

Course Plan



- lec. 1: **Introduction** ●
- lec. 2: **Mathematics** for 3D Games ●●●●●●
- lec. 3: **Scene Graph** ●●
- lec. 4: **Game 3D Physics** ●●● + ●●
- lec. 5: **Game Particle Systems** ●
- lec. 6: **Game 3D Models** ●●
- lec. 7: **Game Textures** ●●
- lec. 8: **Game 3D Animations** ●●●
- lec. 9: **Game 3D Audio** ●
- lec. 10: **Networking** for 3D Games ●
- lec. 11: **Artificial Intelligence** for 3D Games ●
- lec. 12: **Game 3D Rendering Techniques** ●●

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Euler integration methods



init
state

$\mathbf{p} \leftarrow \dots$
 $\vec{\mathbf{v}} \leftarrow \dots$

one
step

$\vec{\mathbf{f}} \leftarrow \text{fun}(\mathbf{p}, \dots)$
 $\vec{\mathbf{a}} \leftarrow \vec{\mathbf{f}} / m$
 $\mathbf{p} \leftarrow \mathbf{p} + \vec{\mathbf{v}} \cdot dt$
 $\vec{\mathbf{v}} \leftarrow \vec{\mathbf{v}} + \vec{\mathbf{a}} \cdot dt$

$t = t + dt$

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Forward Euler *pseudo code*



```
Vec3 position = ...  
Vec3 velocity = ...  
  
void initState() {  
    position = ...  
    velocity = ...  
}  
  
void physicsStep( float dt )  
{  
    Vec3 acceleration = compute_force( position ) / mass;  
    position += velocity * dt;  
    velocity += acceleration * dt;  
}  
  
void main() {  
    initState();  
    while (1) do physicsStep( 1.0 / FPS );  
}
```

Equivalent to...

$$\begin{aligned}\vec{f}_i &= \text{function}(p_i, \dots) \\ \vec{a}_i &= \vec{f}/m \\ \vec{v}_{i+1} &= \vec{v}_i + \vec{a}_i \cdot dt \\ p_{i+1} &= p_i + \vec{v}_i \cdot dt\end{aligned}$$

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Forces

...

$$\vec{f} = \text{function}(p, \dots)$$

...



- Forces are often a function of current positions
 - But not always
- Examples:
 - Gravity
 - Constant, near the surface of a planet
 - But, function of positions in a space simulation
 - Wind
 - Depends on the area exposed in the wind direction
 - Electrical / magnetic forces
 - Archimede's buoyancy
 - Depends on the weight of the submerged volume
 - Mechanical springs
 - simple model: hooke's law – see later
 - shock waves (explosions)
 - Fake / "Magic" control forces
 - added for controlling the evolution of the system, not physically justified

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Forces

$$\vec{f} = \text{function}(\mathbf{p}, \dots)$$



- Forces are often a function of current positions
 - Not always
- Real-world forces can be modelled by things that aren't "forces":
 - Frictions
 - In reality: a force in the opposite direction of motion
 - Its magnitude is proportional to speed (\vec{f} is a function of $\dot{\mathbf{p}}$: difficult to solve!)
 - Can be modelled with velocity drag / damp (see later)
 - Impacts & other violent things
 - In reality: very short, very strong forces
 - Duration $\ll dt$
 - Must be modelled with impulses (see later)
 - Resistance forces
 - E.g.: what prevents your computer to fall through the table
 - E.g.: what prevents a pencil to contract when you push it on the paper
 - In reality, an internal force that contrast an external force (such as gravity)
 - Necessary to model "rigid bodies" and solid bodies
 - Must be modelled by positional constraints (see later)

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Forces: control forces



- Example: the player pressing the forward button
 \Rightarrow a forward force is applied to his/her avatar
 - no physical justification
 - "Don't ask questions, physics engine"
- According to many:
it's better when that's not done much
 - the more physically justified the forces, the better
 - for example: does the car accelerate...
because a **torque** is applied to its two traction wheels VS
because a **force** is applied to its body
 - usually much harder to control
 - see also: gameplay VS cosmetics, control VS realism,
emerging behaviours

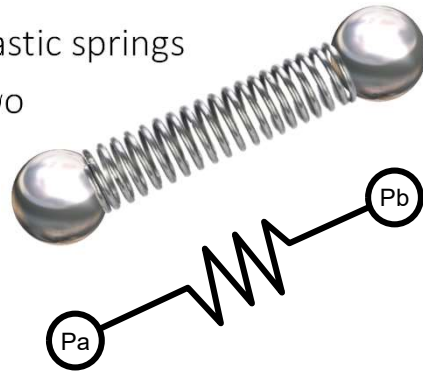
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Forces: Springs (Hooke's law)

- Simplified model for elastic springs
- One spring connects two particles in \mathbf{p}_a and \mathbf{p}_b
- Characterized by:
 1. Rest length ℓ
 2. Stiffness k
- Spring force:

counteracts stretching and compression

$$\vec{f}_a = k(\ell - \|\mathbf{p}_b - \mathbf{p}_a\|) \frac{\mathbf{p}_b - \mathbf{p}_a}{\|\mathbf{p}_b - \mathbf{p}_a\|}$$

$$\vec{f}_b = -\vec{f}_a$$


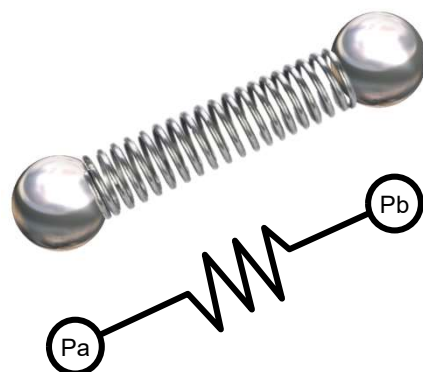
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Forces: springs friction

- A dissipative force
 - Damping factor k_D
- Wants to slow down elongation / shrinking

$$\hat{d} = \frac{\mathbf{p}_b - \mathbf{p}_a}{\|\mathbf{p}_b - \mathbf{p}_a\|}$$

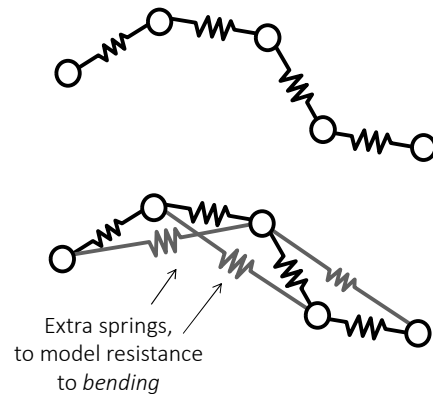
$$\vec{f}_a = k_D(\hat{d} \cdot (\vec{v}_b - \vec{v}_a)) \hat{d}$$



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Mass and Spring systems

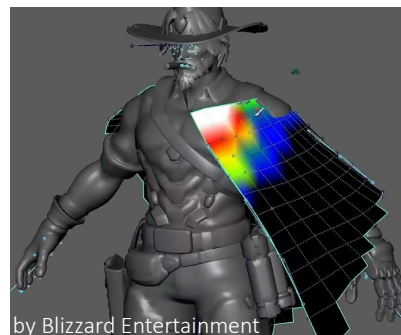
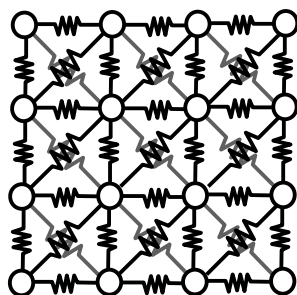
- Useful for deformable objects
- for instance: elastic ropes (or hairs)



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Mass and Spring systems

- For instance: cloth



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Mass and Spring systems can model...



- Elastic deformable objects (aka “soft bodies”)
 - Elastic = go back to original shape
 - Easily modelled as compositions of (ideal) springs.
- Plastic deformable objects? (yes, but not easy)
 - Plastic = assume deformed pose permanently
 - Dynamically change rest-length L in response to large compression/stretching, in certain conditions (not easy)
- Rigid bodies / inextensible ropes ? (they can't)
 - Increase spring stiffness? $k \rightarrow \infty$
 - Makes sense, physically, but...
 - Large $k \Rightarrow$ large $\mathbf{f} \Rightarrow$ instability \Rightarrow unfeasibly small dt needed
 - Doesn't work. How, then? see later

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Continuity of pos and vel



- In real Newtonian physics the state (pos and vel) can only change *continuously*
 - No sudden jump!
- In practice, sometimes is useful to artificially break continuity in the simulations
- Discontinuous changes:
 - in positions: “teleports”
 - in velocity: “impulses”
 - (those are not necessary variations justified by forces)

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Dynamics displacements VS kinematic

...

$$\mathbf{p} = \mathbf{p} + \vec{v} \cdot dt$$

...

aka **dynamic**
displacements

(justified by the
physics)

...

$$\mathbf{p} = \mathbf{p} + d\mathbf{p}$$

...

aka **Kinematic**
displacements

just
“teleportation”

direct and discontinuous change of state (position)

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Impulses VS Forces

...

$$\vec{v} = \vec{v} + (\vec{f} / m) \cdot dt$$

...

...

$$\vec{v} = \vec{v} + (\vec{i} / m)$$

...

a discontinuous change of state (velocity)!

- **Forces** (continuous)
 - Continuous application
 - every frame

- **Impulses**
 - Infinitesimal time
 - una tantum

they model very intense but short forces (such as impacts)

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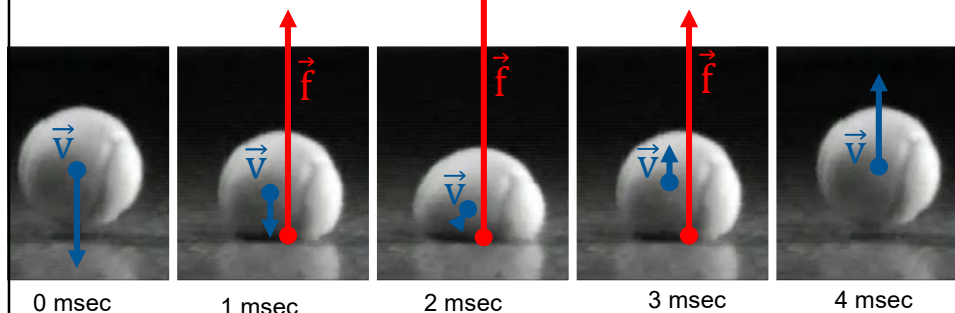
Impulses VS Forces

- **Force** :
 - it determines an **acceleration**
 - **acc** determines a (continuous!) change of **vel**
 - physically correct
- **Impulse** :
 - a (**discontinuous!**) change of **vel**
 - useful to control a simulation (direct change of velocity)
 - a physical interpretation: a force with:
 - application time approaching **zero**
 - magnitude approaching **infinity**
 - Useful to model phenomena with a time scale $\ll dt$
 - ex: a tennis ball rebounding against a tennis racket

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Impulses VS Forces

- what does *truly* happen when it bounces off the ground?



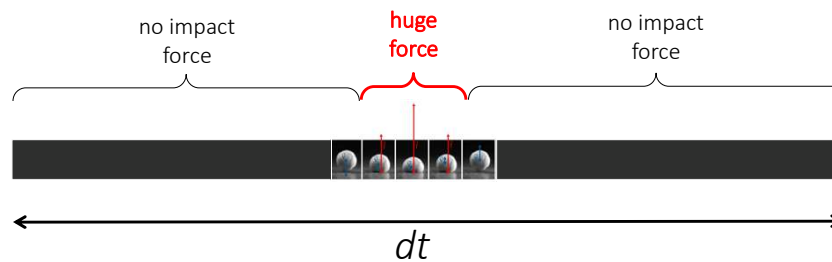
- very strong forces (but not infinite)
- applied for a very short time (but not instantaneous)
- see *collision response* later for details about the impulse based approximations

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Impulses VS Forces



- what does *truly* happen when it bounces off the ground?



- This can only be modelled as an *impulse*, not a force
- See also *collision response*, later

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Effect of integration errors of System Energy



- Because integration errors:
simulated solutions \neq "real" solutions
- In a real system, the total energy cannot increase.
 - Usually, it *decreases* over time, due to dissipations
 - That is, **attrition** turns dynamic energy into heat
- Therefore, a particularly nasty integration error is when the **total energy** of the system *increases* over time
 - e.g.: a pendulum swings faster and faster
- Particularly bad because:
 - Compromises stability
(velocity = big, displacements = crazy, error = crazy)
 - Compromises plausibility
(we can see it's wrong)
- Therefore, a simple way to avoid this:
make sure the simulation always includes **attritions**
 - makes simulation more stable + robust

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Damping VS attrition forces



- We can include attrition as **forces** in our system
 - direction: opposite of current velocity direction
 - magnitude: proportional to a constant, and to speed (speed = magnitude of velocity vector)
 - note: so this force depends on velocity, not just positions.
 - This is the most correct way to model attrition
- Huge simplification: model attrition as “velocity damping”
 - simply, we reduce velocity vectors by a fixed proportion
 - e.g. reduce them all by 2% (drag = 0.02)
 - makes sense!
Higher speed = more attrition = more loss of speed.
Attrition = a “fixed tax” on speed.

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Velocity Damping: the math



- I want to decrease velocity of a percentage for **every second** of (virtual) time
 - e.g.: if 2% then Drag = 0.02
- how should I update velocity for at **every dt** ?

1/FPS sec

$$\vec{v} \leftarrow \vec{v} \cdot (1 - Drag)^{dt}$$

- for small enough Drag, this is well approximated by

$$\vec{v} \leftarrow \vec{v} \cdot (1 - dt \cdot Drag)$$

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Velocity Damping: pseudo-code



```
Vec3 position = ...
Vec3 velocity = ...

void initState(){
    position = ...
    velocity = ...
}

void physicsStep( float dt )
{
    Vec3 acceleration = force( positions ) / mass;
    position += velocity * dt;
    velocity += acceleration * dt;
    velocity *= (1.0 - DRAG * dt);
}

void main(){
    initState();
    while (1) do physicsStep( 1.0 / FPS );
}
```

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Velocity Damping: notes



- Velocity Damping is useful for robustness,
 - avoids energy to increase
- Problems of Velocity Damping
 - tends to exaggerate frictions; even when it makes sense, e.g. in space, no air
 - Crude approximation: attrition forces are not really linear with speed
 - It's attrition with everything...: air, soil.
 - Isotropic force: in reality, attrition force depends of velocity direction
- In practice:
 - low values: hardly noticeable (except in the long run)
 - high values: feels like everything is moving in molasses; (ita: *melassa*) everything quickly grinds to a halt
 - very high values: (e.g. 50% per frame) basically, no inertia anymore (useful to quickly converge to (local) minimal energy states: becomes basically a solver for static problems, not of dynamics)

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Other numerical integrators ("numerical ways to compute integrals")



- Some commonly used alternatives:
 - "Forward" Euler method (the one seen so far)
 - Symplectic Euler method
 - Leapfrog method
 - Verlet method
- These are just variants of each other – let's see them!
 - From the code point of view, no big change
 - They can differ in accuracy / behavior
 - E.g. order of accuracy
 - Note: a more accurate method is also more efficient (larger dt are possible, so fewer steps are necessary)

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Forward Euler Method: limitations



- efficiency / accuracy: not too good
 - error accumulated over time = linear in dt
 - it's only a "first order" method
 - Doubles the steps = halve the dt , only halves the errors (can be better, but no guarantees)
- in practice, scarce stability for large dt
- minor problem: no reversibility, *even in theory*
 - real Newtonian Physics is reversible: flip all velocities and forces \Rightarrow go backward in time.
 - In our simulation (with Euler): this doesn't work exactly
 - Ability to go reverse a simulation would be useful in games! E.g. replays in a soccer game ?
 - Pro tip: basically, reverse time direction never done like this To go backward in time accurately, store states

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Forward Euler

init
state

$\mathbf{p} \leftarrow \dots$
 $\vec{\mathbf{v}} \leftarrow \dots$

one
step

$\vec{\mathbf{f}} \leftarrow \text{fun}(\mathbf{p}, \dots)$
 $\vec{\mathbf{a}} \leftarrow \vec{\mathbf{f}}/m$
 $\mathbf{p} \leftarrow \mathbf{p} + \vec{\mathbf{v}} \cdot dt$
 $\vec{\mathbf{v}} \leftarrow \vec{\mathbf{v}} + \vec{\mathbf{a}} \cdot dt$

$t = t + dt$

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Symplectic Euler

init
state

$\mathbf{p} \leftarrow \dots$
 $\vec{\mathbf{v}} \leftarrow \dots$

one
step

$\vec{\mathbf{f}} \leftarrow \text{fun}(\mathbf{p}, \dots)$
 $\vec{\mathbf{a}} \leftarrow \vec{\mathbf{f}}/m$
 $\vec{\mathbf{v}} \leftarrow \vec{\mathbf{v}} + \vec{\mathbf{a}} \cdot dt$
 $\mathbf{p} \leftarrow \mathbf{p} + \vec{\mathbf{v}} \cdot dt$

$t = t + dt$

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Forward Euler *pseudo code*






Vec3 position = ...	Equivalent to...
Vec3 velocity = ...	$\vec{f}_i \leftarrow \text{function}(p_i, \dots)$
void initState() {	$\vec{a}_i \leftarrow \vec{f}/m$
position = ...	$\vec{v}_{i+1} \leftarrow \vec{v}_i + \vec{a}_i \cdot dt$
velocity = ...	$p_{i+1} \leftarrow p_i + \vec{v}_i \cdot dt$
}	
void physicsStep(float dt)	
{	
Vec3 acceleration = compute_force(position) / mass;	
position += velocity * dt;	
velocity += acceleration * dt;	
}	
void main() {	
initState();	
while (1) do physicsStep(1.0 / FPS);	
}	

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Symplectic Euler *pseudo code* (aka semi-implicit Euler)



Vec3 position = ...	Equivalent to...
Vec3 velocity = ...	$\vec{f}_i \leftarrow \text{function}(p_i, \dots)$
void initState() {	$\vec{a}_i \leftarrow \vec{f}/m$
position = ...	$\vec{v}_{i+1} \leftarrow \vec{v}_i + \vec{a}_i \cdot dt$
velocity = ...	$p_{i+1} \leftarrow p_i + \vec{v}_{i+1} \cdot dt$
}	
void physicsStep(float dt)	
{	
Vec3 acceleration = compute_force(position) / mass;	
velocity += acceleration * dt;	
position += velocity * dt;	 just flip the order
}	
void main() {	
initState();	
while (1) do physicsStep(1.0 / FPS);	
}	

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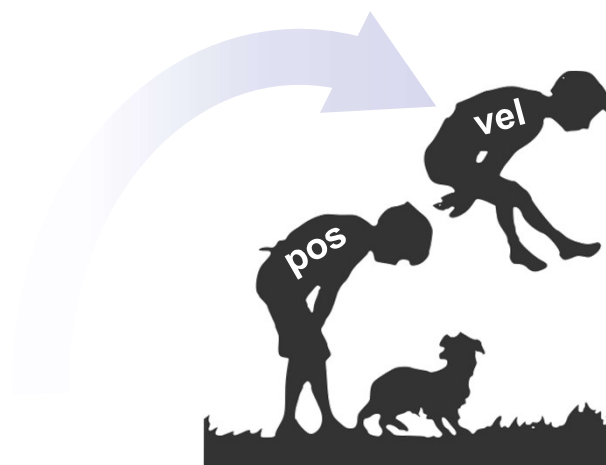
Forward Euler VS Symplectic Euler (warning: over-simplifications)



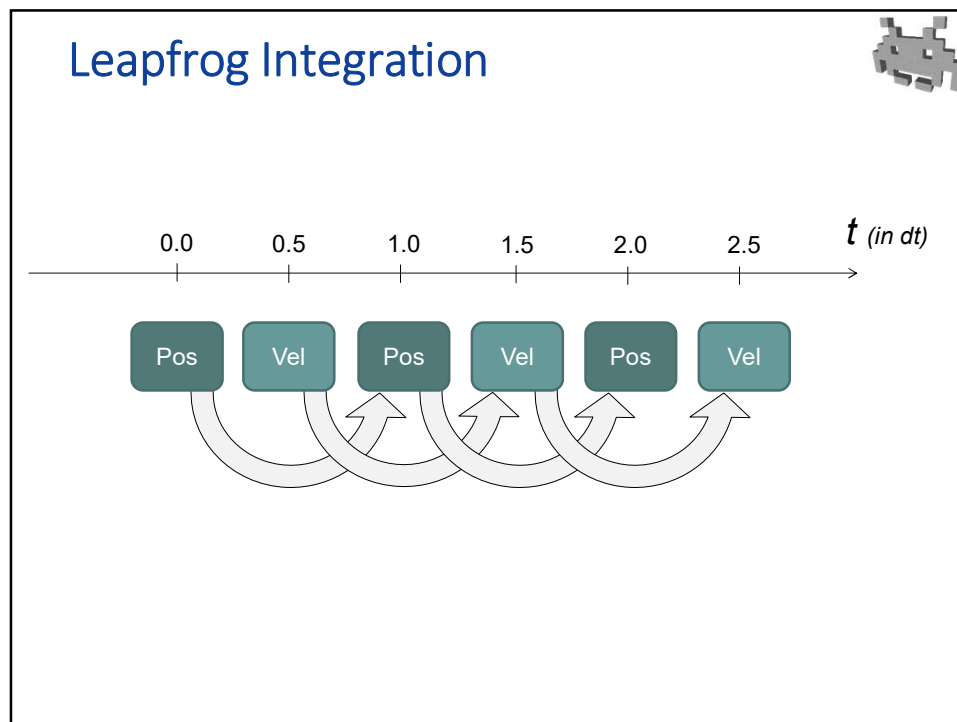
- From the code point of view, they are very similar
- The semantics changes:
 - in Symplectic Euler
the position altered using *next frame* velocity
 - (it's "wrong", in a sense – but works better)
- Similar properties, but better in practice
 - Same order of convergence (still just one ☹)
 - On average, better behavior
 - More stable, more accurate

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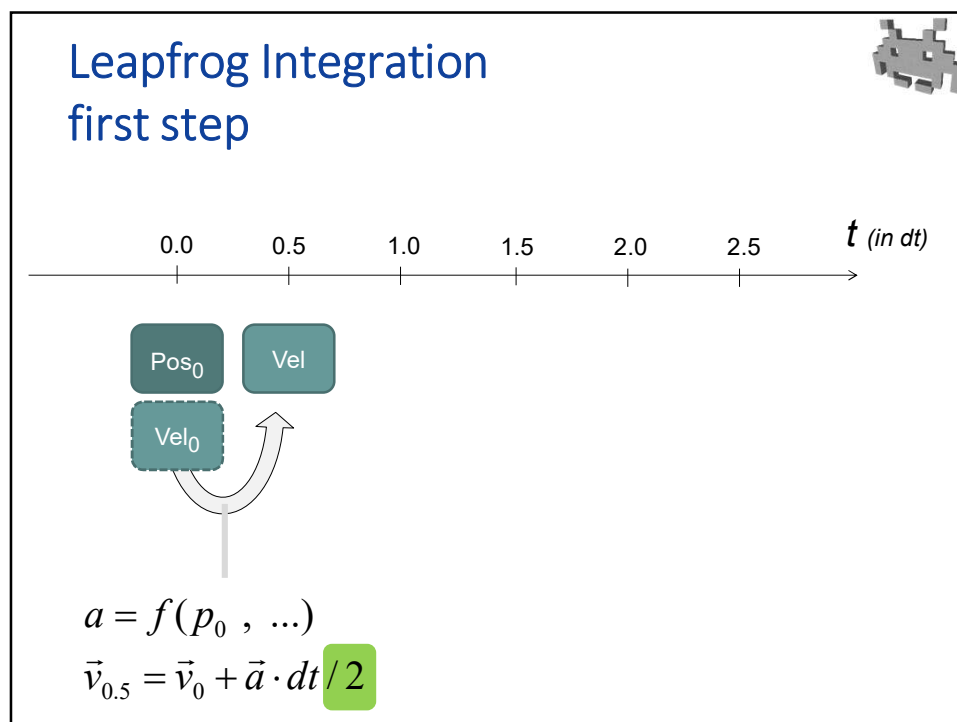
Leapfrog Integration



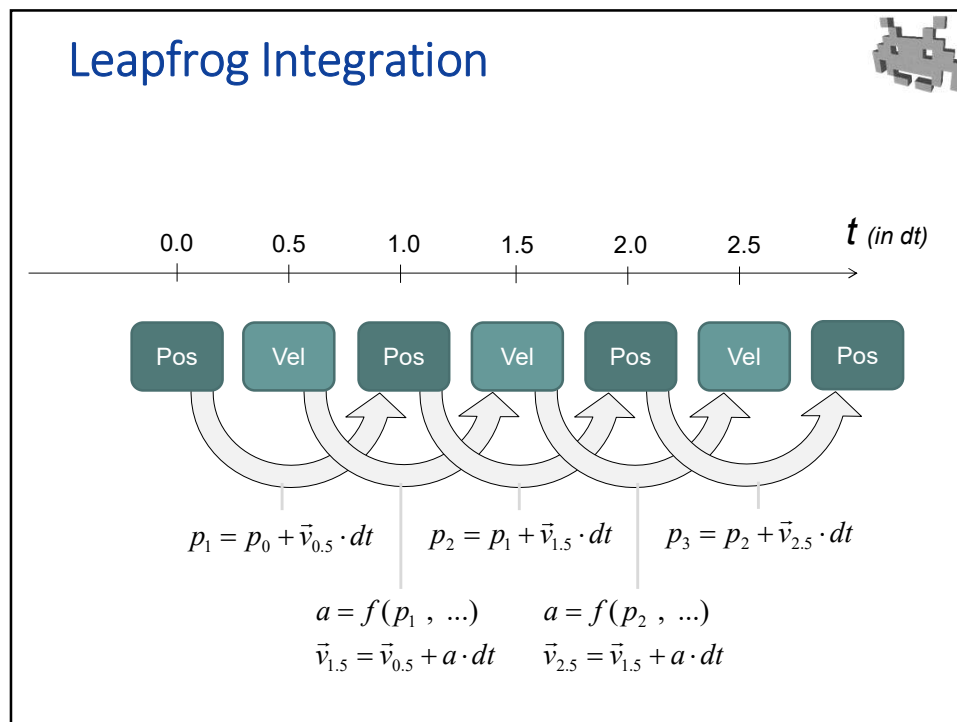
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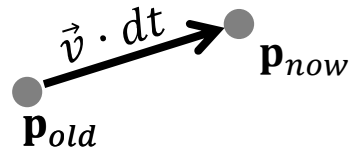
Leapfrog method: pros and cons

- Same cost as Euler – and basically same code
 - Velocity stored in status = velocity “half a dt ago” (and after updating it: “half a frame in the future”)
 - Only real difference: the initialization of speed
- Better theoretical accuracy, for the same dt
 - better asymptotic behavior: it's a second order instead of first!
 - cumulated error: proportional to dt^2 instead of dt
 - error per frame: proportional to dt^3 instead of dt^2
- Bonus: fully reversible!
 - (in theory only. Beware e.g. floating point errors)
- But: requires fixed dt during all the simulation
 - for the theory to work as advertised

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Verlet integration method

- Idea: remove velocity from state
- Current velocity is implicit
- It's defined from:
 - current pos \mathbf{p}_{now}
 - last pos \mathbf{p}_{old}
which we need to record



$$\mathbf{p}_{now} = \mathbf{p}_{old} + \vec{v} \cdot dt \quad \leftarrow \text{Euler \& variants}$$

$$\Rightarrow$$

$$\vec{v} = (\mathbf{p}_{now} - \mathbf{p}_{old})/dt \quad \leftarrow \text{Verlet}$$

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Verlet integration method

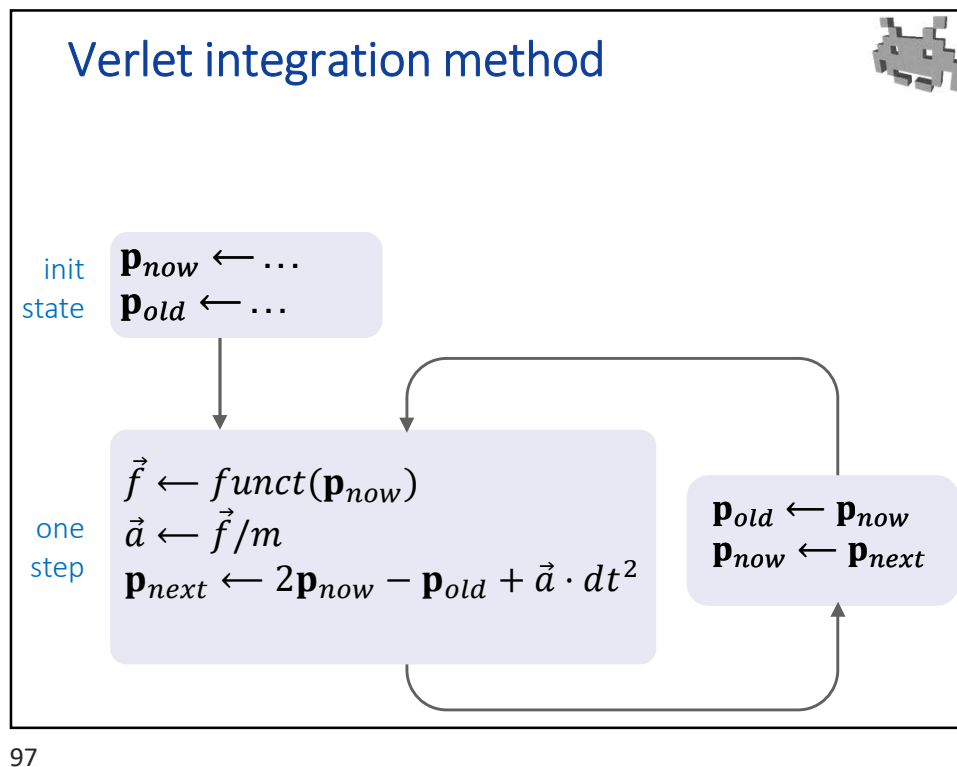
init
state $\mathbf{p}_{now} = \dots$
 $\mathbf{p}_{old} = \dots$

one
step

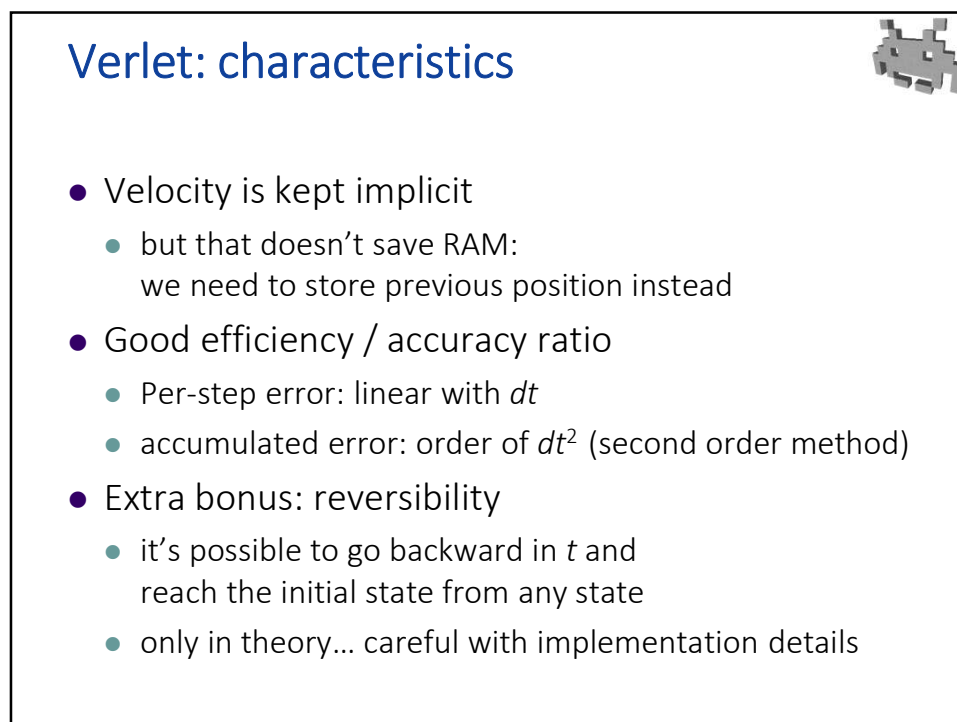
$$\begin{aligned} \vec{f} &= \text{funct}(\mathbf{p}_{now}, \dots) \\ \vec{a} &= \vec{f}/m \\ \vec{v} &= (\mathbf{p}_{now} - \mathbf{p}_{old})/dt \\ \vec{v} &= \vec{v} + \vec{a} \cdot dt \\ \mathbf{p}_{next} &= \mathbf{p}_{now} + \vec{v} \cdot dt \end{aligned}$$

expanding
this...

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Verlet: *caveats*



- ⚠ it assumes a constant dt (time-step duration)
 - if it varies: corrections are needed! (how? – see below)
- ⚠ Q: how to act on **velocity** (which is now implicit)?
 - e.g., how to apply **impulses**
 - A: change \mathbf{p}_{old} instead (how?)
- ⚠ Q: how to act of **positions** w/o impacting velocity?
 - e.g. to apply **teleports** / **kinematic motions**
 - A: displace both \mathbf{p}_{new} and \mathbf{p}_{old} by the same amount
- ⚠ Q: how to apply **velocity damps**?
 - A: act on \mathbf{p}_{old} or \mathbf{p}_{next} (see below)

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Changing the value of dt in Verlet (if it's not constant)



Problem:

if dt now changes to a new dt'
then, all \mathbf{p}_{old} must be updated to some \mathbf{p}'_{old}

Find \mathbf{p}'_{old} : $\vec{v} = (\mathbf{p}_{now} - \mathbf{p}_{old})/dt$ current velocity \vec{v}
 $\vec{v} = (\mathbf{p}_{now} - \mathbf{p}'_{old})/dt'$ and position \mathbf{p}_{now}
 must not change



$$\mathbf{p}'_{old} = \mathbf{p}_{now} \cdot (dt - dt')/dt + \mathbf{p}_{old} \cdot dt'/dt$$

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Velocity damping in Verlet (way 2)

implicit

- Velocity at next frame: $\vec{v} = (\mathbf{p}_{next} - \mathbf{p}_{now})/dt$
- We want to multiply \vec{v} a factor c_{damp}
 - before applying accelerations
 - e.g. 0.98
obtained as $(1 - dt \cdot c_{DRAG})$
- We can do that using a more general formula for \mathbf{p}_{next}

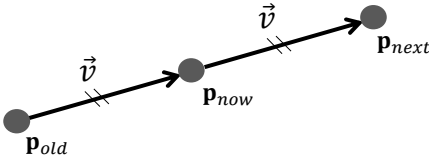
$$\mathbf{p}_{next} = 2 \cdot \mathbf{p}_{now} - 1 \cdot \mathbf{p}_{old} + dt^2 \cdot \vec{a}$$

↓

$$\mathbf{p}_{next} = (1 + c_{damp}) \cdot \mathbf{p}_{now} - c_{damp} \cdot \mathbf{p}_{old} + dt^2 \cdot \vec{a}$$

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Velocity damping in Verlet (way 2) (geometric interpretation)

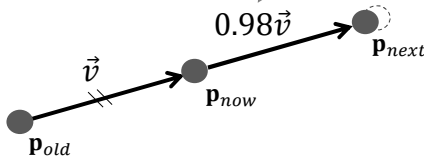


$\mathbf{p}_{next} = 2 \cdot \mathbf{p}_{now} - 1 \cdot \mathbf{p}_{old}$

Equivalently,
 \mathbf{p}_{next} is an **extrapolation**
of $\mathbf{p}_{now}, \mathbf{p}_{old}$:

$\mathbf{p}_{next} = \text{mix}(\mathbf{p}_{old}, \mathbf{p}_{now}, 2)$

a bit shorter



$\mathbf{p}_{next} = 1.98 \cdot \mathbf{p}_{now} - 0.98 \cdot \mathbf{p}_{old}$

Equivalently,
 \mathbf{p}_{next} is a different **extrapolation**
of $\mathbf{p}_{now}, \mathbf{p}_{old}$:

$\mathbf{p}_{next} = \text{mix}(\mathbf{p}_{old}, \mathbf{p}_{now}, 1.98)$

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