

# Interactions with Granular Material

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# Introduction

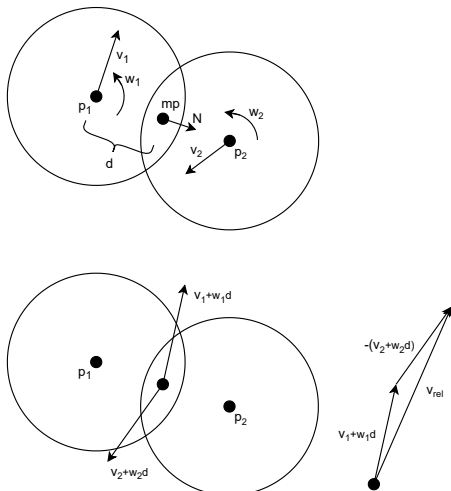
- Simulation of granular materials is important for many reasons
  - ▶ Industrial processes, graphics, geology, etc
- Simulations generally take the following approaches:
  - ▶ *Eulerian grid*-based approaches treat space as a grid of cells, similar to fluid dynamics solvers
    - ★ Key disadvantages: grid structure doesn't preserve matter; inter-granular forces are difficult to model in this paradigm
  - ▶ *Lagrangian particle*-based approaches simulate each individual particle
    - ★ Key disadvantage: computational complexity grows linearly in the number of particles
  - ▶ *Hybrid* approaches, such as *Particle-In-Cell* (PIC), attempt to merge the two
    - ★ Key disadvantages: energy dissipation, instability
    - ★ Key paper: "The Affine Particle-In-Cell Method" [1]

# Particle Types

- Event-Driven (ED)
  - ▶ Can accurately and efficiently simulate continuous particle motion
  - ▶ Properly models particle rigidity and exchange of momentum/energy in collisions
  - ▶ Efficiency decreases when:
    - ★ Simultaneous collisions occur between more than two bodies
    - ★ Collisions are common and contact is prolonged
  - ▶ Unsuitable for sand simulation due to the formation of static piles
- Molecular Dynamics (MD)
  - ▶ Force-based approach that allows slight particle overlap
  - ▶ Simple to program and extend
  - ▶ Downsides:
    - ★ Requires extremely small timesteps to enforce rigidity, otherwise particles will compress
    - ★ Significantly slower than ED
  - ▶ Approach taken in "Particle-Based Simulation of Granular Materials"  
[2]

# Circle Collision Description

- $p_1, p_2$ : Circle centers
- $v_1, v_2$ : Linear velocities
- $w_1, w_2$ : Angular velocities
- $mp$ : Collision midpoint
- $N$ : "Line of centers"  
(normalized  $p_2 - p_1$ )
- $d$ : Distance from center to collision (lever arm length)
- $w_1 d, w_2 d$ : shorthand for rotational velocity  $d$  units away from center with angular velocity  $w$  (at  $mp$ )
- $v_{rel}$ : relative velocity at  $mp$
- Force calculation uses amount of overlap  $\xi = 2(r - d)$  and  $v_{rel}$



# A Third Approach

- Physics engines work similarly to ED, but with discrete timesteps
- Rather than solving each collision as an event, they iteratively resolve constraints (such as non-interpenetration) on each frame
- An excellent short series on physics simulation is "Physically Based Modeling" [3]




Live demo of four approaches:

- Single-particle MD
- Single-particle rigid
- Polysphere MD
- "Bouncy" rigid polyspheres

Characteristics to look for:

- (In)compressibility
- Nonzero Angle of Repose
- Stability/Oscillations
- *Anisotropic force chains*

## Further Reading

-  C. Jiang, C. Schroeder, A. Selle, J. Teran, and A. Stomakhin, “The affine particle-in-cell method,” *ACM Trans. Graph.*, vol. 34, no. 4, Jul. 2015. [Online]. Available: <https://doi.org/10.1145/2766996>
-  N. Bell, Y. Yu, and P. Mucha, “Particle-based simulation of granular material,” 01 2005, pp. 77–86.
-  A. Witkin and D. Baraff, “Physically based modeling,” Available at <https://graphics.pixar.com/pbm2001/> (2001).