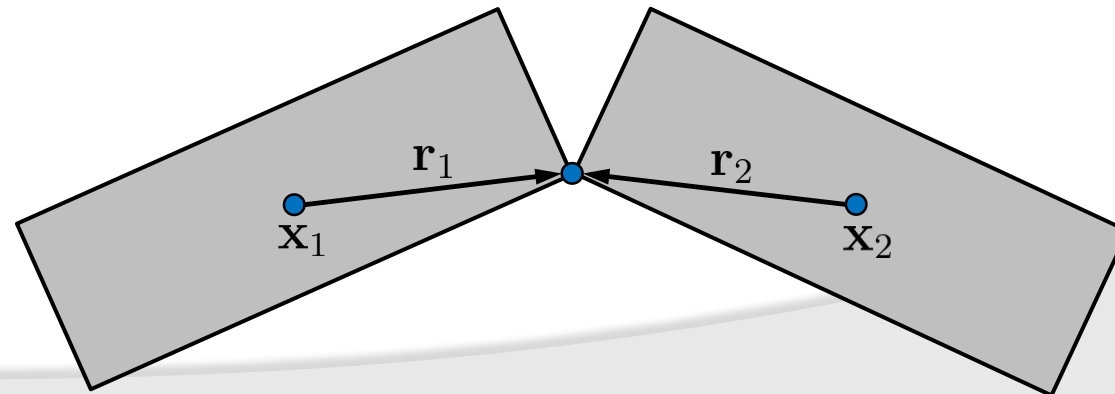


# Position Correction

- To apply the mass weighted correction at the positions  $\mathbf{r}_1$  and  $\mathbf{r}_2$  we need the generalized inverse masses:

$$w_1 = \frac{1}{m_1} + (\mathbf{r}_1 \times \mathbf{n})^T \mathbf{I}_1^{-1} (\mathbf{r}_1 \times \mathbf{n})$$

$$w_2 = \frac{1}{m_2} + (\mathbf{r}_2 \times \mathbf{n})^T \mathbf{I}_2^{-1} (\mathbf{r}_2 \times \mathbf{n})$$



# Position Correction

- ⦿ The Lagrange multiplier is then determined by:

$$\lambda = \frac{-c}{w_1 + w_2}$$

- ⦿ We get the mass weighted position change as:

$$\mathbf{p} = \lambda \mathbf{n}$$

# Position Correction

- Finally, we apply the position change:

$$\mathbf{x}_1^* := \mathbf{x}_1^* + \frac{\mathbf{p}}{m_1}$$

$$\mathbf{x}_2^* := \mathbf{x}_2^* - \frac{\mathbf{p}}{m_2}$$

$$\mathbf{q}_1^* := \mathbf{q}_1^* + \frac{1}{2} (0, \mathbf{I}_1^{-1}(\mathbf{r}_1 \times \mathbf{p})) \mathbf{q}_1^*$$

$$\mathbf{q}_2^* := \mathbf{q}_2^* - \frac{1}{2} (0, \mathbf{I}_2^{-1}(\mathbf{r}_2 \times \mathbf{p})) \mathbf{q}_2^*$$

# Velocity Update

- We update the velocities as:

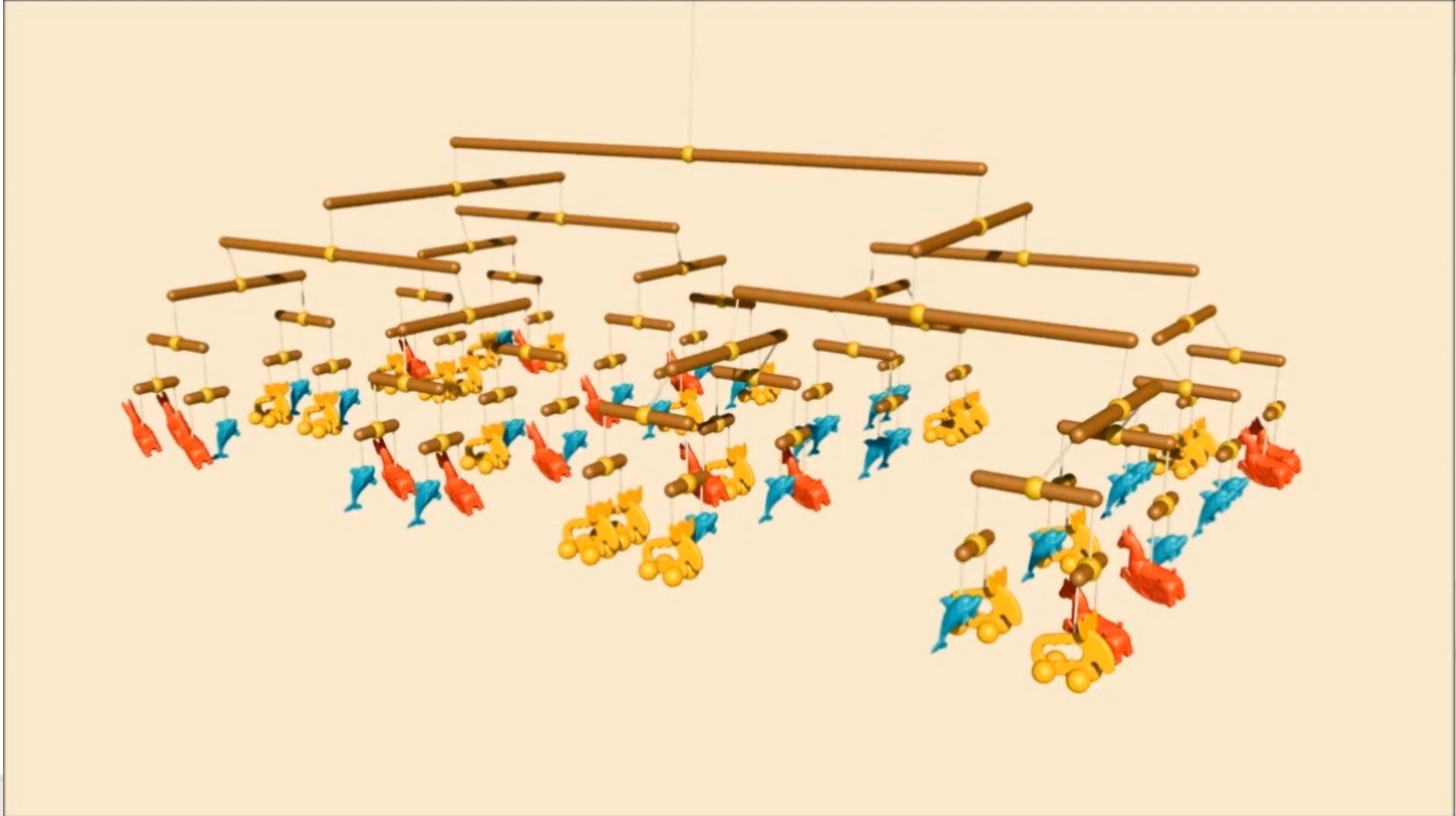
$$\mathbf{v}^{n+1} = \frac{1}{\Delta t} (\mathbf{x}^{n+1} - \mathbf{x}^n)$$

$$\Delta \mathbf{q} = \mathbf{q}^{n+1} (\mathbf{q}^n)^{-1}$$

$$\boldsymbol{\omega}^{n+1} = \frac{2}{\Delta t} (\Delta \mathbf{q}_x, \Delta \mathbf{q}_y, \Delta \mathbf{q}_z)$$

- To avoid a rotation flip we negate  $\boldsymbol{\omega}^{n+1}$ , if  $\Delta \mathbf{q}_w < 0$ .

# Results

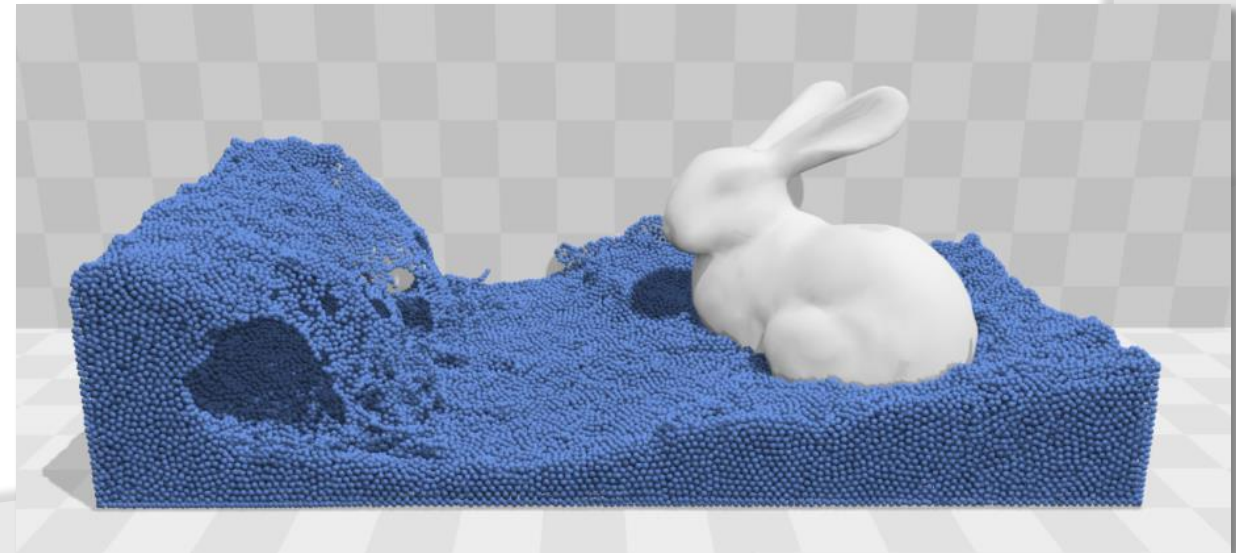
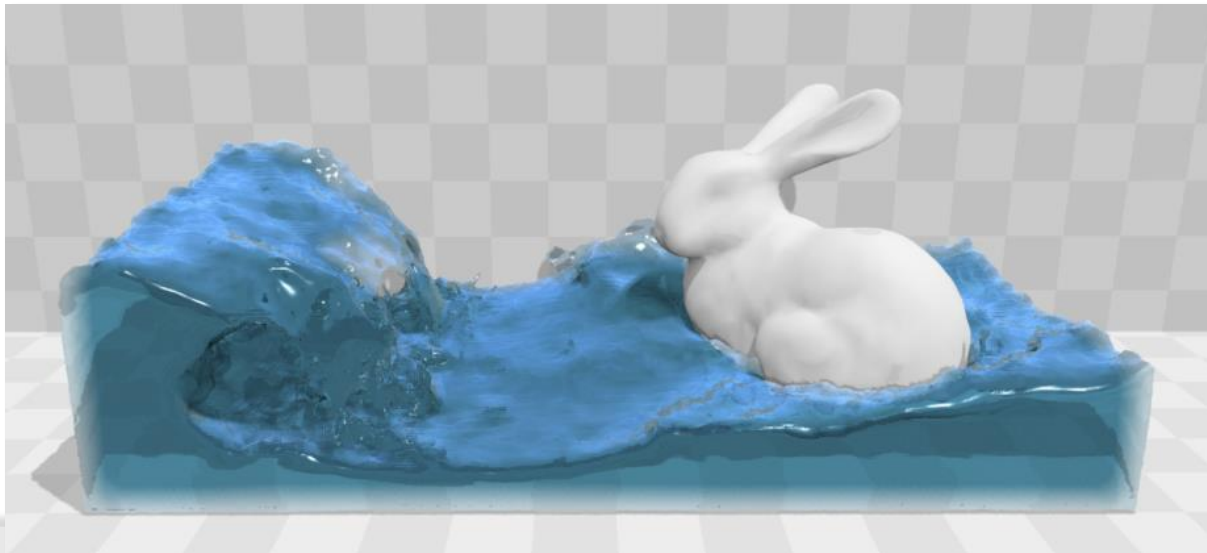


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# Position Based Fluids

- Density constraint:  $C_i = \frac{\rho_i}{\rho_0} - 1 = \frac{1}{\rho_0} \sum_j m_j W_{ij} - 1 = 0$
- Gradient:  $\nabla C_i = \frac{1}{\rho_0} \sum_j m_j \nabla W_{ij}$



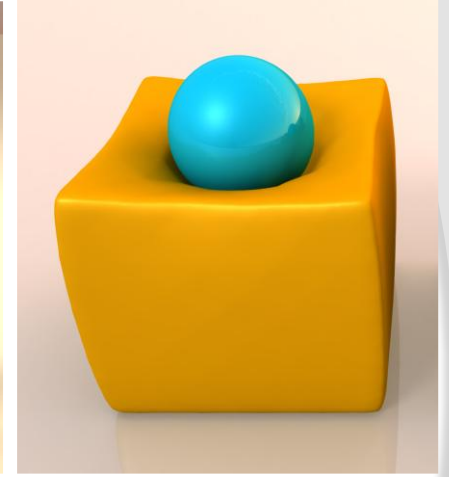
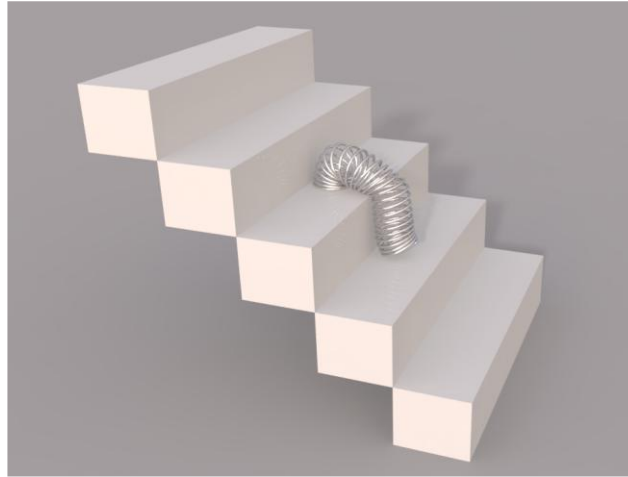
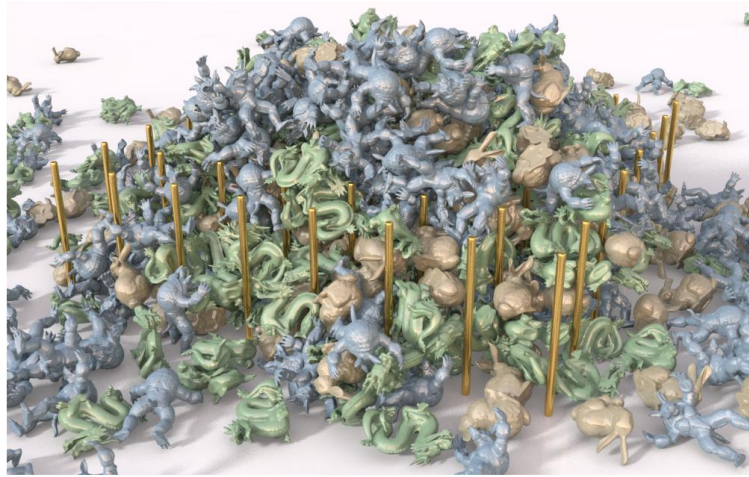
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# Conclusion

- ◎ Position-based methods:
  - are fast, stable and simple to implement,
  - provide a high level of control,
  - can simulate deformable solids (1D, 2D, 3D), multibody systems, fluids and granular materials,
  - support complex physical effects like anisotropy or plasticity and
  - provide visual plausibility.

# Thank you for your attention!



<https://github.com/InteractiveComputerGraphics/PositionBasedDynamics>

