

Adam Mally CIS 700-006 Spring 2017 University of Pennsylvania

Advanced Character Physics

- Much of these slides' content is based on the paper "<u>Advanced Character Physics</u>" by Thomas Jakobsen
- Discusses algorithms developed for IO Interactive's game "Hitman: Codename 47"
 - Released in November of 2000!
 - Ragdoll physics
 - Cloth sim
 - Rigid body collisions

Particle Systems

- The basis for most physics simulations is the **particle system**
- Easy to compute the motion of simple points in space
- <u>A fun demonstration</u>

Particle Systems: Euler integration

- Particles usually have three core attributes
 - Position
 - Velocity
 - Acceleration
- We use a repeating update function to simulate the motion of our particles:

- $v' = v + a \cdot \Delta t$
- The equations above are simple <u>Euler</u> <u>integration</u>

Particle Systems: Verlet integration

- Euler integration is rather unstable when our time step is too large
- What if we tried Verlet integration?
 - Velocity is implicit since we store our previous position p*

$$p' = 2p - p^* + a \cdot \Delta t^2$$

p* = p

• More difficult for position and velocity to become desynchronized

Particle Systems: Verlet integration

$$p' = 2p - p* + a \cdot \Delta t^2$$

p* = p

- The logic behind this becomes clearer if we rewrite it slightly:
- $p' = p + (p p^*) + a \cdot \Delta t^2$
- The statement within the parentheses is an approximation of the particle's current velocity, with the Δt already multiplied in

Particle Systems: Runge-Kutta Integration

- Pronounced (sort of) "roon-geh koot-tuh"
 - They're German last names
 - We're not disassembling a ladder
- A method that can be used to numerically solve differential equations, provided we can define our equation in the form:

y' = f(t, y) $y(t_0) = y_0$

Particle Systems: Runge-Kutta Integration

- A bit more complex to write out than Verlet or Euler integration, but is numerically stable for larger time steps
 - Useful if you want to simulate something in real time

$$y_{n+1} = y_n + h/6(k_1 + 2k_2 + 2k_3 + k_4)$$

$$k_1 = f(t_n, y_n)$$

$$k_2 = f(t_n + h/2, y_n + k_1 \cdot h/2)$$

$$k_3 = f(t_n + h/2, y_n + k_2 \cdot h/2)$$

$$k_4 = f(t_n + h/2, y_n + k_3 \cdot h)$$

- h is our time step
- We're evaluating the slope of our function at four points along our time interval: k₁ is essential Euler integration, k₂ and k₃ are midpoint slopes, and k₄ is the slope at the end of the interval

Handling collisions

- We have methods for moving our particles around in a vacuum
- How can we make particles interact with everything else in the scene?
 - Let's not worry about particle-particle interaction just yet

Handling collisions

- We just project our particles back out of the colliding objects
- Translate them along the normal vector of the surface through which they moved
- What will be the effective motion of our particles when they collide with surfaces?
 - Euler integration case?
 - Verlet integration case?

Handling collisions

- We just project our particles back out of the colliding objects
- Translate them along the normal vector of the surface through which they moved
- What will be the effective motion of our particles when they collide with surfaces?
 - Euler integration case?
 - Verlet integration case?
- They'll slide along the surfaces they hit, generally

Handling collisions: Example

- What if we want to force our particles to stay inside a 1000x1000 box?
- How might we code this?

Handling collisions: Example

- What if we want to force our particles to stay inside a 1000³ box?
- How might we code this?
- for(Particle p : particles)

p = min(max(p, vec3(0)), vec3(1000))

Particle Constraints

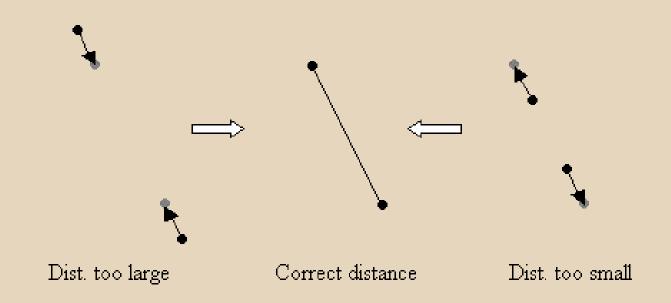
- What if we force certain particles to stay within certain distances of one another?
- Let's say we have two particles, and we want them to always remain 100 units away from each other

 $\circ |p_1 - p_2| = 100$

- If we integrate as before, it's likely that the particles will soon break our constraint
- How do we fix this?

Particle Constraints

• We just move both particles closer if they're too far away, and push them away if they're too close



Particle Constraints

- The end result is as though we've attached a spring with infinite stiffness to the particle pair
 - Visually, the particles are *always* 100 units away
 - We fix their positions between frames
- But what if we want to impose additional constraints on our particles?
 - Let's say they both have to fit inside a box with side lengths at least 100 units long
- We may repeatedly invalidate one constraint, then the other

Particle Constraints: Relaxation

- Let's just repeatedly solve for our constraints until they are both satisfied
 while(!c1.satisfied() && !c2.satisfied()) { //Force particles to be 100 units away //Force particles to fit in the box
 }
- Seems naïve, but it will converge to the desired result
- Known as Jacobi iteration, or just "relaxation"

Particle Constraints: Relaxation

- Can improve relaxation by making it adaptive based on the amount of state change applied by the previous loop of relaxation
- Can change to a FOR loop to stop it early to avoid long or infinite iteration
 - May not be exactly physically correct, but should be close enough
- What happens if we make our springs less stiff?

- Let's say we have a collection of particles where each particle is attached to at least two others via springs
- We relax every particle each frame
 As it turns out, we only need to perform one iteration in relaxation
- Hence, our algorithm becomes O(n), which is great for real time purposes

- Let's discuss how to represent springs in code
- We need some sort of constraint structure:

struct Constraint {
 Particle *p1, *p2;
 float restLength;
};

- We also need to write the code that applies our spring constraints to the particles
- Given a constraint c between two particles:

```
Particle& p1 = *(c.p1); Particle& p2 = *(c.p2);
```

```
vec3 diff = p2 - p1;
```

```
float dist = sqrt(diff.x^2 + diff.y^2 + diff.z^2); // Costly!
```

```
float diffNorm = (dist - c.restLength) / dist;
```

p1 += diff * 0.5 * diffNorm;

p2 -= diff * 0.5 * diffNorm;

• We use this code on all constraints for N iterations

• We can handle particles of nonuniform mass as well

```
Particle& p1 = *(c.p1); Particle& p2 = *(c.p2);
vec3 diff = p2 - p1;
float dist = sqrt(diff.x^2 + diff.y^2 + diff.z^2); // Costly!
float diffNorm = (dist - c.restLength) /
   (dist * (1/p1.mass + 1/p2.mass));
p1 += diff * 1/p1.mass * diffNorm;
```

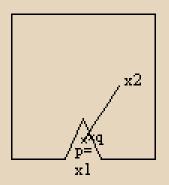
```
p2 -= diff * 1/p2.mass * diffNorm;
```

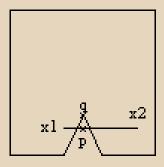
Rigid Bodies

- We can represent rigid bodies as collections of particles constrained by infinitely stiff springs
- This is fine as long as the world is composed of convex objects
- We have to handle concave shapes specially

Rigid Bodies

- We're going to hand wave much of this part now because it is hard! Yay!
- Say we have a line that intersects a concave scene
- We need to know how far it sticks into the collider
- If we know this distance as well as the shortest path to make the line leave the collision, we can translate both endpoints to satisfy our collision constraints





- Joints holding together rigid bodies
 - A pin joint is a single particle that is shared by two rigid bodies
 - A hinge joint is a pair of particles that are shared by two rigid bodies

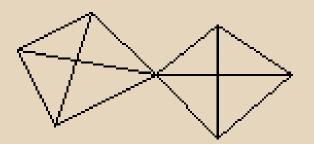


Figure 7a. A pin joint.

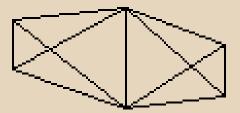


Figure 7b. A hinge.

• How might we simulate an entire human body using particles and constraints?

• How might we simulate an entire human body using particles and constraints?

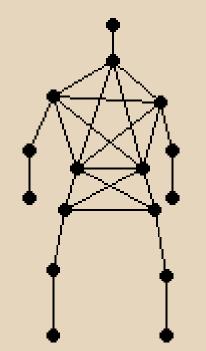


Figure 9. The particle/stick configuration used in Hitman for representing the human anatomy.

 How can we determine the strength of friction that should be applied to an object colliding with a surface?

 How can we determine the strength of friction that should be applied to an object colliding with a surface?

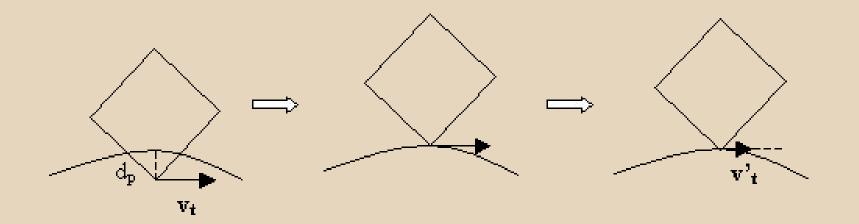


Figure 10. Collision handling with friction (projection and modification of tangential velocity).