## COSC3500 Forum Presentation

Max Bo

22nd of October

#### Outline

I)	escri	nt	ion

Demo

Integration

Barnes-Hut

Correctness

Performance Optimizations

Parallelization

Results

???

## Description

The task was to create a stock-standard, 2-dimensional gravitational *n*-body simulator.

All bodies were to be assumed to be point masses. The simulation was to be accurate, maintaining a constant total energy, and exhibiting phenomena such as apsidal precession.

## Demo

# Integration I

$$F=G\frac{m_1m_2}{r^2}$$

$$a_i = F(x_i)$$
$$v_{i+1} = v_i + a_i \, \Delta t$$

# Integration II

Dehen and Read note that the Euler method 'performs very poorly in practice', further noting that 'errors are proportional to  $\Delta t^2$ '. They contrast it with the second-order *Leapfrog* symplectic integrator, which is 'heavily used in collisionless N-body applications'.

$$x_i = x_{i-1} + v_{i-1/2} \Delta t$$
  
 $a_i = F(x_i)$   
 $v_{i+1/2} = v_{i-1/2} + a_i \Delta t$ 

which only requires a single acceleration calculation per every two half timesteps

# Integration III

and a 'kick-drift-kick' form

$$v_{i+1/2} = v_i + a_i \frac{\Delta t}{2}$$

$$x_{i+1} = x_i + v_{i+1/2} \Delta t$$

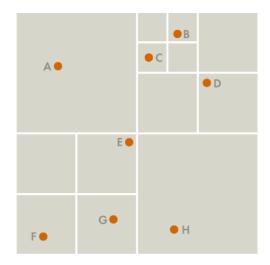
$$v_{i+1} = v_{i+1/2} + a_{i+1} \frac{\Delta t}{2}$$

that is stable with variable timstepping, but incurs an additional acceleration calculation per every two half timesteps.

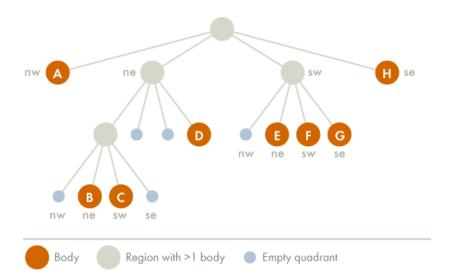
# Integration IV

Runge-Kutta

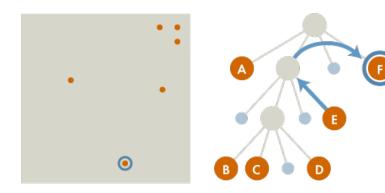
#### Barnes-Hut I



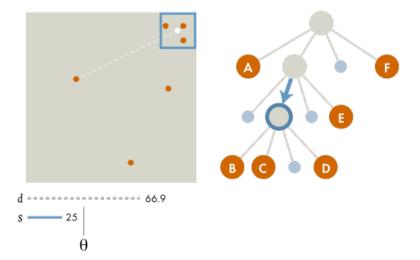
#### Barnes-Hut II



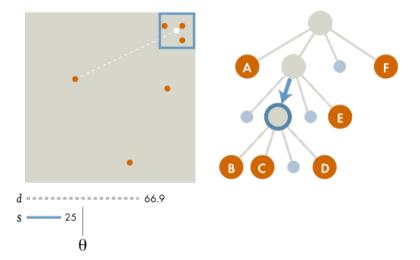
## Barnes-Hut III



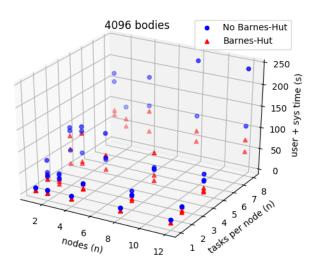
#### Barnes-Hut IV



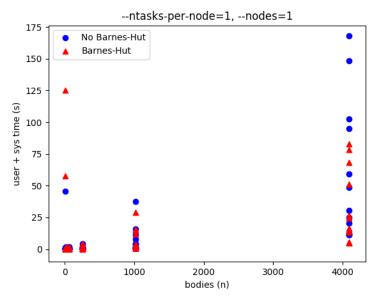
#### Barnes-Hut V



#### Barnes-Hut VI



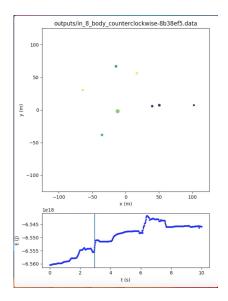
#### Barnes-Hut VII



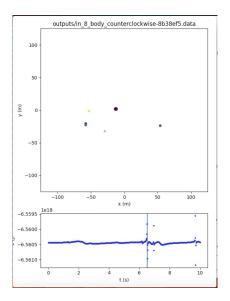
#### Correctness I

$$U = -G \frac{mM}{R}$$
$$E_{k} = \frac{1}{2} m v^{2}$$

#### Correctness II



#### Correctness III



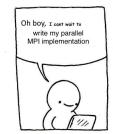
# Performance Optimizations I

Structure of vectors of properties VS vectors of stuctures as properties

# Performance Optimizations II

$$a_i = F(x_i)$$
$$v_{i+1} = v_i + a_i \, \Delta t$$

#### Parallelization I









#### Parallelization II

```
while (True) {
    quadtree = QuadTree(bodies);
    #pragma omp parallel for shared(bodies)
    for (size_t i = 0; i < bodies.size(); i++) {</pre>
        auto& body = sbodies[i];
        quadtree.calculate_force(body);
    }
 MPI Scatter
```

#### Parallelization III

```
#pragma omp parallel for shared(bodies)
  for (size_t i = 0; i < sbodies.size(); i++) {</pre>
      auto& body = sbodies[i];
      if (step % 2 == LEAP) {
           body.leap(timestep);
      else {
           body.frog(timestep);
MPI Gather
```

#### Parallelization IV

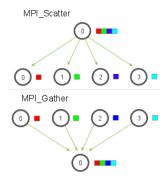
Total CPU time on rank 0 was 89.820000 Total CPU time on rank 1 was 37.040000 Total CPU time on rank 2 was 36.750000 Total CPU time on rank 3 was 36.960000

#### Parallelization V

# MPI\_Allgather

$$(b/n)n^2 = bn$$

#### Parallelization VI



$$2\frac{b}{n}n=2b$$

## Results I

