Assignment 12

Introduction

The task was to implement a 3-D n-body simulation. Although there are quite fast and clever algorithms for this problem like the Barnes-Hut-Algorithm, we decided to implement the naive version, which has a runtime complexity of $\mathcal{O}(n^2)$.

In the naive approach, every body compares its position directly to every other body, calculates the force between them and then updates its velocity and position accordingly.

Implementation

Each particle gets represented by a struct containing information about the mass, velocity and position of the body. Note that position and velocity consist of three components (one for every dimension).

```
typedef struct body_struct {
    float position[3];
    float velocity[3];
    int mass;
} body;
```

Following functions which are necessary for the simulation are provided in n_body.c:

Initialization

Every body gets initialized with a random mass, a velocity of 0 and a random position in the predefined space.

```
void initialize_bodies(body* bodies, int size){
    srand(time(NULL));
    unsigned int seed = rand();
    for (int i = 0; i<size; ++i){
        for (int j = 0; j<3; ++j){
            bodies[i].position[j] = rand_r(&seed) % MAX_DIM;
            bodies[i].velocity[j] = 0;
        }
        bodies[i].mass = ((rand_r(&seed) % MAX_MASS) + 500) * 1000;
    }
}</pre>
```

Updating velocity

The velocities get updated as described in the introduction section. Note that the force on an axis gets set to 0 as soon as bodies get too close, since we wanted to avoid absurdly high accelarations which can occur since we did not implement any collision model.

```
void update_velocities_naive(body* bodies, int size){
    for (int i = 0; i<size; ++i){
        for (int j = 0; j < size; ++j){
            if(i != j){
                float numerator = G * (bodies[i].mass + bodies[j].mass);
                for (int k = 0; k<3; ++k){
                     float radius = calculate_radius_squared(bodies[i].position[k],
bodies[j].position[k]);
                     float force;
                    if (fabs(radius) < 0.5){</pre>
                         force = 0;
                     } else{
                         force = numerator / radius;
                    bodies[i].velocity[k] += force/bodies[i].mass;
                }
            }
        }
    }
}
```

Updating position

The positions also get updated as described in the introduction section. Note. that the velocity of a body gets set to 0 as soon as the simulation boundaries are reached.

```
void update_positions(body* bodies, int size){
  for (int i = 0; i<size; ++i){
    for (int j = 0; j<3; ++j){
      float update = bodies[i].position[j] + bodies[i].velocity[j];
      if (update > MAX_DIM){
        bodies[i].velocity[j] = 0;
      }else if(update<0){
        bodies[i].velocity[j] = 0;
      } else{
        bodies[i].position[j] = update;
      }
    }
}</pre>
```

Saving the data to a file

In order to visualize the simulation it is necessary to write the position of each particle to a file. To achive this, the following function is provided:

```
void print_bodies_to_file(body* bodies, int size){
    FILE* file = fopen (OUT_FILE, "a");
    for (int j = 0; j<3; ++j){
        for (int i = 0; i<size; ++i){
            fprintf(file, "%f ", bodies[i].position[j]);
        }
        fprintf(file, "\n");
    }
    fprintf(file, "\n");
}</pre>
```

Putting it all together

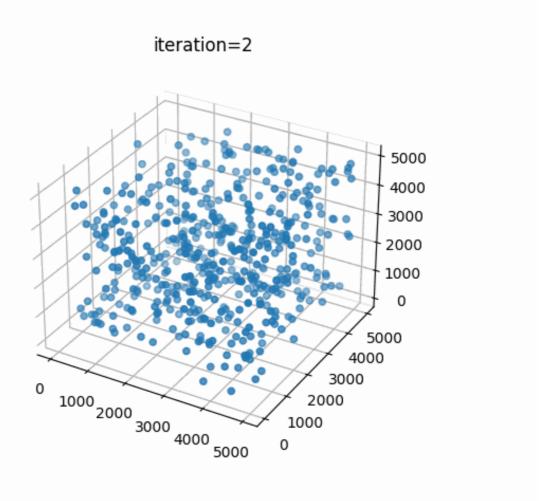
The main simulation loop is defined in the main function in simulation.c. Note that nothing is written to the file if the program is compiled with the TIME_MES flag.

```
for (int i = 0; i<steps; ++i){
    update_velocities_naive(bodies, size);
    update_positions(bodies, size);
    #ifndef TIME_MES
    print_bodies_to_file(bodies, size);
    #endif
}</pre>
```

Visualization of the simulation

We visualized the simulation with a custom python script using matplotlib. In order to run the script, one need to have matplotlib and numpy installed.

Here is an example of the generated plot:



Some observations

- Some points are very fast. This is due to some points are initialized with a very low mass and therefore get accellerated quite violently.
- The simulation probably needs some fine tuning in terms of parameters (e.g. the masses of particles and the gravitaional constant).
- A collision model would be very benefitial because we could avoid using workarounds (described in here) which currently lead to odd behaviour of the particles.

Parallel version

In order to parallelize the code, we added two simple statements in the most computationally intensive functions:

```
void update_positions(body* bodies, int size){
    #pragma omp parallel for collapse(2)
    for (int i = 0; i<size; ++i){
        for (int j = 0; j<3; ++j){
            ...
        }
    }
}</pre>
```

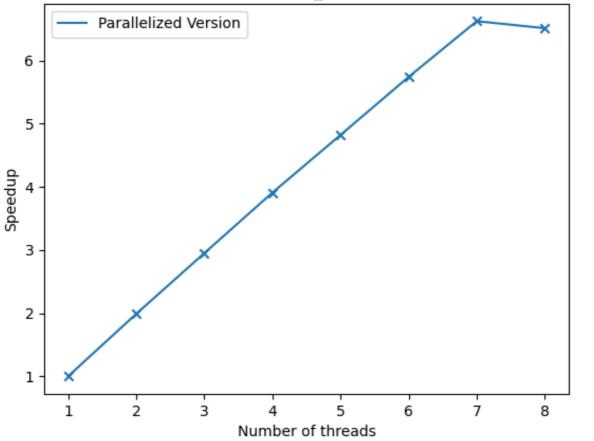
and

```
void update_velocities_naive(body* bodies, int size){
    #pragma omp parallel for collapse(2)
    for (int i = 0; i<size; ++i){
        for (int j = 0; j<size; ++j){
            ...
        }
    }
}</pre>
```

Benchmark

The program has been tested with $n_bodies = 5000$ and timesteps = 100.





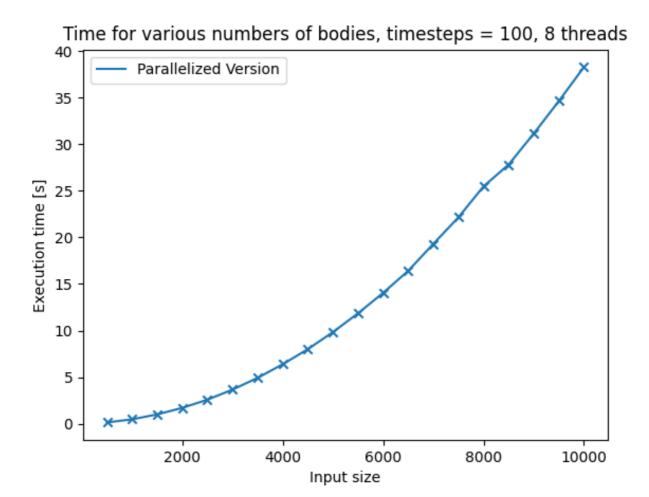
| Number of threads | Parallelized Version | Reference |
|-------------------|----------------------|-----------|
| 1 | 74.01 | 74.0 |
| 2 | 37.25 | 74.06 |
| 3 | 25.08 | 73.98 |
| 4 | 18.93 | 73.99 |
| 5 | 15.34 | 73.97 |
| 6 | 12.89 | 74.0 |
| 7 | 11.18 | 74.05 |
| 8 | 11.36 | 73.97 |

We are quite happy with this (nearly linear) speedup and especially with the effort to speedup ratio.

Time measurements for various inputs

Varriyng the number of bodies

Measurements have been taken with 100 timesteps and 8 threads for various numbers of bodies.

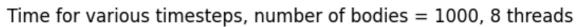


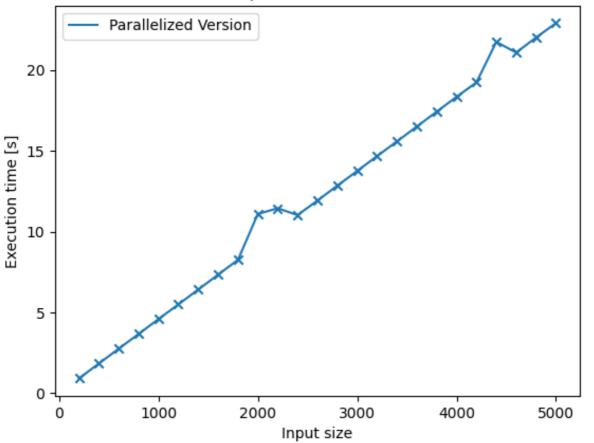
| Number of bodies | Execution time |
|------------------|----------------|
| 500 | 0.14 |
| 1000 | 0.47 |
| 1500 | 1.0 |
| 2000 | 1.7 |
| 2500 | 2.58 |
| 3000 | 3.66 |
| 3500 | 4.92 |
| 4000 | 6.38 |
| 4500 | 8.0 |
| 5000 | 9.83 |
| 5500 | 11.85 |
| 6000 | 14.04 |
| 6500 | 16.41 |
| 7000 | 19.3 |
| 7500 | 22.15 |
| 8000 | 25.47 |
| 8500 | 27.81 |
| 9000 | 31.14 |
| 9500 | 34.63 |
| 10000 | 38.27 |

One can clearly see the exponentially rising runtime (which fits our asymptotic complexity of $\mathcal{O}(n^2)$.

Varriyng number of timesteps

Measurements have been taken with 1000 bodies and 8 threads for various numbers of timesteps.





| Input size | Parallelized Version |
|------------|----------------------|
| 200 | 0.92 |
| 400 | 1.84 |
| 600 | 2.76 |
| 800 | 3.67 |
| 1000 | 4.59 |
| 1200 | 5.5 |
| 1400 | 6.42 |
| 1600 | 7.34 |
| 1800 | 8.26 |
| 2000 | 11.1 |
| 2200 | 11.45 |
| 2400 | 11.03 |
| 2600 | 11.93 |
| 2800 | 12.85 |
| 3000 | 13.76 |
| 3200 | 14.68 |
| 3400 | 15.59 |
| 3600 | 16.51 |
| 3800 | 17.43 |
| 4000 | 18.35 |
| 4200 | 19.24 |
| 4400 | 21.75 |
| 4600 | 21.09 |
| 4800 | 22.02 |
| 5000 | 22.91 |

This curve also fullfills our expecation, since the program has a linear runtinme complexity $(\mathcal{O}(n))$ in the number of timesteps.