



IMT Atlantique

Bretagne-Pays de la Loire

École Mines-Télécom

GALAX Project

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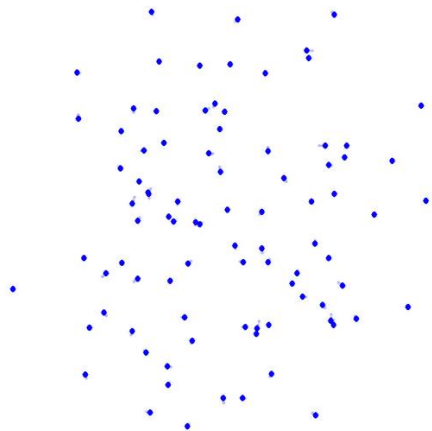
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CHAPTER 1

Results



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n-body problem

CPU Model

FPS: ~50

omp+xsimd+algorithm

```
average distance vs reference: 0.00101288; min error : 0; max error : 0.00745216
average distance vs reference: 0.00296431; min error : 0; max error : 0.0221624
average distance vs reference: 0.00526584; min error : 0; max error : 0.0437569
average distance vs reference: 0.00661514; min error : 0; max error : 0.052601
average distance vs reference: 0.00721774; min error : 0; max error : 0.0594342
average distance vs reference: 0.00809052; min error : 0; max error : 0.0716785
average distance vs reference: 0.00893563; min error : 0; max error : 0.0706503
average distance vs reference: 0.00966297; min error : 4.76837e-07; max error : 0.0601539
average distance vs reference: 0.0107886; min error : 9.53674e-07; max error : 0.076121
average distance vs reference: 0.0116452; min error : 9.53674e-07; max error : 0.107231
average distance vs reference: 0.0125096; min error : 9.53674e-07; max error : 0.187115
average distance vs reference: 0.0136752; min error : 2.13248e-06; max error : 0.246346
average distance vs reference: 0.0151134; min error : 2.38419e-06; max error : 0.201754
average distance vs reference: 0.0156335; min error : 2.05095e-06; max error : 0.182085
average distance vs reference: 0.0166952; min error : 2.4314e-06; max error : 0.21645
average distance vs reference: 0.0193012; min error : 4.26496e-06; max error : 0.326813
average distance vs reference: 0.0217572; min error : 4.26496e-06; max error : 0.421306
average distance vs reference: 0.0233909; min error : 5.37897e-06; max error : 0.623463
average distance vs reference: 0.0249607; min error : 5.76658e-06; max error : 0.746107
average distance vs reference: 0.0281325; min error : 6.48569e-06; max error : 0.506702
average distance vs reference: 0.0307454; min error : 7.26688e-06; max error : 0.494735
```

Method	N=10000
Naive	0
OMP PARFOR	4.5
XSIMD-v1	30
XSIMD-v2	50

CHAPTER 2

Method



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2.1 Main idea

Goal: find acceleration a_{ij} between every 2 particles i, j

= a symmetrical matrix whose diagonal line is all zero

```
for (int i = 0; i < n_particles; i++)  
{  
    for (int j = 0; j < n_particles; j++)  
    {  
        if (i != j)  
        {  
            ...  
        }  
    }  
}
```

```
#pragma omp parallel for
    for (int i = 0; i < n_particles; i++)
    {
        for (int j = i+1; j < n_particles; j++)
            ...
    }
```

Reduce half of the calculations:
change the square to the triangle

Use “omp parrallel for”

Use “Xsimd” to make vectorised calculation

```
for (int i = 0; i < n_particles; i++)  
{  
    b_type rposx_i = b_type::load_unaligned(&particles.x[i]);  
    ...  
    auto not_zero = 1 if batch > threshold;  
    dij = xs::select(not_zero, dij, 0);  
    ...  
    raccx_i.store_unaligned(&accelerationsx[i]);  
}
```

Optimization of “if”:

```
b_type c = xs::rsqrt(dij);  
dij = xs::fmin(10, 10.0 * dij * dij * dij);
```

2.3 XSIMD-v2

In version 1, check if the position of j is 0 for every batch.

for (int $j = 0$; $j < n_particles$; $j += 1$)

In version 2:

split the loop in version 1 to 2 loops:

first loop:

All the batches of j are complete

for (int $j = 0$; $j < n_particles - inc + 1$; $j += 1$)

second loop:

The previous value is the real value, the latter value does not exist, so it is filled with 0.

for (int $j = n_particles - inc + 1$; $j < n_particles$; $j += 1$)

Idea: Use multiple computer to multiply the calculation.

Core problem: data load/save conflit

We explain you...

CHAPTER 3

Other tricks



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IF or MIN ?

IF LOOP

real 0m0,202s
user 0m0,194s
sys 0m0,000s

MIN

real 0m0,308s
user 0m0,291s
sys 0m0,004s

#FMIN

real 0m0,370s
user 0m0,360s
sys 0m0,000s

Vscode Liveshare :(-> Any other program:(

“-O3” during compilation -> didn't change anything

Barnes-Hut Algorithm -> Octree ??

<https://www.cs.princeton.edu/~appel/papers/nbody.pdf>

CHAPTER 4

Demo



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Thanks for listening



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