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Pipelining the Fast Multipole Method over a Runtime System

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CSE 2013 - February 26

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Outline

- ▶ Overview of the FMM
- ▶ FMM on a runtime system
- ▶ Results and performance analysis
- ▶ Conclusion and Perspectives

Chebyshev Fast Multipole Method Overview

N-body problems

Example, interactions between planets in a galaxy:



$N(N - 1)$ interactions → **Complexity is $O(N^2)$.**

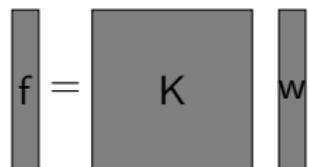
- ▶ Potentials (Kernel $\approx 1/r$)
 - ▶ Electrostatic → Molecular dynamics
 - ▶ Gravitational → Astrophysics
 - ▶ but also electromagnetism, fluid dynamics, VLSI capacitance,

...

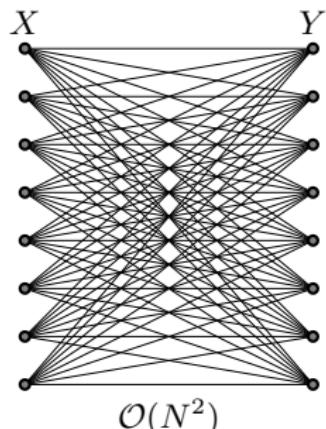
FMM : Motivations

- ▶ Pairwise interactions among N bodies.

$$f_i = \sum_{j=1}^N K(x_i, x_j) w_j \quad \text{for } i = 1, \dots, N.$$



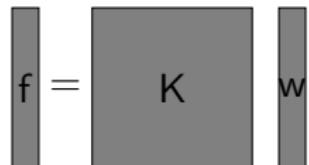
- ▶ Direct computation $\mathcal{O}(N^2)$



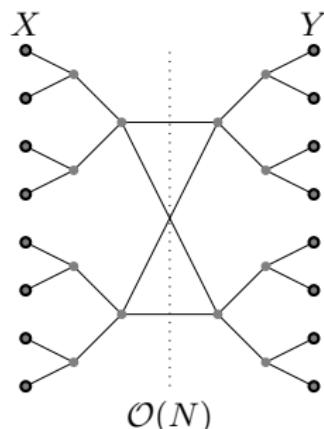
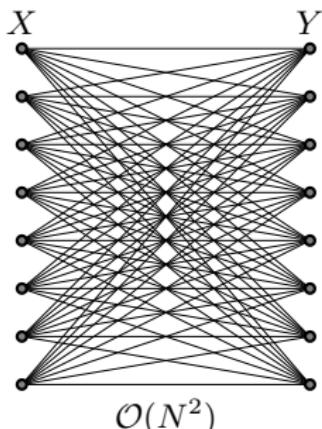
FMM : Motivations

- ▶ Pairwise interactions among N bodies.

$$f_i = \sum_{j=1}^N K(x_i, x_j) w_j \quad \text{for } i = 1, \dots, N.$$



- ▶ Direct computation $\mathcal{O}(N^2)$ → FMM $\mathcal{O}(N)$.

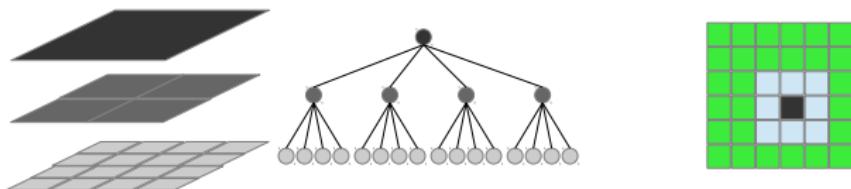


FMM : Overview

- ▶ Potential decomposition

$$f_i = f_i^{near} + f_i^{far}$$

- ▶ Involve hierarchical space decomposition with an tree (**octree**)



- ▶ Near field is computed by direct interactions.
- ▶ Different approaches to approximate the far field
 - ▶ Expansions (Cartesian or Spherical)
 - ▶ Interpolation (Chebyshev FMM)

Chebyshev FMM

Idea: Interpolate the kernel on a Chebyshev grid

$$\frac{1}{|\mathbf{x} - \mathbf{y}|} \sim \sum_{m=1}^{\ell} S_{\ell}(\mathbf{x}, \bar{x}_m) \sum_{n=1}^{\ell} \frac{1}{|\bar{x}_m - \bar{y}_m|} S_{\ell}(\mathbf{y}, \bar{y}_n)$$

Chebyshev FMM

Idea: Interpolate the kernel on a Chebyshev grid

$$\frac{1}{|\mathbf{x} - \mathbf{y}|} \sim \sum_{m=1}^{\ell} S_{\ell}(\mathbf{x}, \bar{x}_m) \sum_{n=1}^{\ell} \frac{1}{|\bar{x}_m - \bar{y}_m|} S_{\ell}(\mathbf{y}, \bar{y}_n)$$

S_{ℓ} is the interpolation polynomial (the higher ℓ the more accurate is the far field)

$$S_{\ell}(x, x_m) = \frac{1}{\ell} + \frac{2}{\ell} \sum_{n=1}^{\ell-1} T_n(\bar{x}_m) T_n(x)$$

Chebyshev FMM

Idea: Interpolate the kernel on a Chebyshev grid

$$\frac{1}{|\mathbf{x} - \mathbf{y}|} \sim \sum_{m=1}^{\ell} S_{\ell}(\mathbf{x}, \bar{x}_m) \sum_{n=1}^{\ell} \frac{1}{|\bar{x}_m - \bar{y}_m|} S_{\ell}(\mathbf{y}, \bar{y}_n)$$

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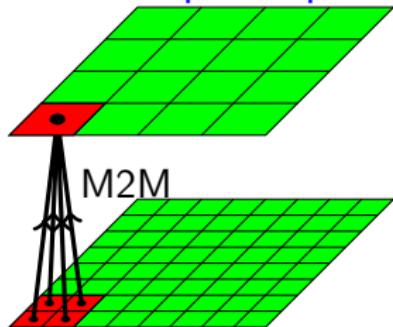
$$S_{\ell}(x, x_m) = \frac{1}{\ell} + \frac{2}{\ell} \sum_{n=1}^{\ell-1} T_n(\bar{x}_m) T_n(x)$$

For all $i = 1, \dots, N$.

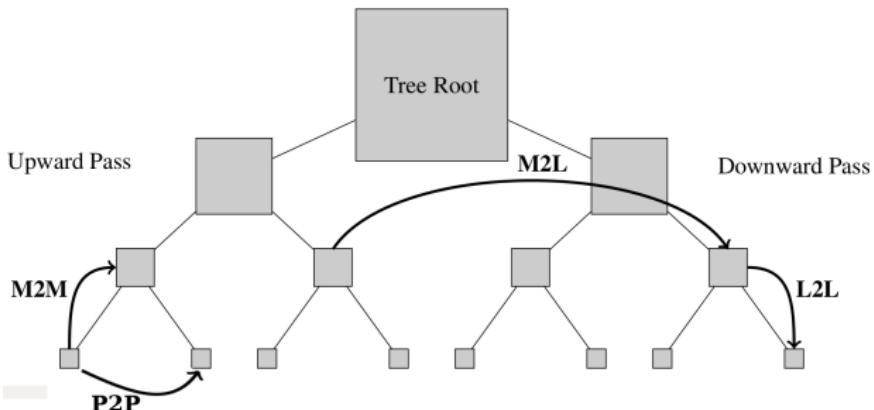
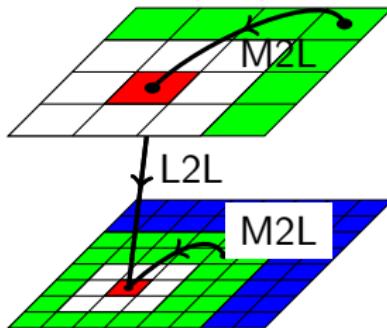
$$f_i = \sum_{j=1}^N \frac{1}{|x_i - y_j|} w_j$$
$$\sim \underbrace{\sum_{m=1}^{\ell} S_{\ell}(x_i, \bar{x}_m)}_{\text{L2L}} \underbrace{\sum_{n=1}^{\ell} \frac{1}{|\bar{x}_m - \bar{y}_n|}}_{\text{M2L}} \underbrace{\sum_{j=1}^{N^{\text{far}}} S_{\ell}(y_j, \bar{y}_n) w_j}_{\text{M2M}} + \underbrace{\sum_{j=1}^{N^{\text{near}}} \frac{1}{|x_i - y_j|} w_j}_{\text{P2P}}$$

FMM : $O(N)$ algorithm

Upward pass



Downward pass

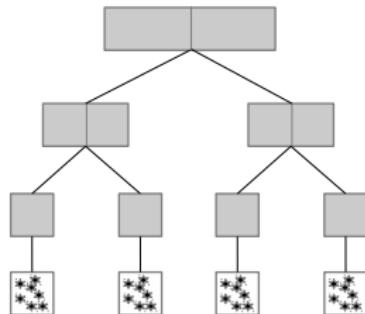


FMM on a runtime

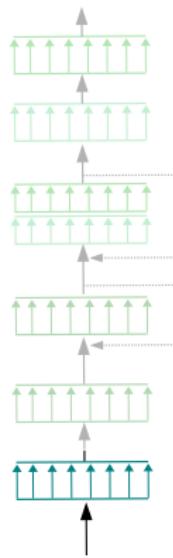
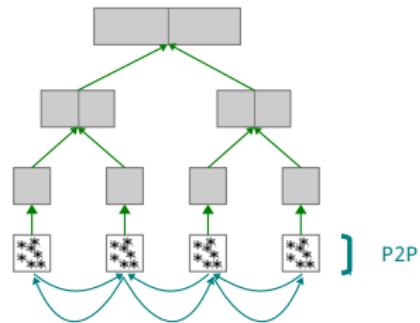
Fork-join parallelization : Description

Straightforward parallelization:

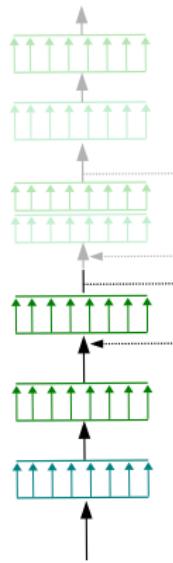
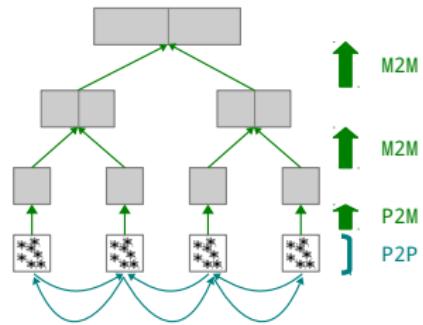
- ▶ Using fork-join model
- ▶ Progress level by level
- ▶ For example using *OpenMP for*



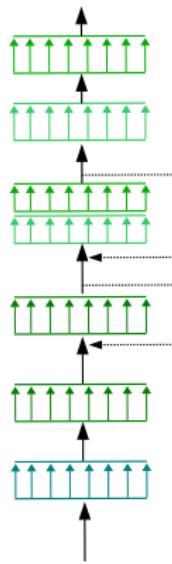
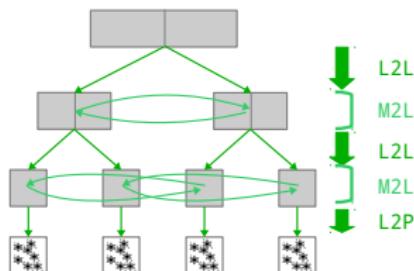
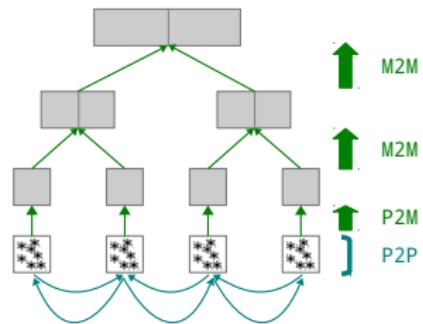
Fork-join parallelization : Simplified illustration



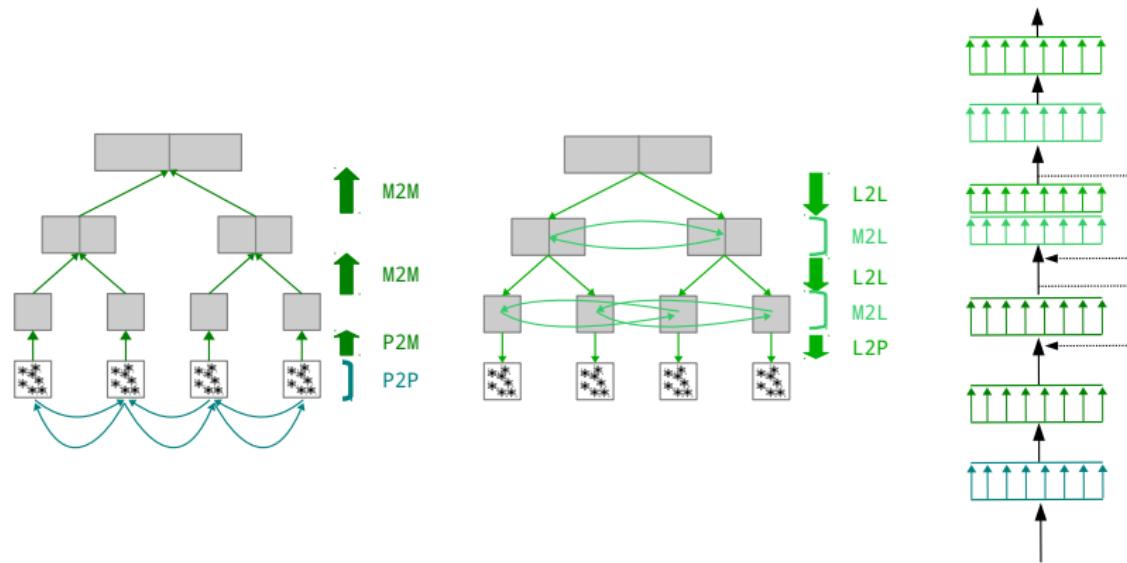
Fork-join parallelization : Simplified illustration



Fork-join parallelization : Simplified illustration

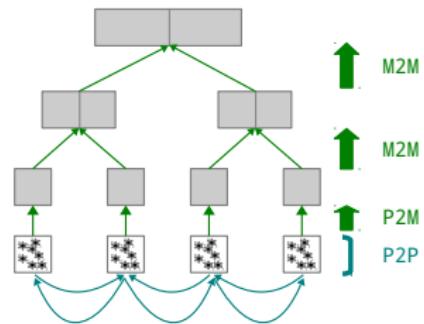


Fork-join parallelization : Simplified illustration

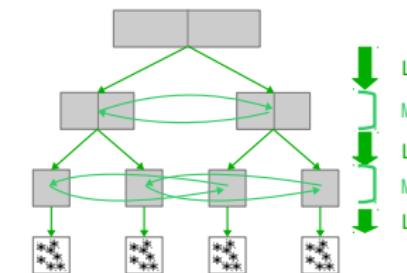


Limits: Synchronizations (+ bottleneck at the top of the tree).

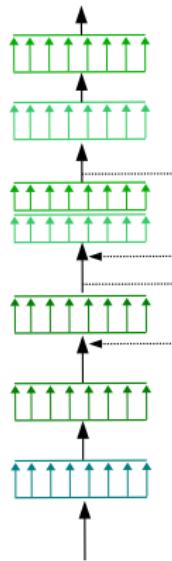
Fork-join parallelization : Simplified illustration



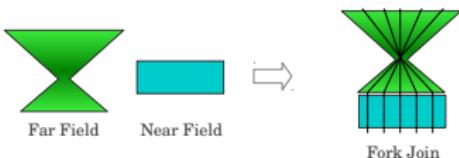
M2M
M2M
P2M
P2P



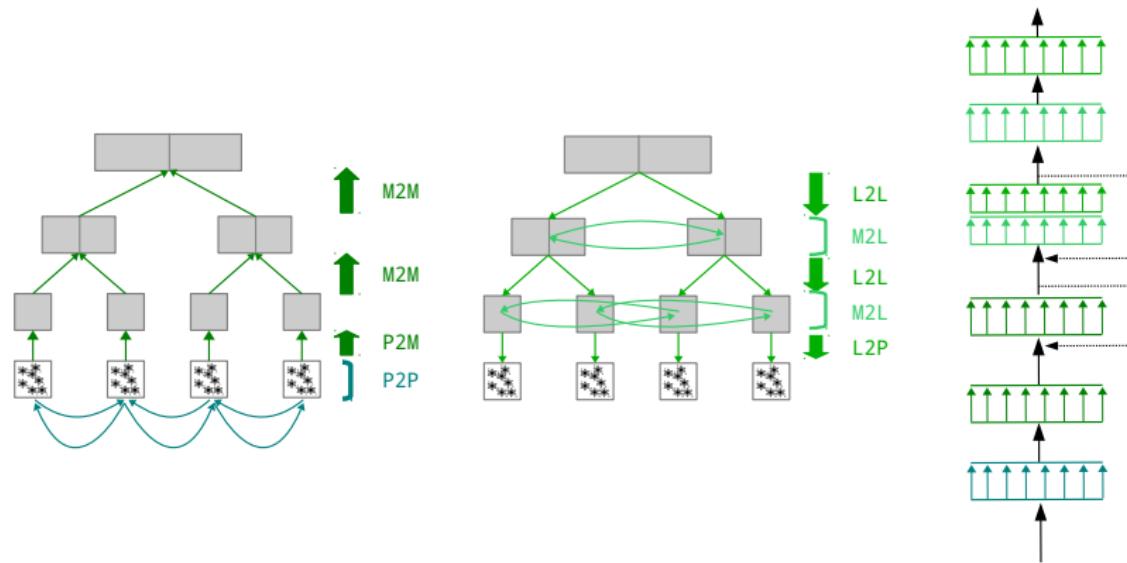
L2L
M2L
L2L
M2L
L2P



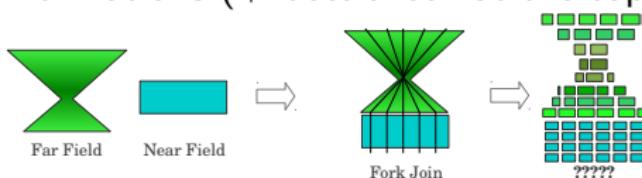
Limits: Synchronizations (+ bottleneck at the top of the tree).



Fork-join parallelization : Simplified illustration



Limits: Synchronizations (+ bottleneck at the top of the tree).



StarPU : Description

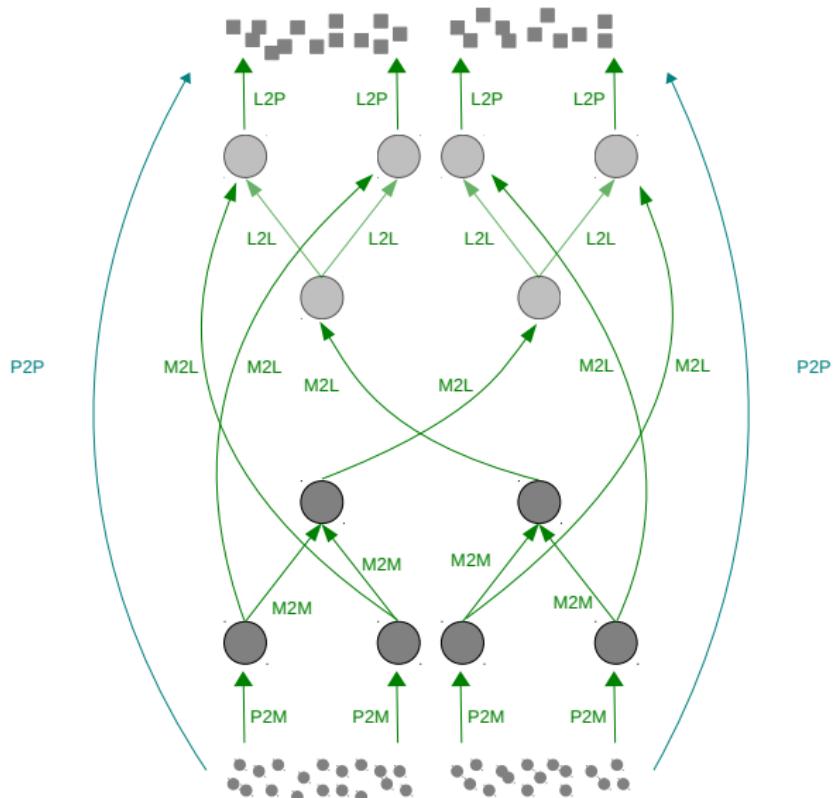
StarPU runtime

- ▶ Advanced tasks scheduling (customizable scheduler)
- ▶ Tasks dependencies expression
- ▶ Gives an abstraction of the architecture (useful for accelerators)
- ▶ <http://runtime.bordeaux.inria.fr/StarPU/>

Insert_task function:

- ▶ `insert_task (operator, {READ,WRITE}, data, ...)`

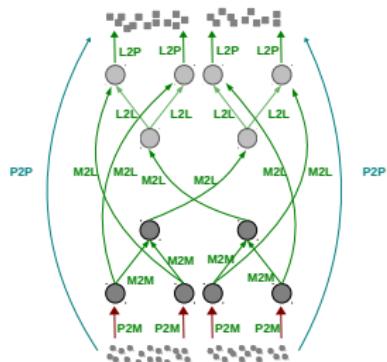
Exploiting the parallelism of the FMM



StarPU : Simplified example

Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
insert_task(P2M, READ, particles[1], WRITE, cell[2][1])
insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

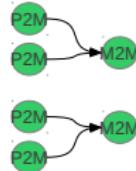
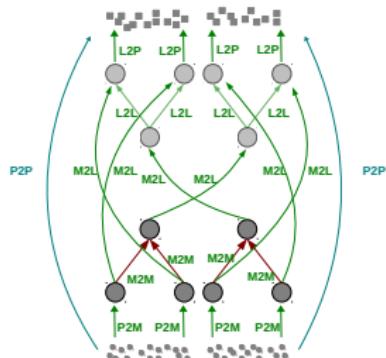


StarPU : Simplified example

Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
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insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```



StarPU : Simplified example

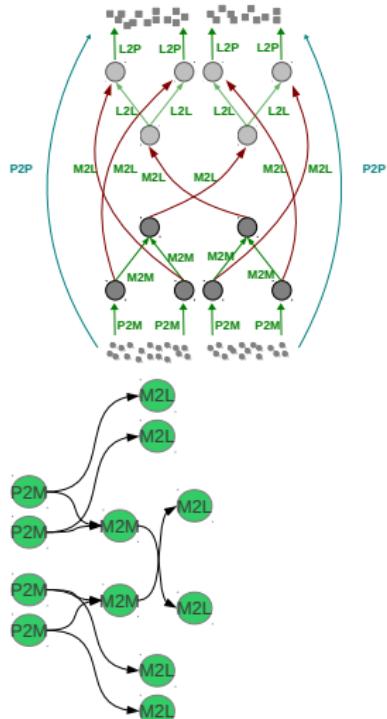
Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
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insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[1][0].interaction, WRITE, cell[1][0])
insert_task(M2L, READ, cell[1][1].interaction, WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[2][0].interaction, WRITE, cell[2][0])
insert_task(M2L, READ, cell[2][1].interaction, WRITE, cell[2][1])
insert_task(M2L, READ, cell[2][2].interaction, WRITE, cell[2][2])
insert_task(M2L, READ, cell[2][3].interaction, WRITE, cell[2][3])
```



StarPU : Simplified example

Cells access → cell[level][position]

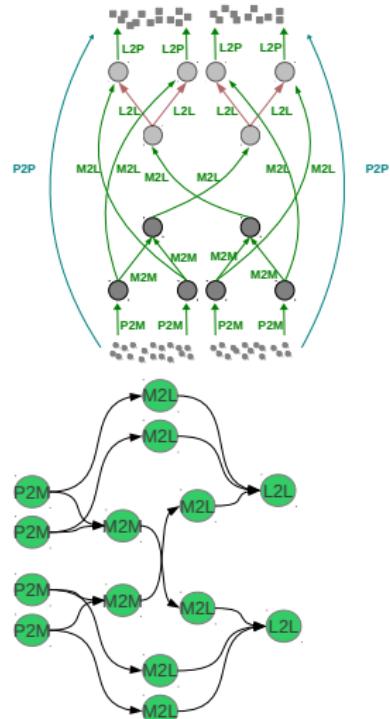
```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
insert_task(P2M, READ, particles[1], WRITE, cell[2][1])
insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[1][0].interaction, WRITE, cell[1][0])
insert_task(M2L, READ, cell[1][1].interaction, WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[2][0].interaction, WRITE, cell[2][0])
insert_task(M2L, READ, cell[2][1].interaction, WRITE, cell[2][1])
insert_task(M2L, READ, cell[2][2].interaction, WRITE, cell[2][2])
insert_task(M2L, READ, cell[2][3].interaction, WRITE, cell[2][3])
```

```
insert_task(L2L, READ, cell[1][0], WRITE, cell[2][0], cell[2][1])
insert_task(L2L, READ, cell[1][1], WRITE, cell[2][2], cell[2][3])
```



StarPU : Simplified example

Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
insert_task(P2M, READ, particles[1], WRITE, cell[2][1])
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insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

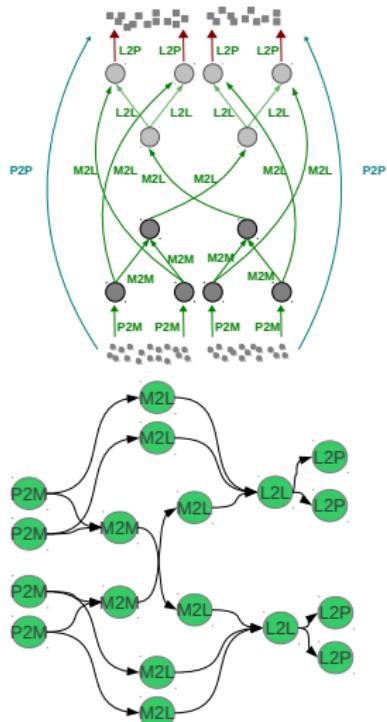
```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[1][0].interaction, WRITE, cell[1][0])
insert_task(M2L, READ, cell[1][1].interaction, WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[2][0].interaction, WRITE, cell[2][0])
insert_task(M2L, READ, cell[2][1].interaction, WRITE, cell[2][1])
insert_task(M2L, READ, cell[2][2].interaction, WRITE, cell[2][2])
insert_task(M2L, READ, cell[2][3].interaction, WRITE, cell[2][3])
```

```
insert_task(L2L, READ, cell[1][0], WRITE, cell[2][0], cell[2][1])
insert_task(L2L, READ, cell[1][1], WRITE, cell[2][2], cell[2][3])
```

```
insert_task(L2P, READ, cell[2][0], WRITE, particles[0])
insert_task(L2P, READ, cell[2][1], WRITE, particles[1])
insert_task(L2P, READ, cell[2][2], WRITE, particles[2])
insert_task(L2P, READ, cell[2][3], WRITE, particles[3])
```



StarPU : Simplified example

Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
insert_task(P2M, READ, particles[1], WRITE, cell[2][1])
insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```

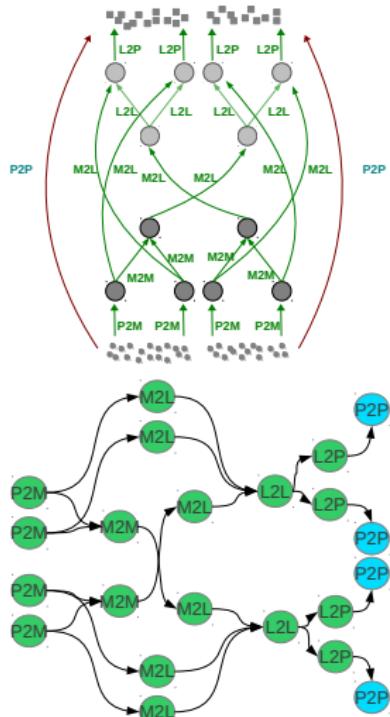
```
insert_task(M2L, READ, cell[1][0].interaction, WRITE, cell[1][0])
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```

```
insert_task(M2L, READ, cell[2][0].interaction, WRITE, cell[2][0])
insert_task(M2L, READ, cell[2][1].interaction, WRITE, cell[2][1])
insert_task(M2L, READ, cell[2][2].interaction, WRITE, cell[2][2])
insert_task(M2L, READ, cell[2][3].interaction, WRITE, cell[2][3])
```

```
insert_task(L2L, READ, cell[1][0], WRITE, cell[2][0], cell[2][1])
insert_task(L2L, READ, cell[1][1], WRITE, cell[2][2], cell[2][3])
```

```
insert_task(L2P, READ, cell[2][0], WRITE, particles[0])
insert_task(L2P, READ, cell[2][1], WRITE, particles[1])
insert_task(L2P, READ, cell[2][2], WRITE, particles[2])
insert_task(L2P, READ, cell[2][3], WRITE, particles[3])
```

```
insert_task(P2P, READ, neighbours[0], WRITE, particles[0])
insert_task(P2P, READ, neighbours[1], WRITE, particles[1])
insert_task(P2P, READ, neighbours[2], WRITE, particles[2])
insert_task(P2P, READ, neighbours[3], WRITE, particles[3])
```



StarPU : Simplified example

Cells access → cell[level][position]

```
insert_task(P2M, READ, particles[0], WRITE, cell[2][0])
insert_task(P2M, READ, particles[1], WRITE, cell[2][1])
insert_task(P2M, READ, particles[2], WRITE, cell[2][2])
insert_task(P2M, READ, particles[3], WRITE, cell[2][3])
```

```
insert_task(M2M, READ, cell[2][0], cell[2][1], WRITE, cell[1][0])
insert_task(M2M, READ, cell[2][2], cell[2][3], WRITE, cell[1][1])
```

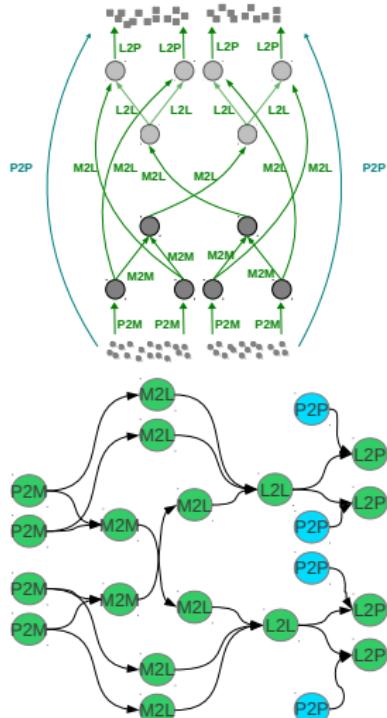
```
insert_task(M2L, READ, cell[1][0].interaction, WRITE, cell[1][0])
insert_task(M2L, READ, cell[1][1].interaction, WRITE, cell[1][1])
```

```
insert_task(M2L, READ, cell[2][0].interaction, WRITE, cell[2][0])
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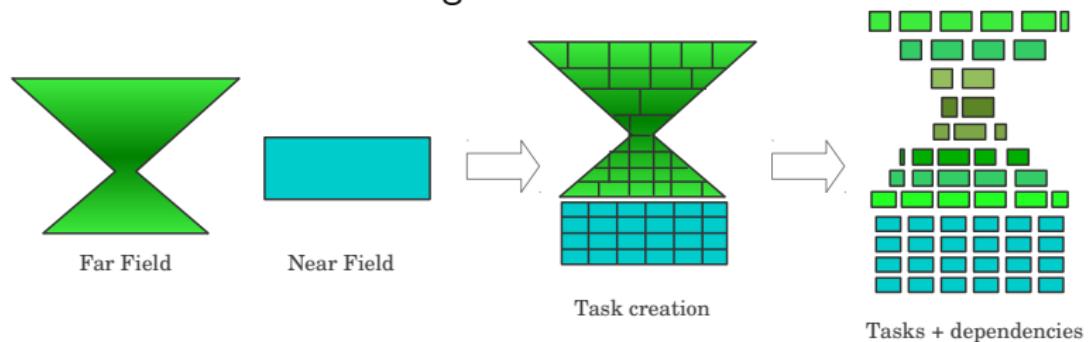
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insert_task(P2P, READ, neighbours[3], WRITE, particles[3])
```

```
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insert_task(L2P, READ, cell[2][3], WRITE, particles[3])
```



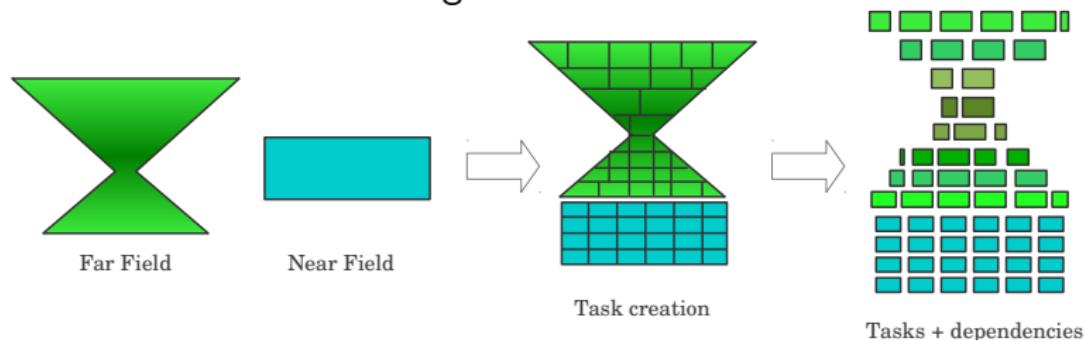
StarPU : Tasks creation and execution

Creation of the tasks using insert function:

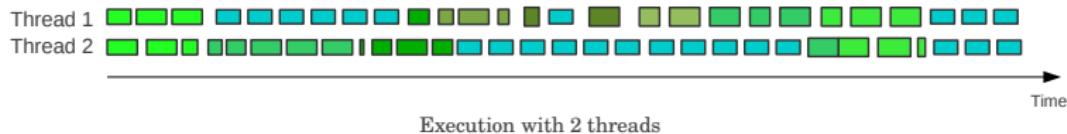


StarPU : Tasks creation and execution

Creation of the tasks using insert function:



Execution of the tasks (trace):



StarPU : Using GPU kernels

Two GPU kernels : P2P and M2L

- ▶ P2P & M2L represent more than 90% of the work

Scheduling:

- ▶ Schedulers assign tasks to processing units.
- ▶ StarPU proposes several schedulers (but none fits our needs).
- ▶ We developed a new scheduler:
 - ▶ Based on Eager
 - ▶ Enable setting priorities to operators
 - ▶ Different priorities for GPU or CPU

Results & Performance

Experimental Setup

ScalFMM:

- ▶ Developed in C++ (OpenMP, MPI, and now StarPU)
- ▶ Available at <http://scalfmm-public.gforge.inria.fr/>

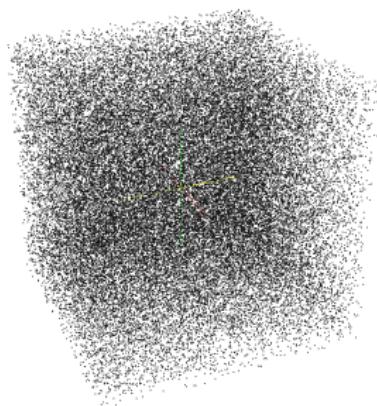
Test cluster:

- ▶ Heterogeneous 12 CPU and 3 GPU
- ▶ CPU dual-socket hexa-core Intel X5650 Nehalem (2.67 GHz)
- ▶ GPU Nvidia M2070 (or M2090)

In the following results Intel Compiler (12.0.5) is used.

Test case

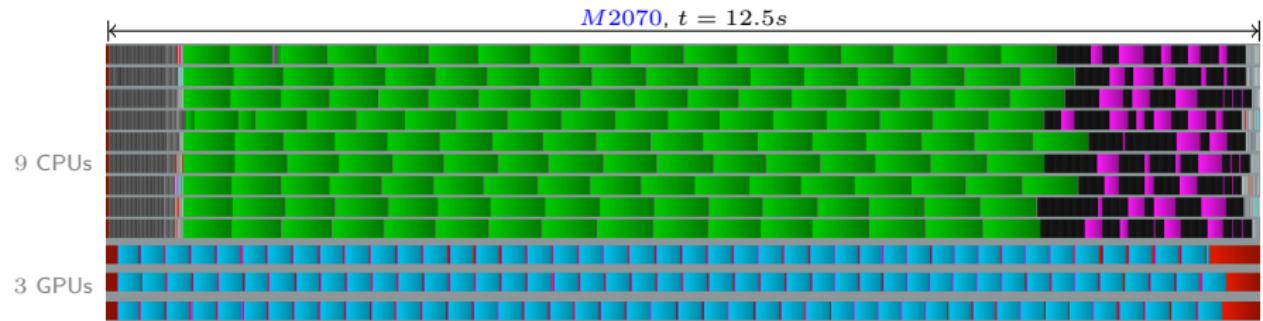
- ▶ Uniform distribution of particles in 3D inside a cube
- ▶ Particles 30.10^6
- ▶ Granularity $n_g = 2000$
- ▶ Octree height $h = \{6, 7\}$
- ▶ $Acc(\ell) = 10^{-\ell}$, with $\ell = \{2 - 7\}$
- ▶ $GPU = \{0, 1, 2, 3\}$





StarPU Execution traces M2070-M2090

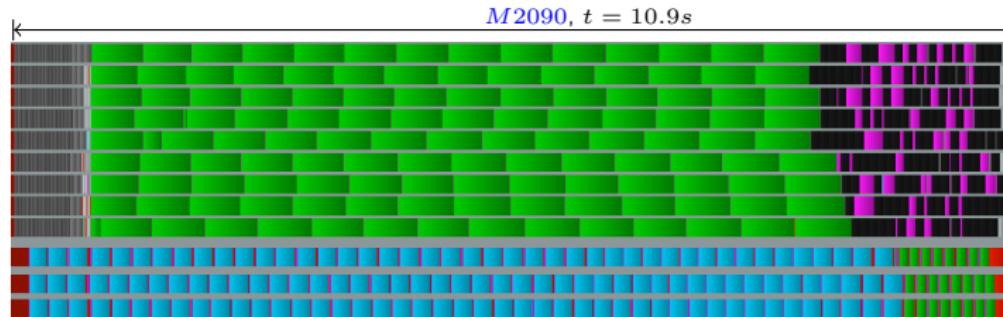
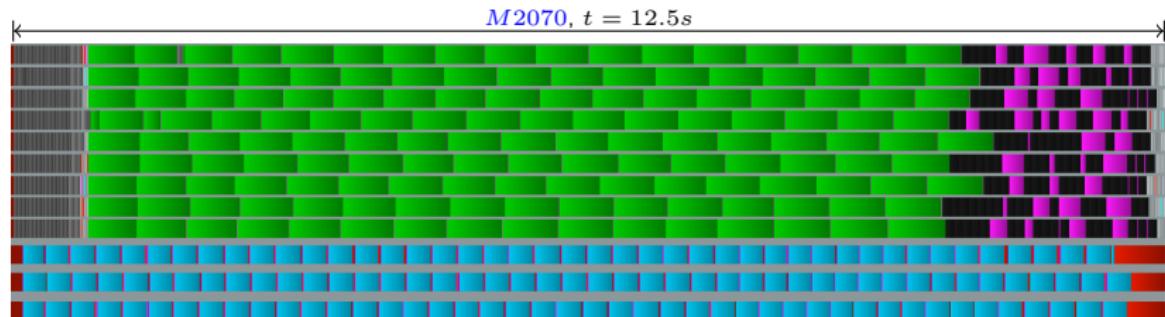
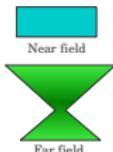
Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$,
 $Acc(\ell = 5) = 10^{-5}$



Legend : P2P, P2M, M2M, M2L, L2L, L2P, Idle, Moving data .

StarPU Execution traces M2070-M2090

Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$,
 $Acc(\ell = 5) = 10^{-5}$



Legend : P2P, P2M, M2M, M2L, L2L, L2P, Idle, Moving data .

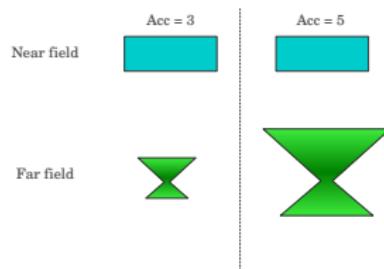
StarPU Execution traces M2070

Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$



StarPU Execution traces M2070

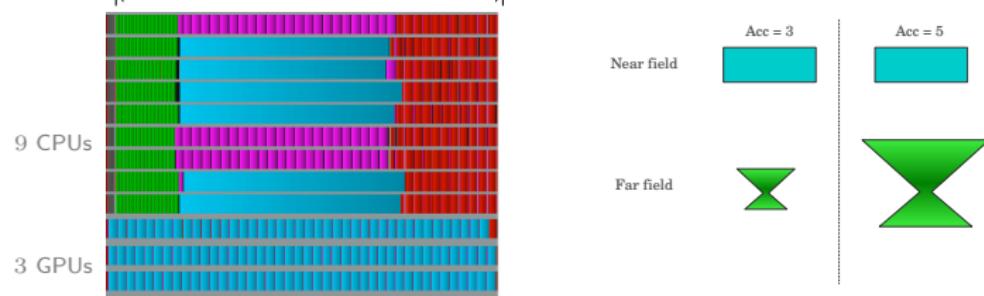
Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$



StarPU Execution traces M2070

Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$

$\text{Acc}(\ell = 3) = 10^{-3}$, $t = 11.6\text{s}$

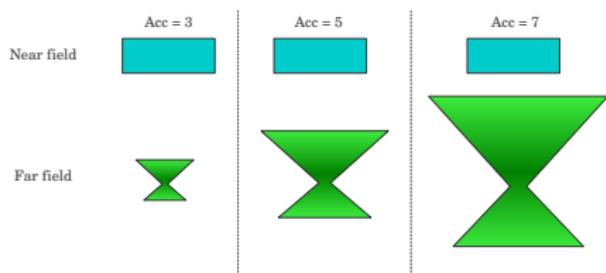
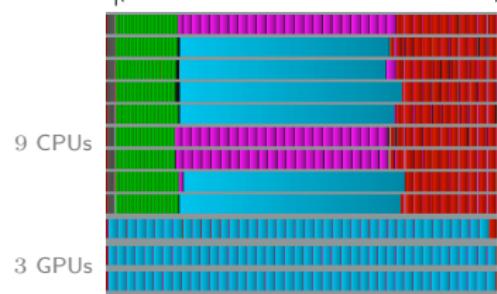


Legend : P2P, P2M, M2M, M2L, L2L, L2P, Idle, Moving data .

StarPU Execution traces M2070

Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$

$\text{Acc}(\ell = 3) = 10^{-3}$, $t = 11.6\text{s}$

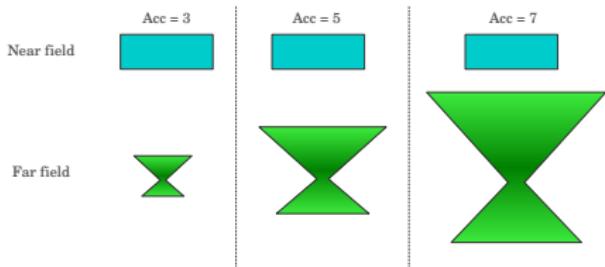
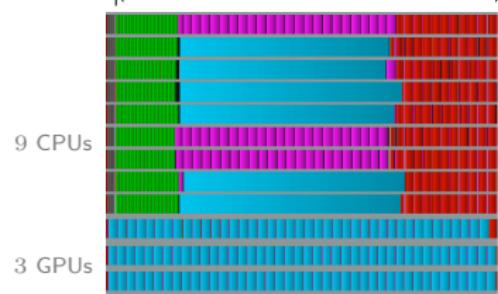


Legend : P2P, P2M, M2M, M2L, L2L, L2P, Idle, Moving data .

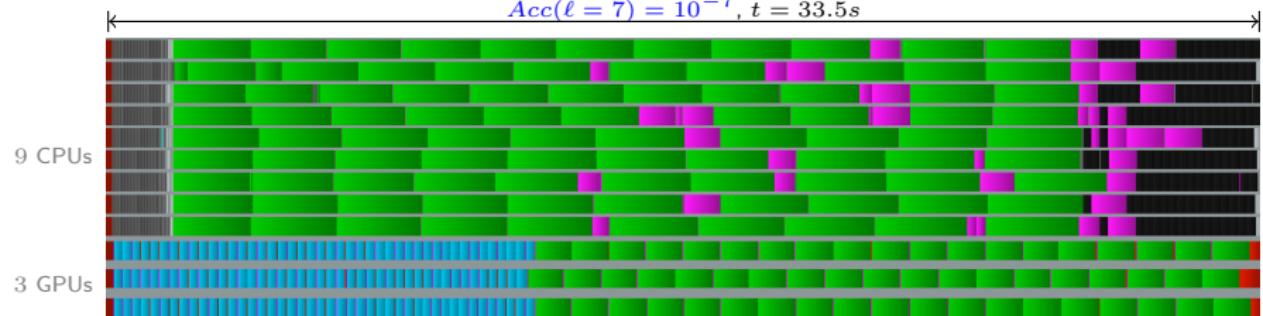
StarPU Execution traces M2070

Particles $30 \cdot 10^6$, Octree height $h = 6$, Granularity $n_g = 2000$

$\text{Acc}(\ell = 3) = 10^{-3}$, $t = 11.6\text{s}$



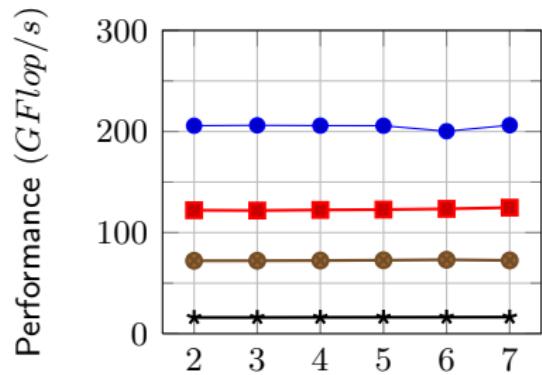
$\text{Acc}(\ell = 7) = 10^{-7}$, $t = 33.5\text{s}$



Legend : P2P, P2M, M2M, M2L, L2L, L2P, Idle, Moving data .

Results M2070

30.10^6 particles inside a box, granularity $n_g = 2000$, for 0 to 3 GPU and accuracy from 2 to 7.

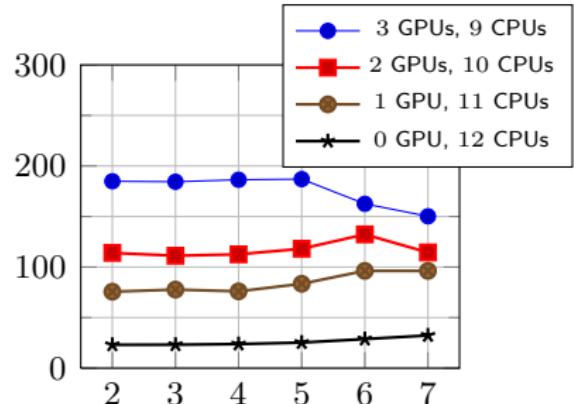


Accuracy $\ell (10^{-\ell})$

$h = 6$

Near field

Far field



Accuracy $\ell (10^{-\ell})$

$h = 7$

Near field

Far field

Conclusion

- ▶ Runtime optimization of parallel mapping was demonstrated.
- ▶ Modern heterogeneous computing platforms lead to hard DAG scheduling.
- ▶ Approach like Fork-join may be insufficient.
- ▶ Code is hardware “independent.” Schedulers determine the best processor to execute a given task.

Acknowledgement

- ▶ Inria & FastLA associate team.
- ▶ StarPU development team.
- ▶ ScalFMM <http://scalfmm-public.gforge.inria.fr/>

Thanks - Questions?

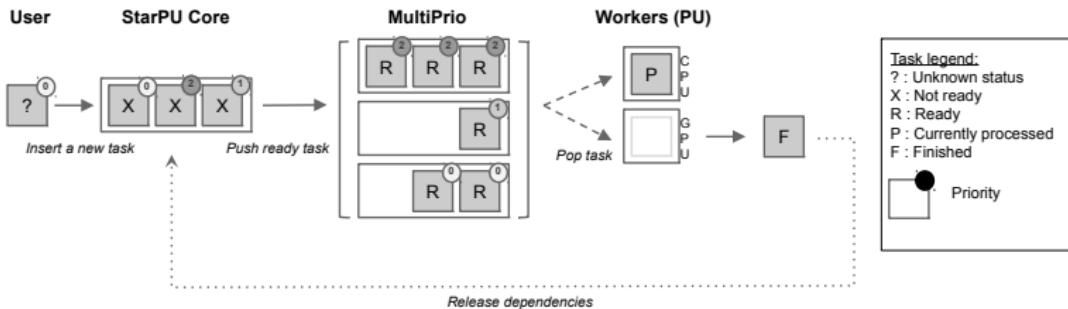


Annexes

Annexes



Annexe : Scheduler



Annexe : Using CPU and GPU

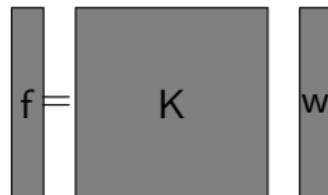
```
function M2L()
    forall the levels lv from 2 to leaf_level do
        foreach Block bl in blocks[lv] do
            insert_task( cpu_m2l, gpu_m2l, READ, bl.interactions,
                         WRITE, bl);
```

- ▶ StarPU is able to compute a M2L on the CPU or on the device (after moving the data)

Chebyshev FMM : Direct Matrix Vector

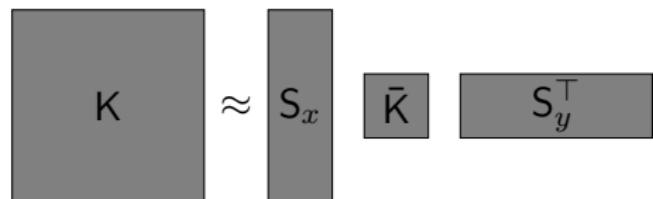
- ▶ $f = K w$

- ▶ $(K)_{ij} = K(x_i, y_j) = \frac{1}{|x_i - y_j|}$



Interpolation based Matrix Vector (low rank approximation)

- ▶ $K \approx S_x \bar{K} S_y^T$
- ▶ $S_x, S_y \text{ dim } N \times \ell^3$
- ▶ $\bar{K} \text{ dim } \ell^3 \times \ell^3$



Second low rank approximation (SVD)

- ▶ $\bar{K} \approx U V^T$
- ▶ $U, V \text{ dim } \ell^3 \times r$

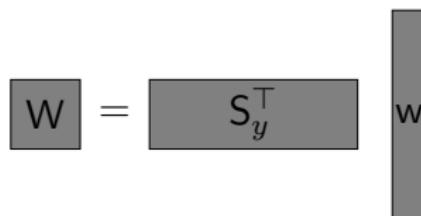


Chebyshev FMM : Interpolation based Matrix Vector

$$f = K w \approx S_x \bar{K} S_y^\top w$$

P2M/M2M:

- ▶ $W = S_y^\top w$



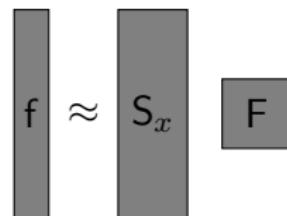
M2L:

- ▶ $F = \bar{K} W$



L2L/L2P:

- ▶ $f \approx S_x F$



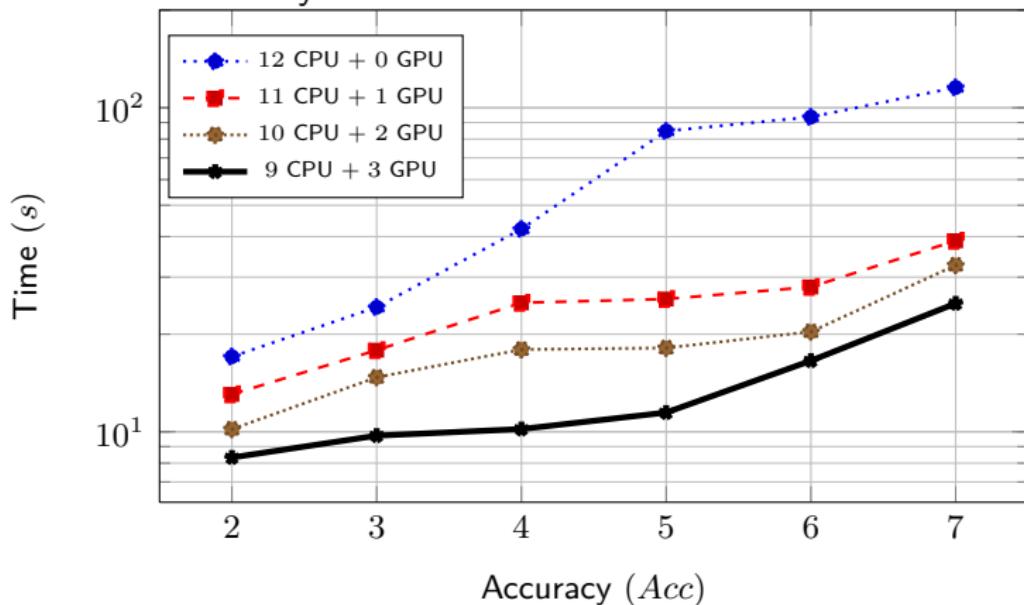
Task : Introduction

A task is composed by:

- ▶ A block of instructions
- ▶ Some parameters
- ▶ An environment
- ▶ Tasks is native in many "libraries" (OpenMP 3.1, StarPU)

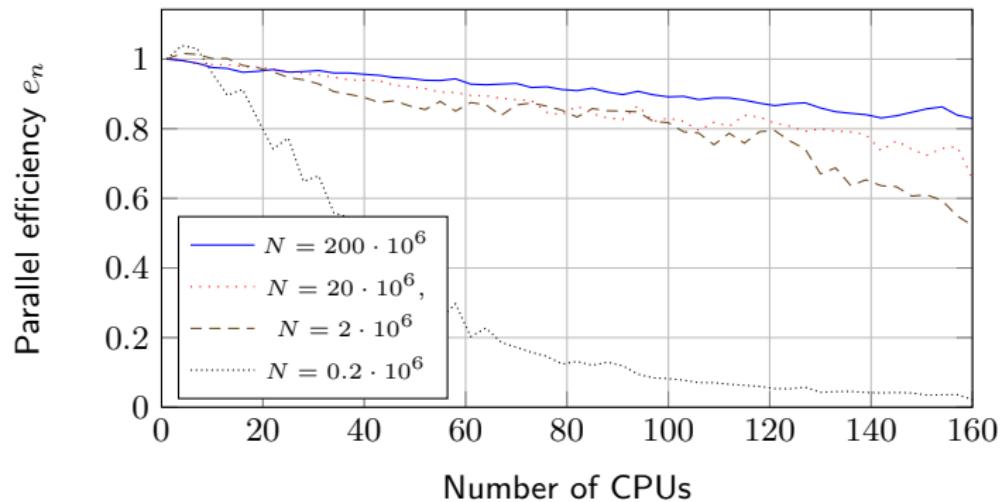
Results M2070

30.10^6 particles inside a box, granularity $n_g = 2000$, for 0 to 3 GPU and accuracy from 2 to 7.



160 Cores efficiency

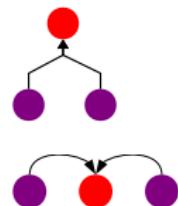
$$\begin{aligned} & \{N = 0.2 \cdot 10^6, n_g = 25, h = 5\}, \quad \{N = 2 \cdot 10^6, n_g = 200, h = 6\} \\ & \{N = 20 \cdot 10^6, n_g = 1000, h = 7\}, \quad \{N = 200 \cdot 10^6, n_g = 3000, h = 8\} \end{aligned}$$



StarPU : Creating tasks

Examples:

```
insert_task(m2m, READ, children, WRITE, cell);
```



```
insert_task(m2l, READ, interactions, WRITE, cell);
```

Order of insertion matters:

- ▶ It leads to the DAG
- ▶ Different orders can create different DAG