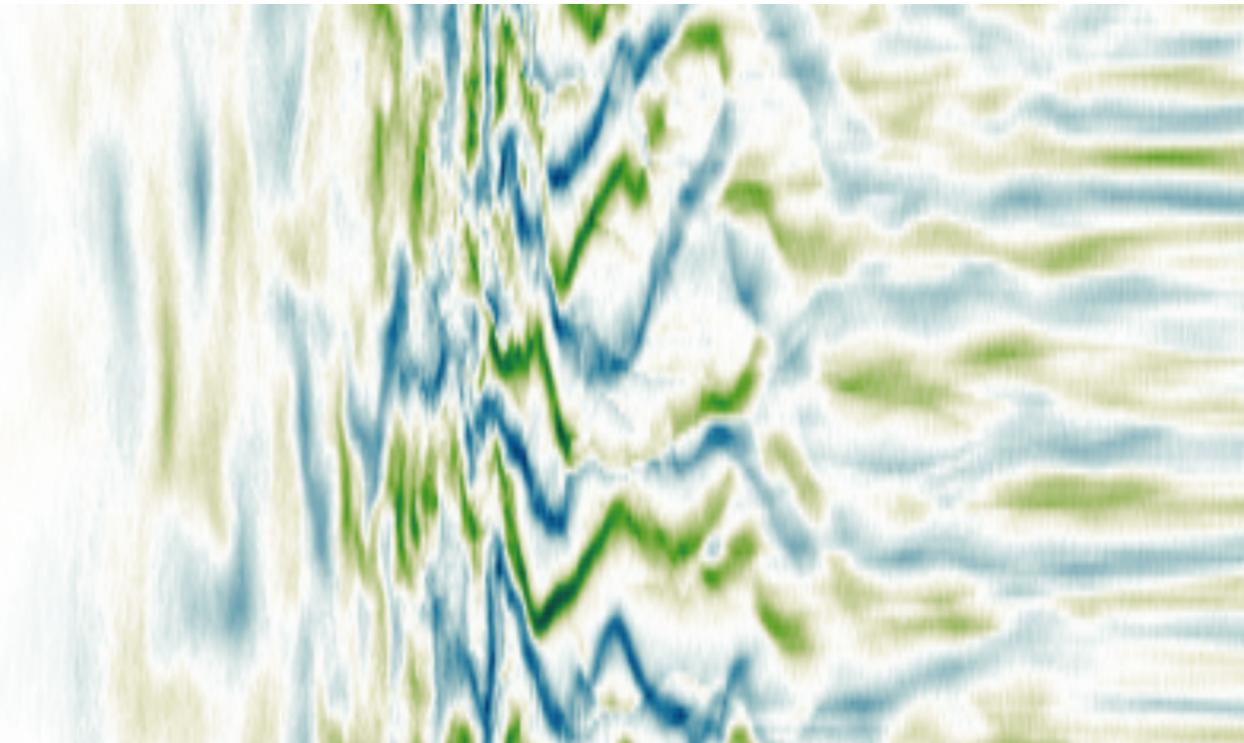


The Particle-In-Cell (PIC) simulation of plasmas

Mickael Grech, LULI, CNRS

mickael.grech@gmail.com

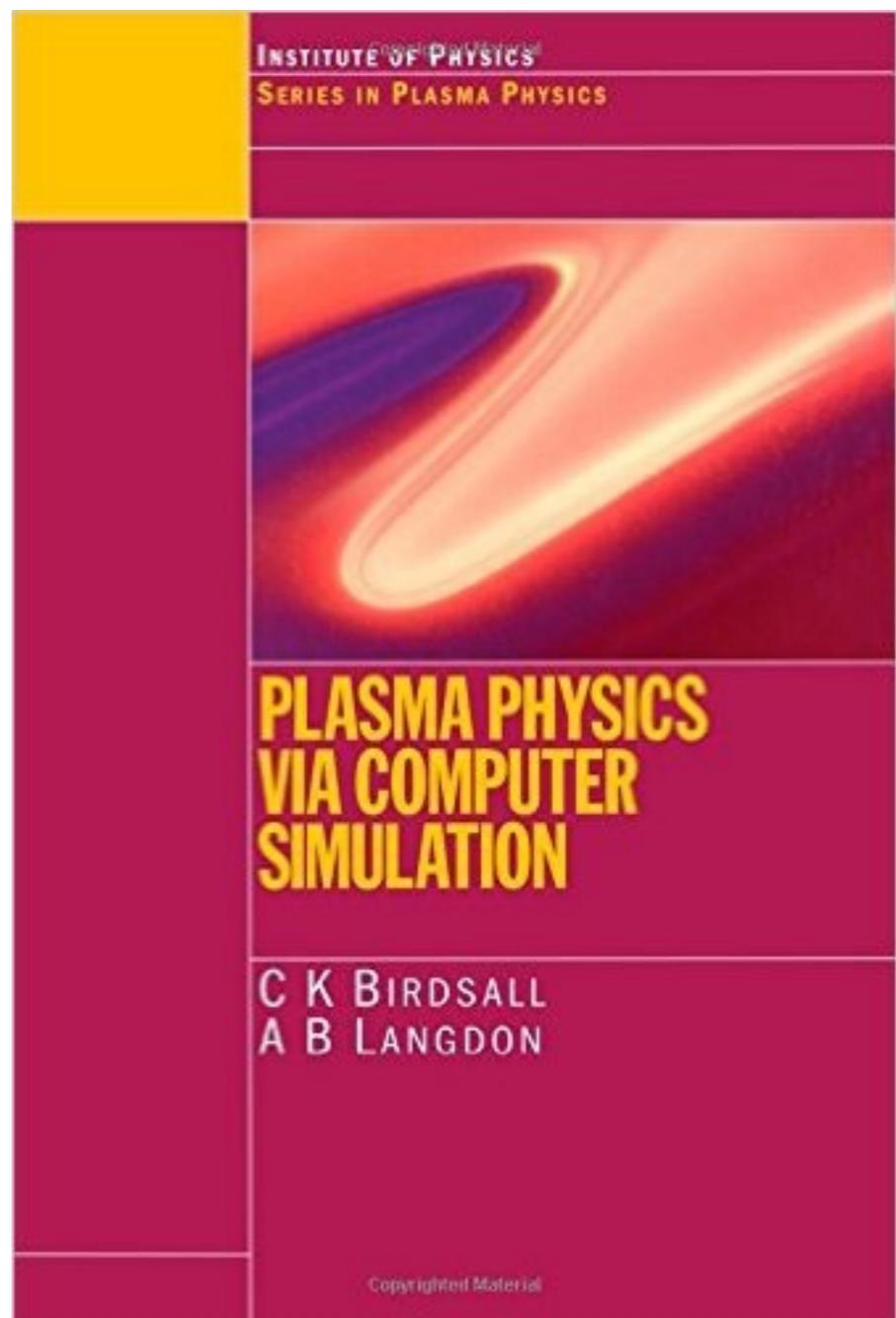
Les Houches, Mai 2019



References

References

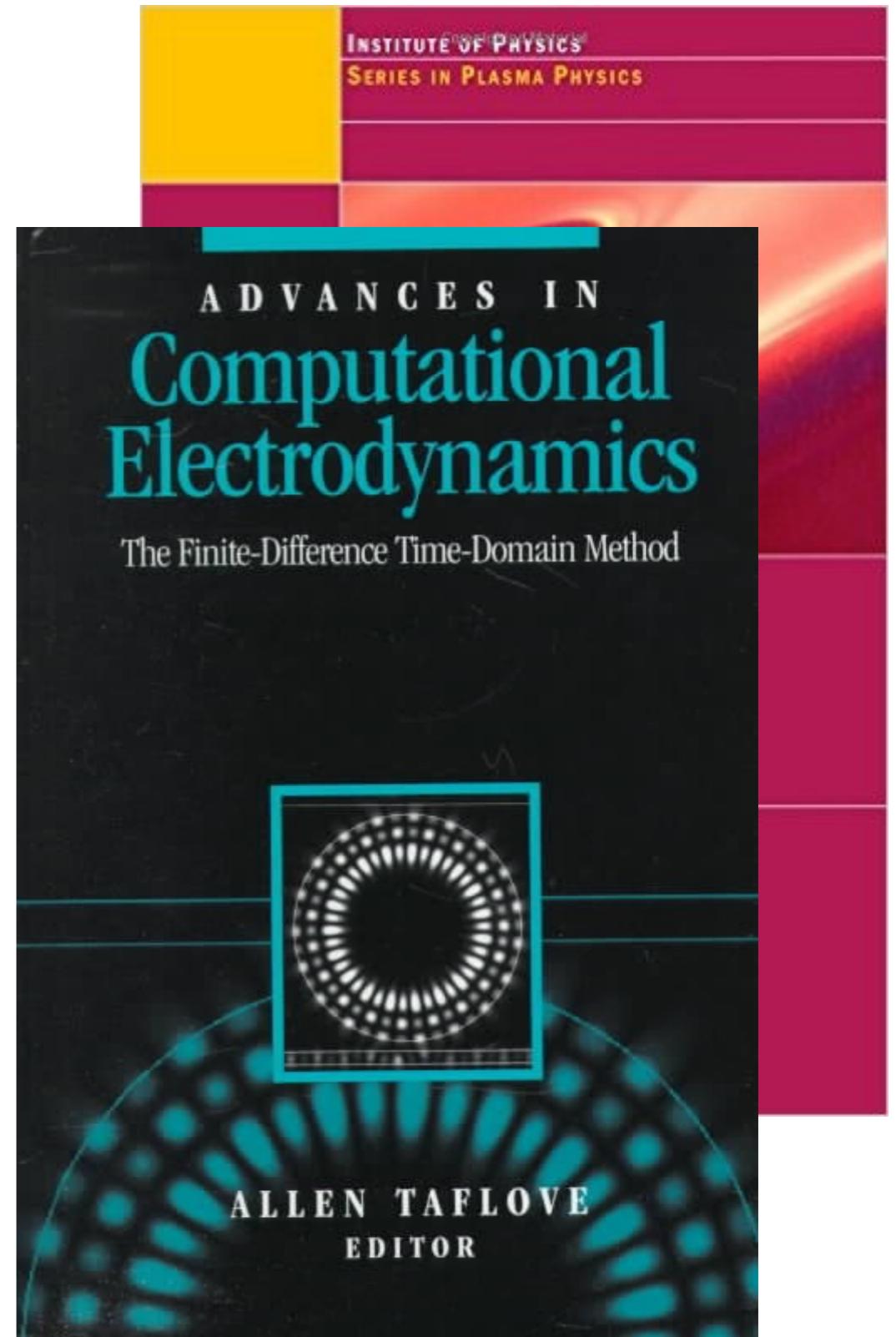
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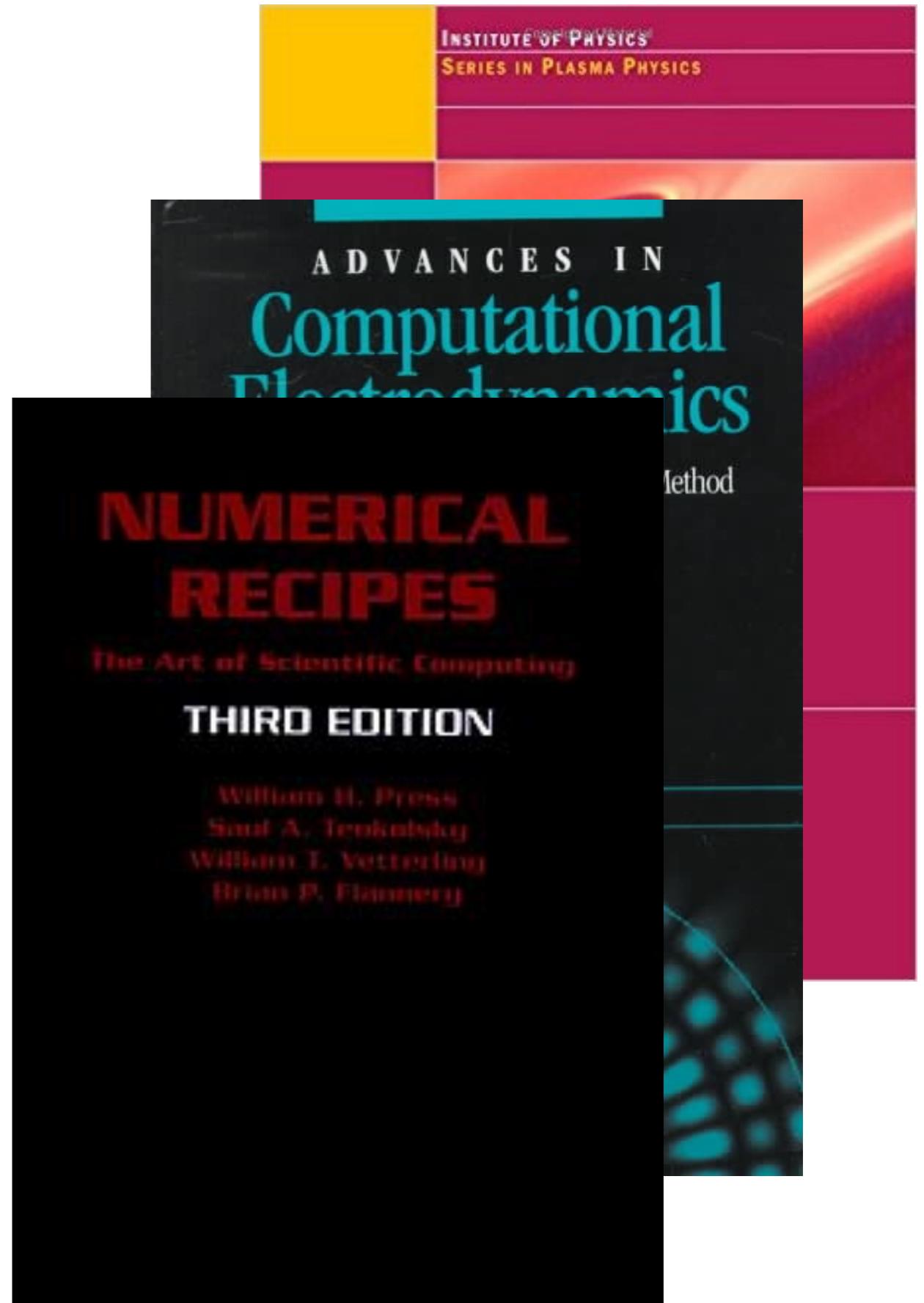


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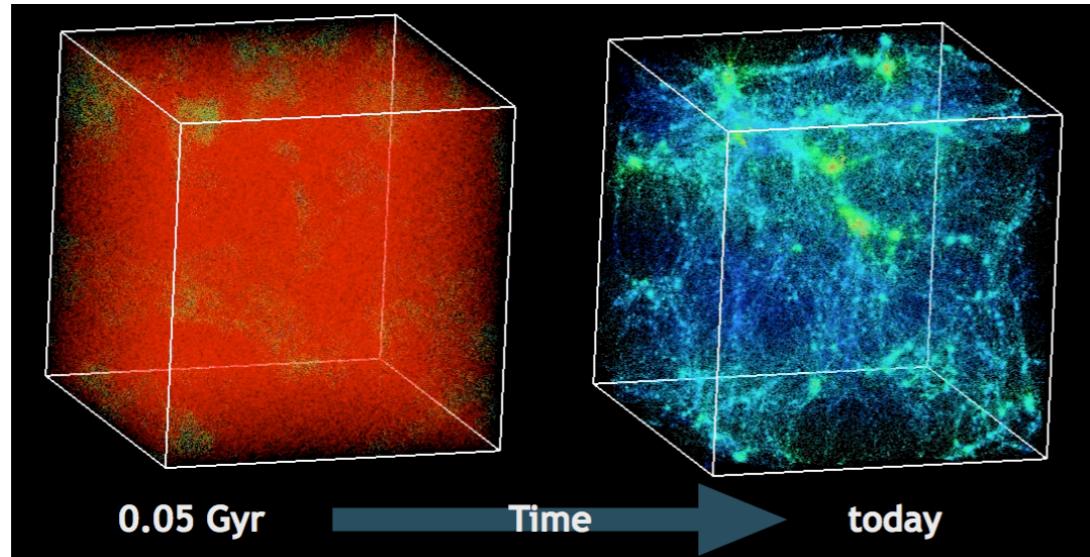
Numerical Recipes
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The Particle-In-Cell (PIC) method is a central tool for simulation over a wide range of physics studies

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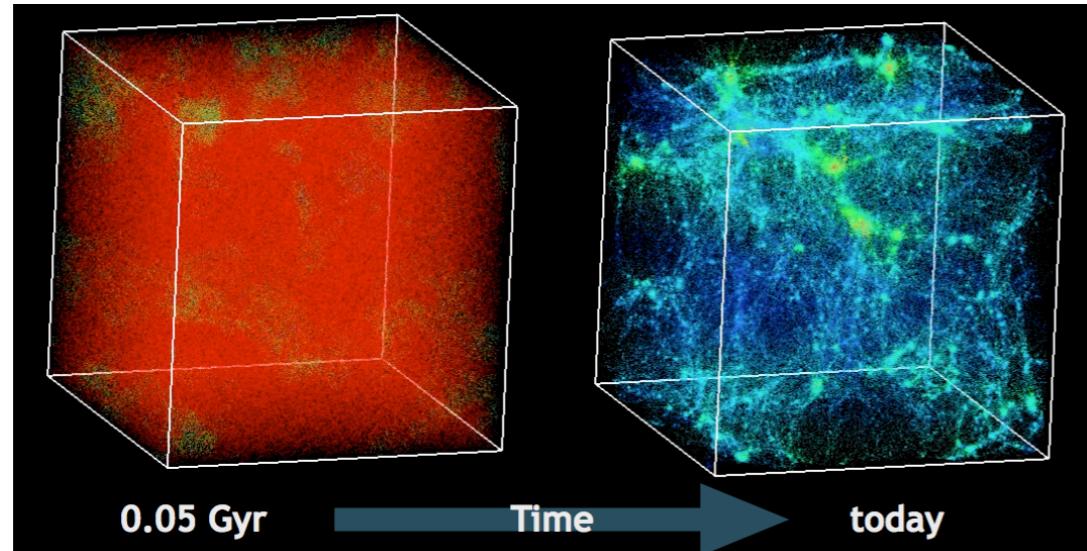
Cosmology



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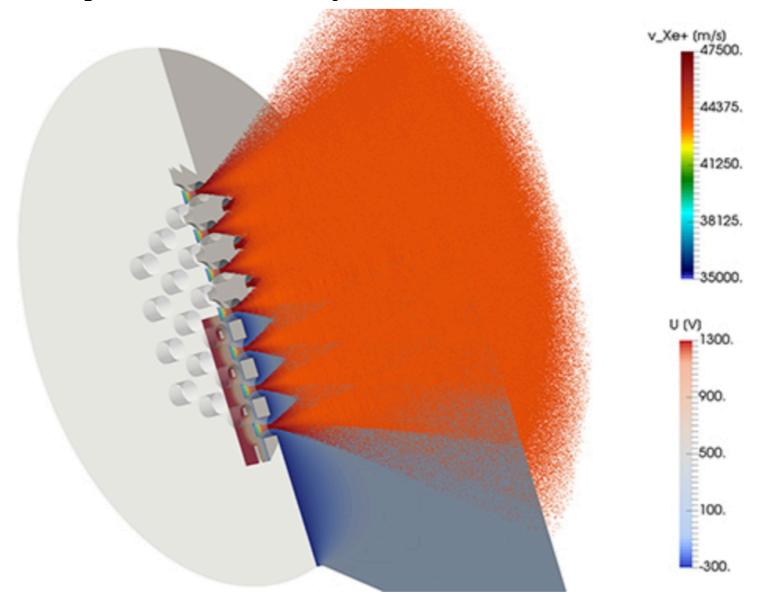
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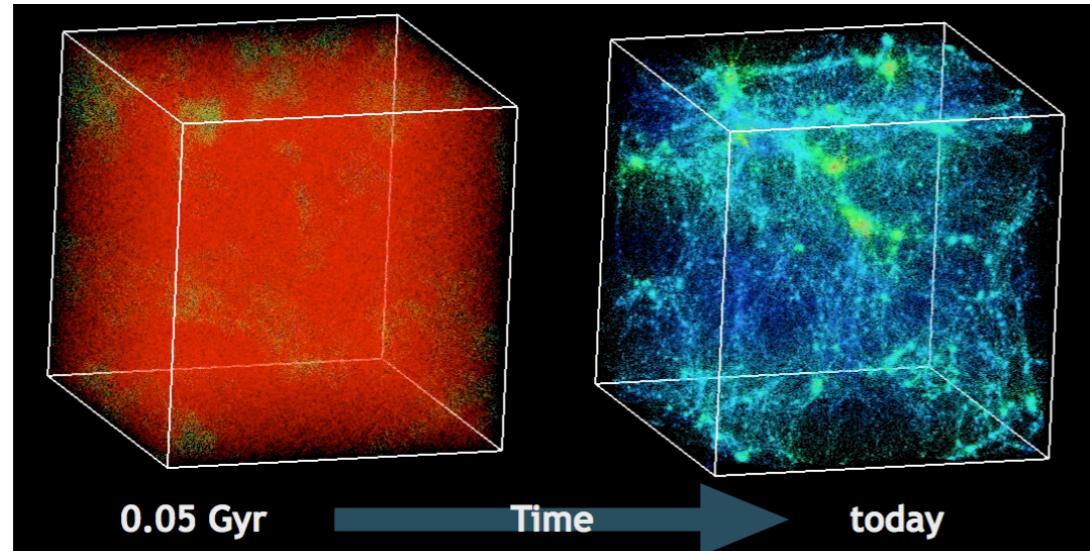
Space propulsion (Plasma thruster)



source: Gauss Center for Supercomputing

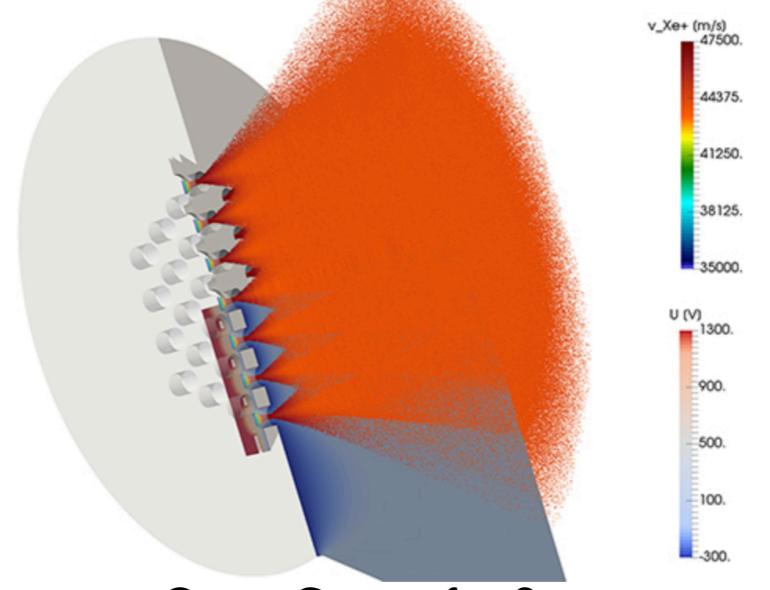
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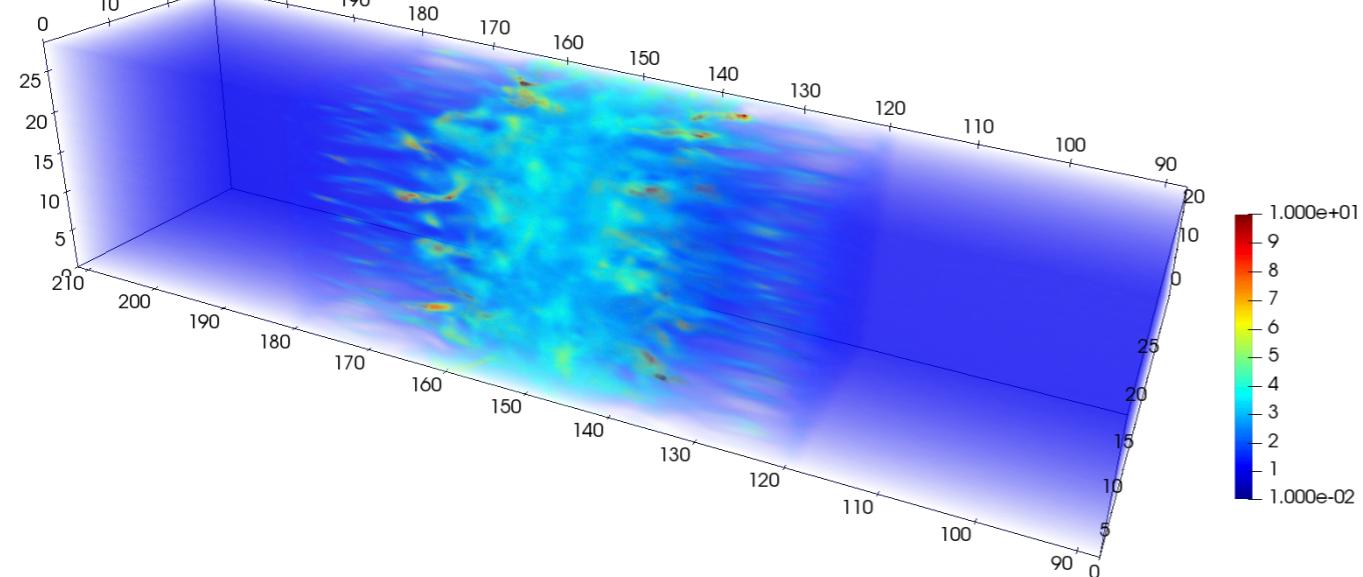
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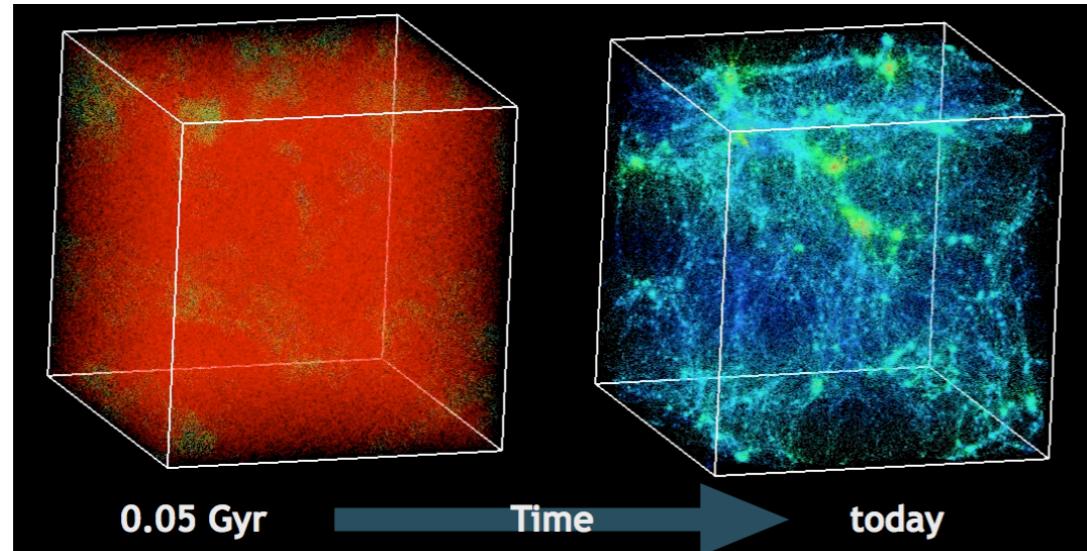
Space plasmas & astrophysics



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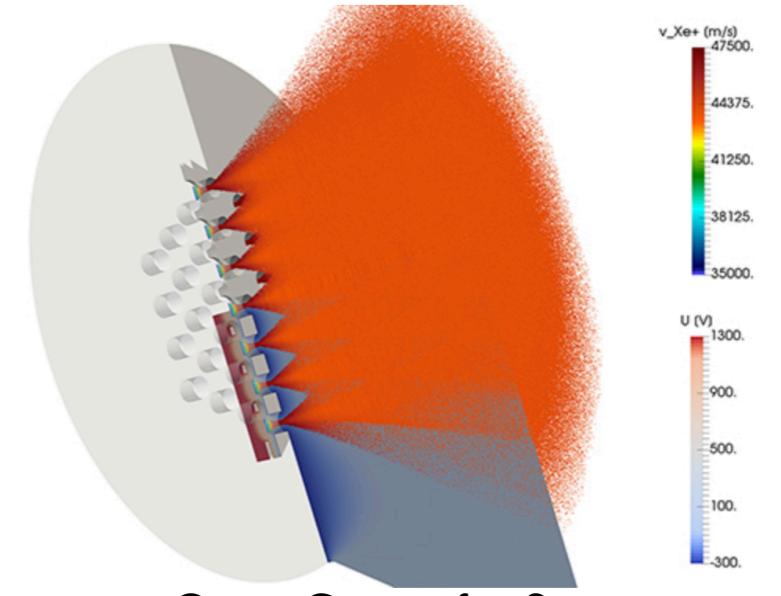
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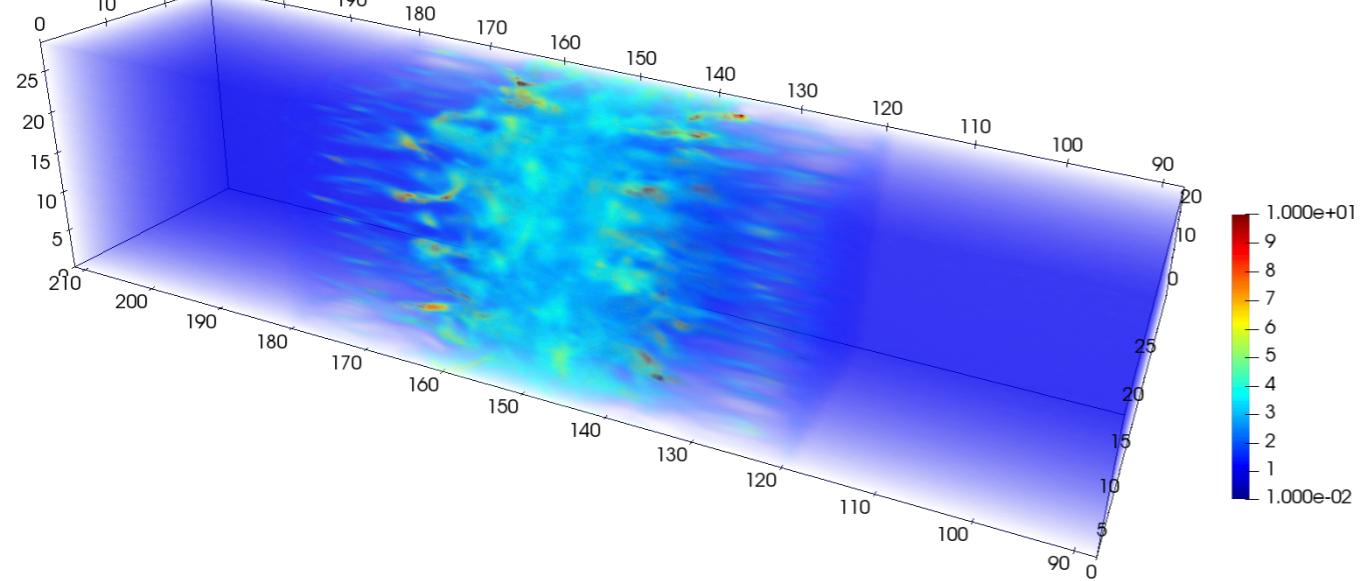
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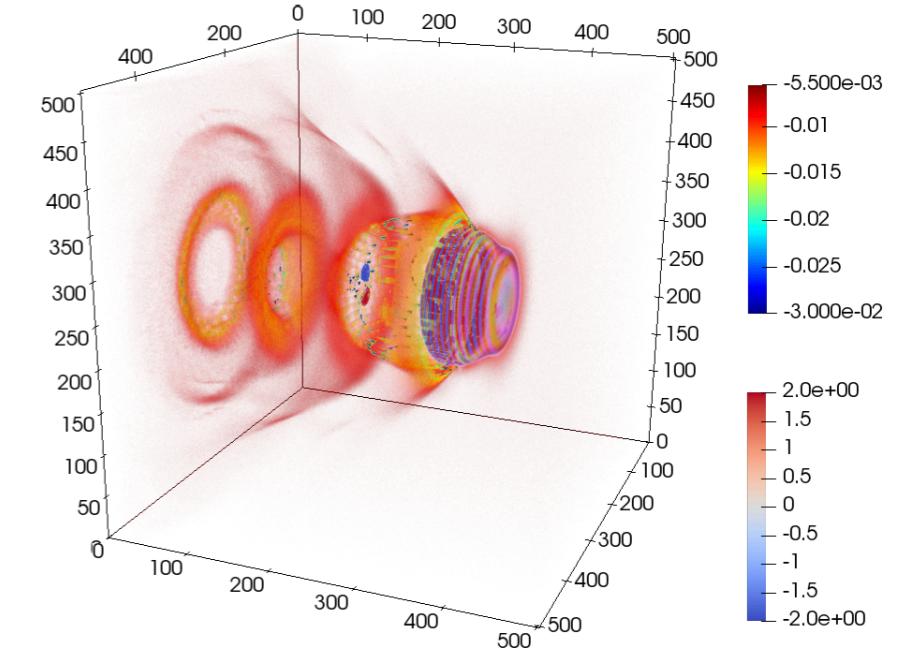
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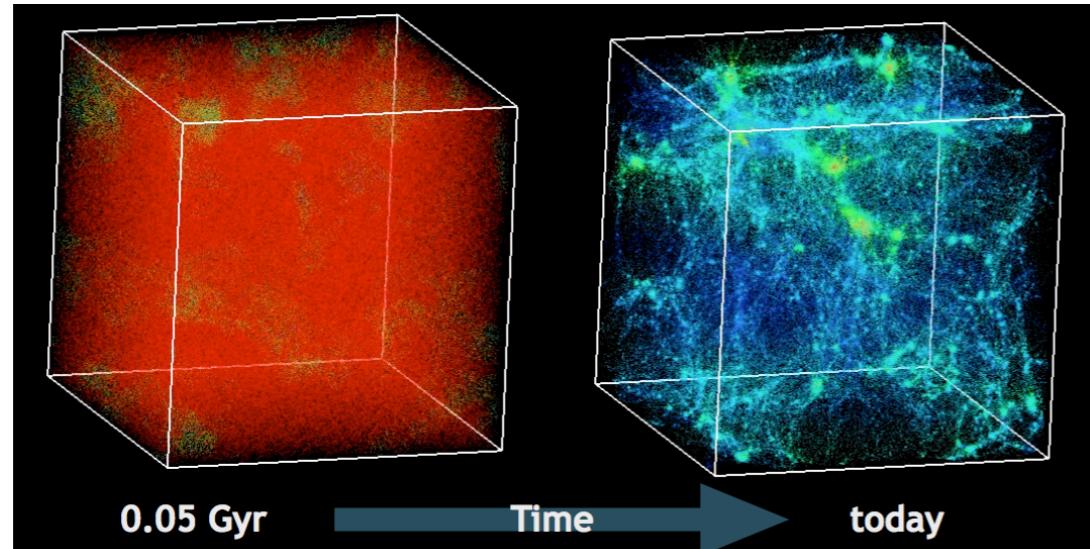
Laser plasma interaction



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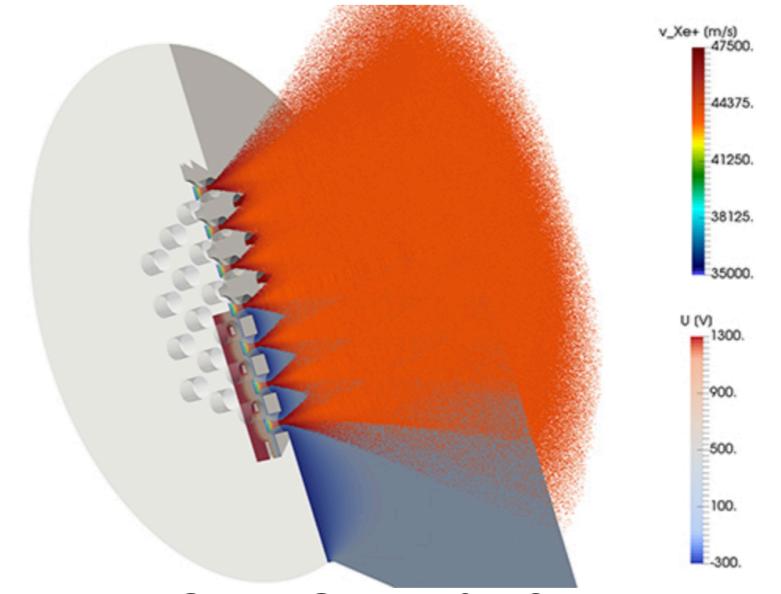
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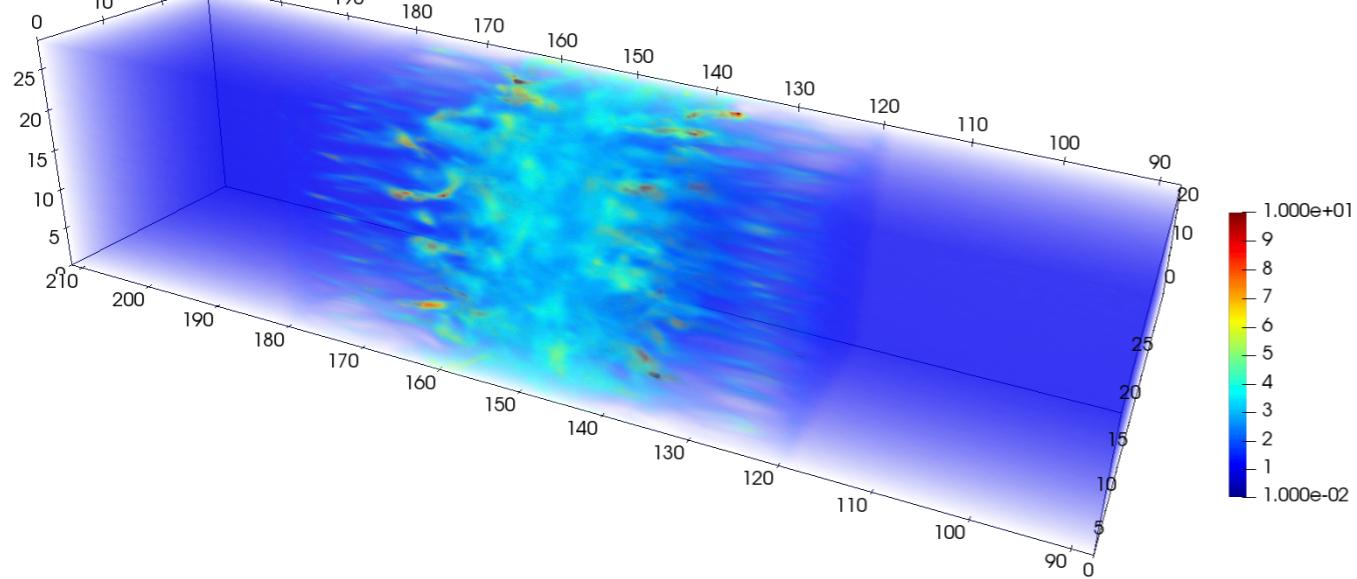
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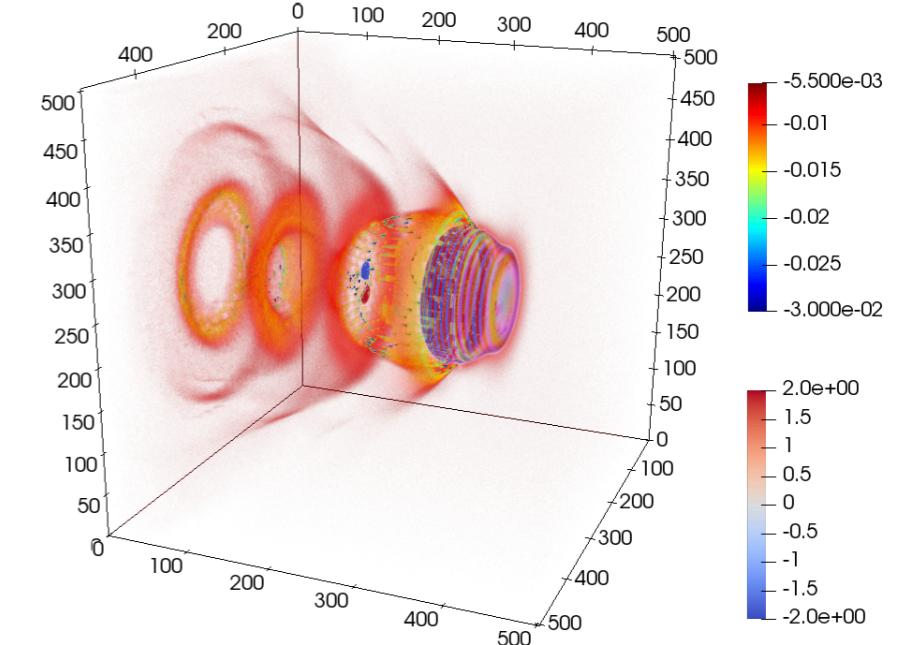
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- Conceptually simple
- Efficiently implemented on (massively) parallel super-computers

Our starting point is the Vlasov-Maxwell description
for a *collisionless* plasma

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$$\mathbf{F}_L = q_s \left(\mathbf{E} + \frac{\mathbf{p}}{m_s \gamma} \times \mathbf{B} \right)$$

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Velocity	c
Charge	e
Mass	m_e
Momentum	$m_e c$
Energy, Temperature	$m_e c^2$
Time	ω_r^{-1}
Length	c/ω_r
Number density	$n_r = \epsilon_0 m_e \omega_r^2 / e^2$
Current density	$e c n_r$
Pressure	$m_e c^2 n_r$
Electric field	$m_e c \omega_r / e$
Magnetic field	$m_e \omega_r / e$
Poynting flux	$m_e c^3 n_r / 2$

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The value of ω_r is not defined *a priori*, and acts as a scaling factor.

2nd Remark

The Particle-In-Cell method integrates Vlasov Equation along the trajectories of so-called *quasi-particles*

Vlasov Eq. is a **partial differential equation** (PDE) in $N_s + N_v$ phase-space:

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The **PIC ansatz** consists in decomposing the distribution fct:

$$f_s(t, \mathbf{x}, \mathbf{p}) = \sum_{p=1}^N w_p S(\mathbf{x} - \mathbf{x}_p(t)) \delta(\mathbf{p} - \mathbf{p}_p(t))$$

↑ ↑
Shape-function Dirac-distribution

2nd Remark

The Particle-In-Cell method **integrates Vlasov Equation along the trajectories of so-called quasi-particles**

Injecting this *ansatz* in Vlasov Eq., multiplying by \mathbf{p} and integrating over all momenta \mathbf{p}

$$\sum_{p=1}^{N_s} w_p \frac{\mathbf{p}_p}{m_s \gamma_p} \mathbf{p}_p \cdot [\partial_{\mathbf{x}_p} S(\mathbf{x} - \mathbf{x}_p) + \partial_{\mathbf{x}} S(\mathbf{x} - \mathbf{x}_p)] \\ + \sum_{p=1}^{N_s} w_p S(\mathbf{x} - \mathbf{x}_p) [\partial_t \mathbf{p}_p - q_s (\mathbf{E} + \mathbf{v}_p \times \mathbf{B})] = 0$$

Let us now integrate in space:

$$\sum_{p=1}^{N_s} w_p \frac{\mathbf{p}_p}{m_s \gamma_p} \mathbf{p}_p \cdot \cancel{\int d\mathbf{x} [\partial_{\mathbf{x}_p} S(\mathbf{x} - \mathbf{x}_p) + \partial_{\mathbf{x}} S(\mathbf{x} - \mathbf{x}_p)]} \\ + \sum_{p=1}^{N_s} w_p \cancel{\int d\mathbf{x}} S(\mathbf{x} - \mathbf{x}_p) [\partial_t \mathbf{p}_p - q_s (\mathbf{E} + \mathbf{v}_p \times \mathbf{B})] = 0$$

Finally leading to solving for all \mathbf{p} :

$$\partial_t \mathbf{p}_p = q_s (\mathbf{E}_p + \mathbf{v} \times \mathbf{B}_p) \quad \text{with} \quad (\mathbf{E}, \mathbf{B})_p \equiv \int d\mathbf{x} (\mathbf{E}, \mathbf{B})(\mathbf{x}) S(\mathbf{x} - \mathbf{x}_p)$$

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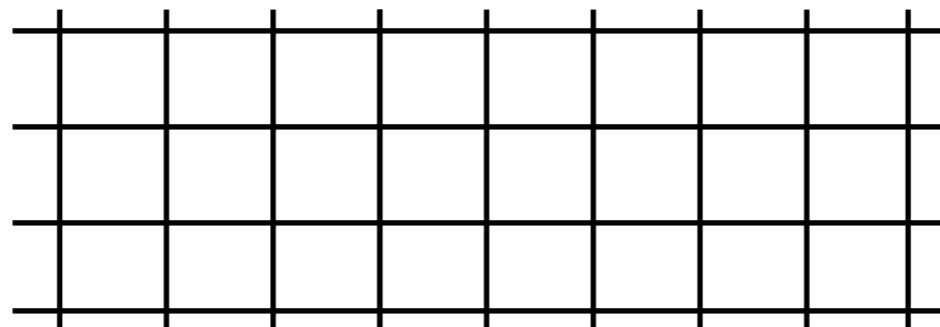
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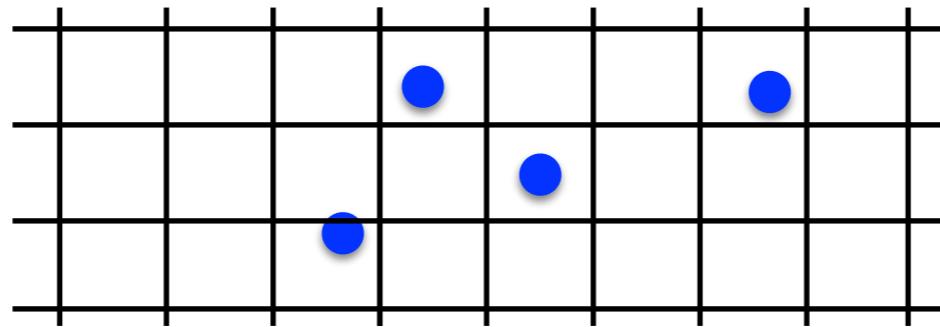
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If at time $t=0$, Poisson & Gauss Eqs. are satisfied, and if current deposition is made in a way that **conserve charge**, then solving **only Maxwell-Ampère & Maxwell-Faraday** ensures that both Eqs. remain satisfied at later time.

Initialization of a PIC simulation

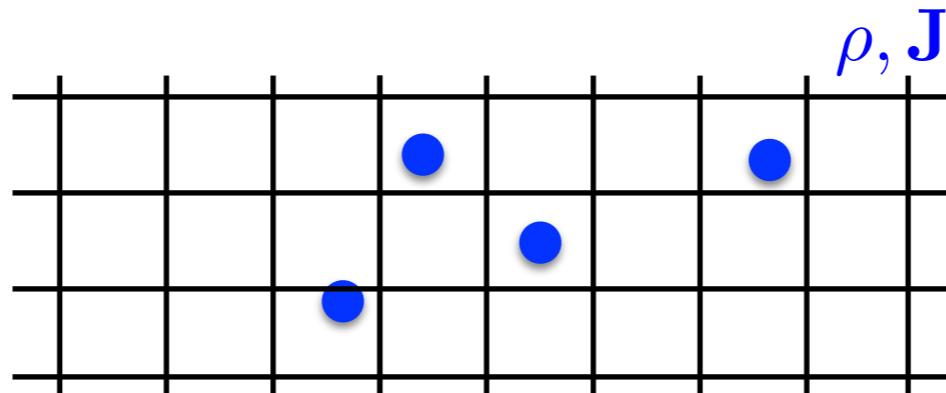


Initialization of a PIC simulation



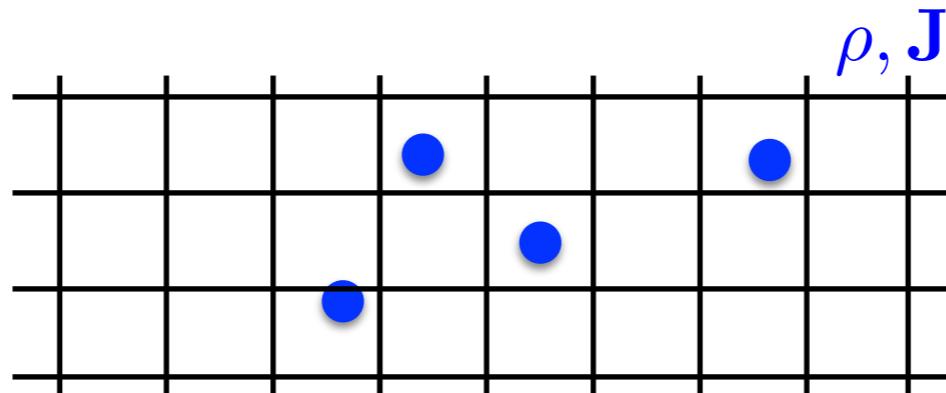
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e.g. defining the species density, velocity and temperature profiles

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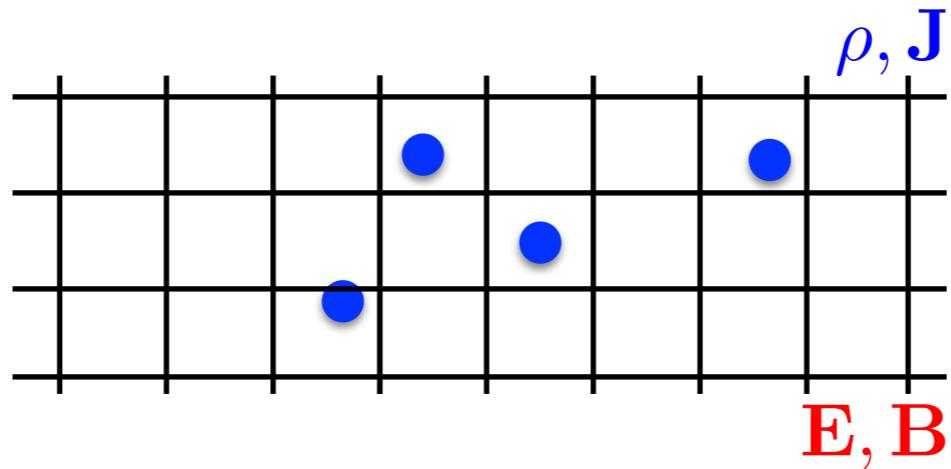
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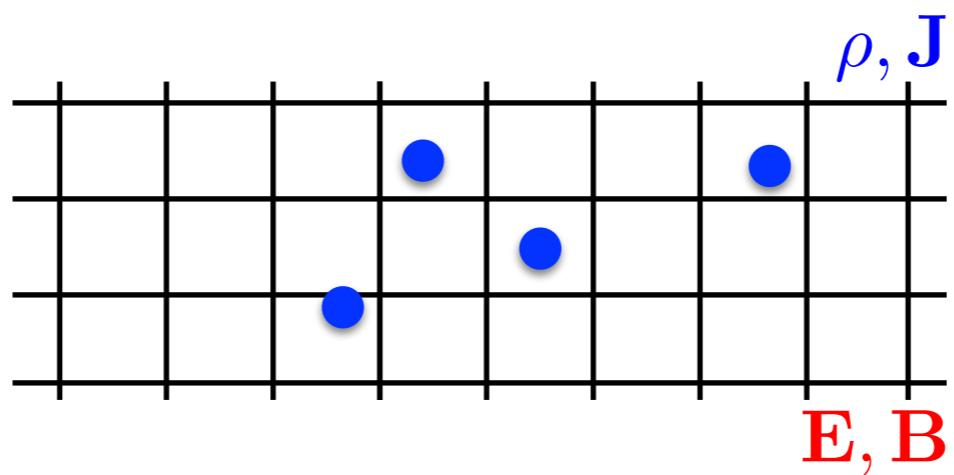
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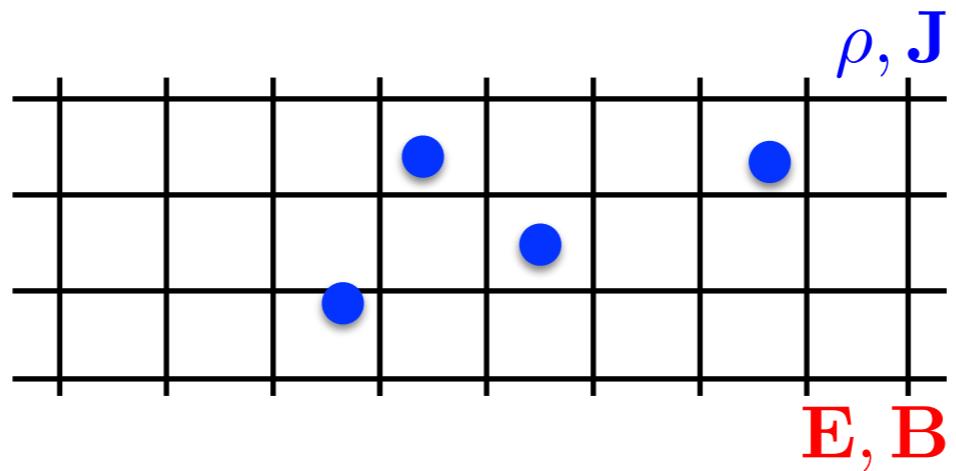


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- 4) add any (user defined) external fields provided they are divergence-free

The Particle-In-Cell loop



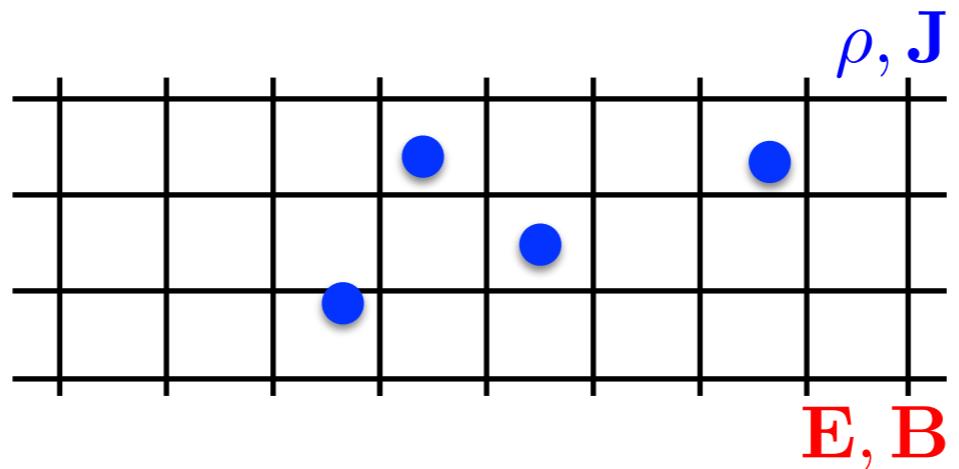
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Gather fields at particle position

$$[\mathbf{E}, \mathbf{B}] \rightarrow [\mathbf{E}_p, \mathbf{B}_p]$$

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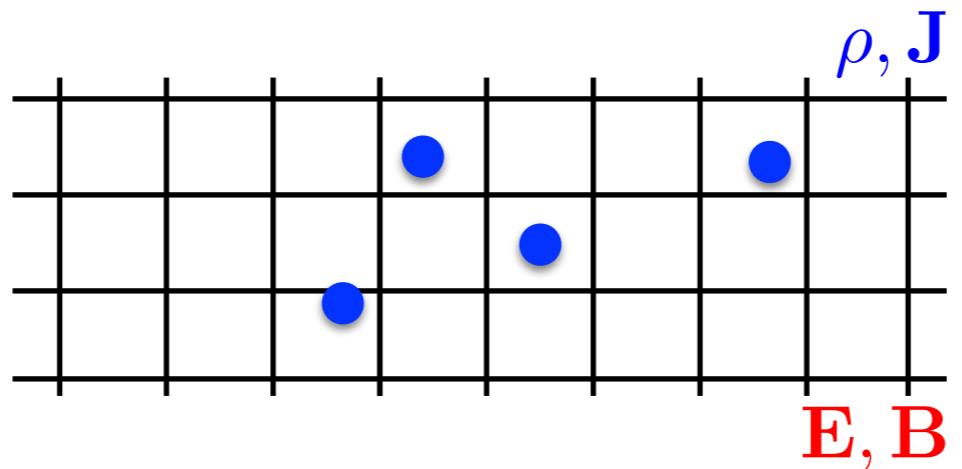
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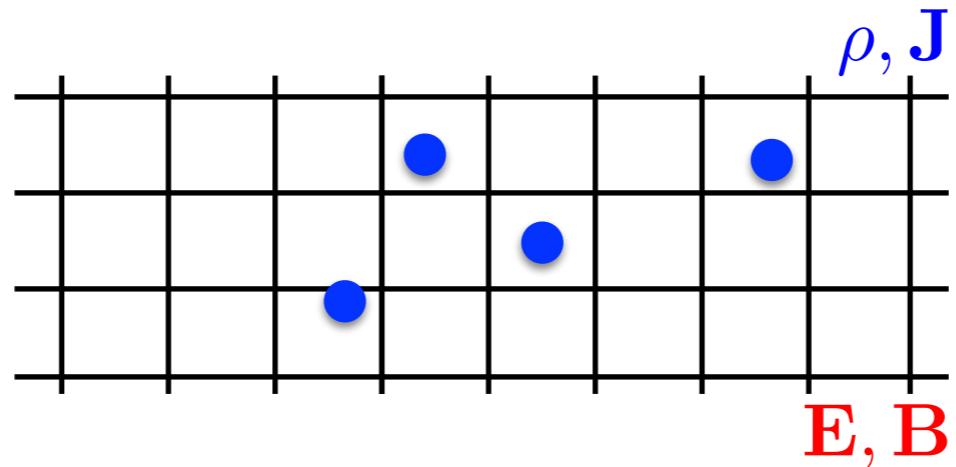
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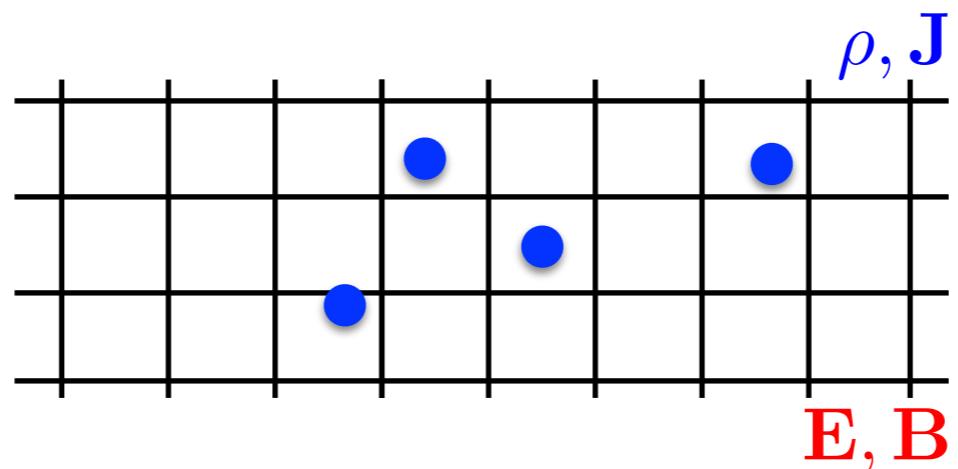
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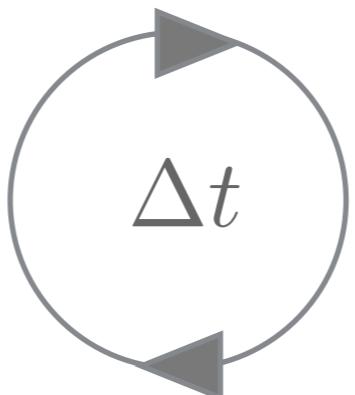
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Outlines

- Numerical approach: how to build up your PIC code
- Parallelization: getting ready for the super-computers
- Additional modules: beyond the *collisionless* plasma
- Some physics highlights: what you can do with a PIC code

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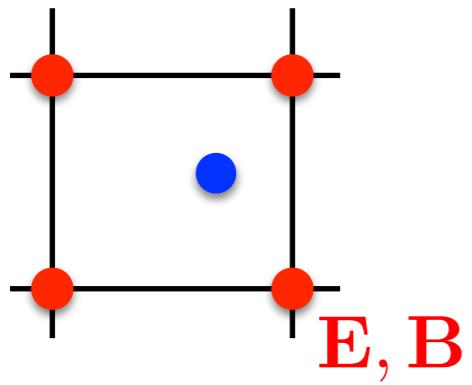
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Step 1

Field gathering: interpolation at particle position

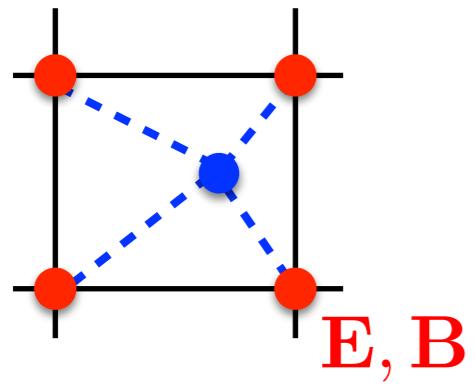
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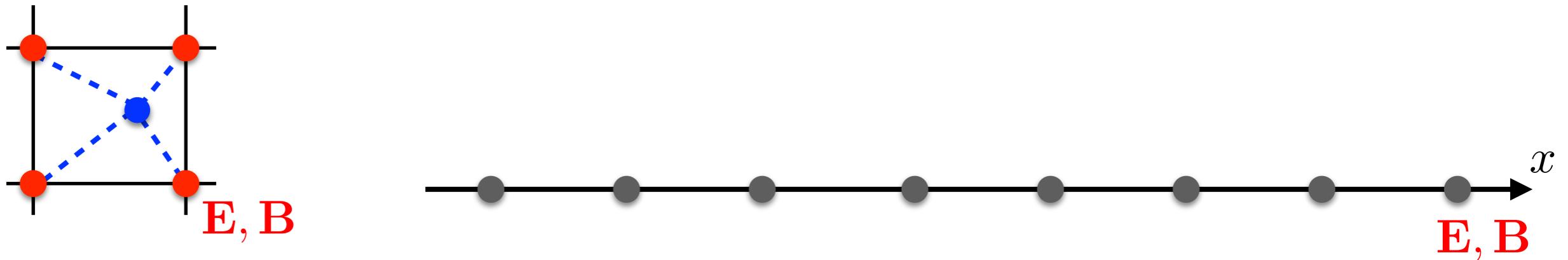
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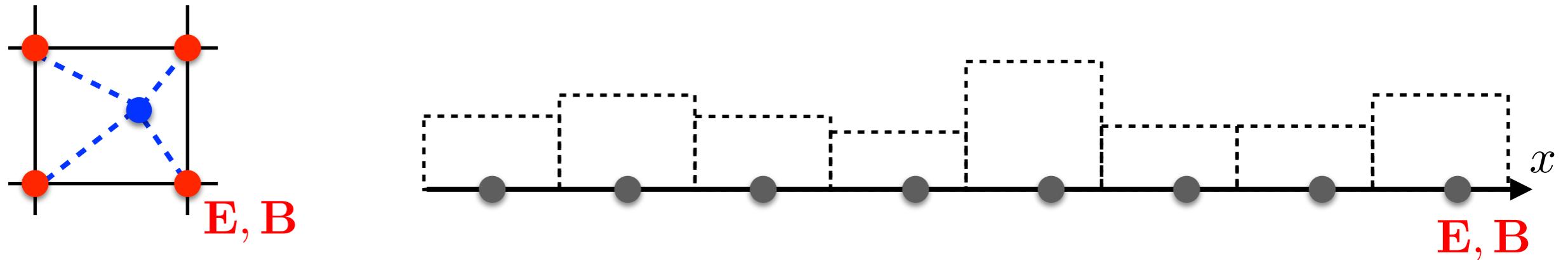
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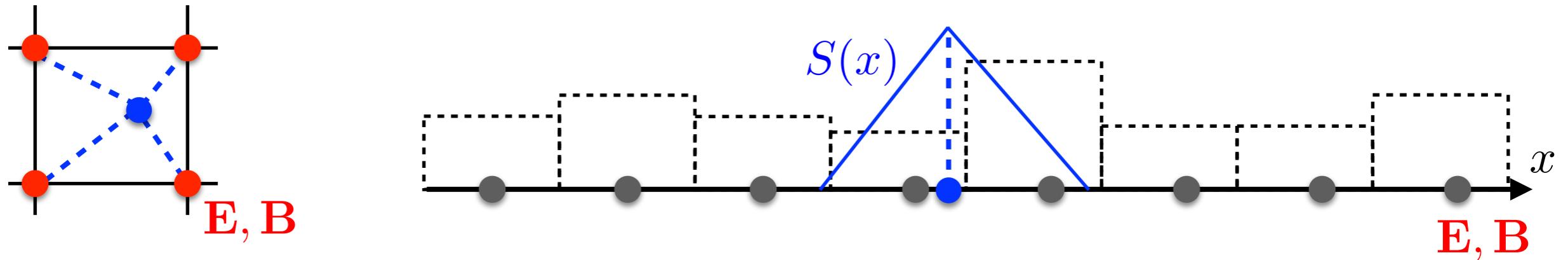
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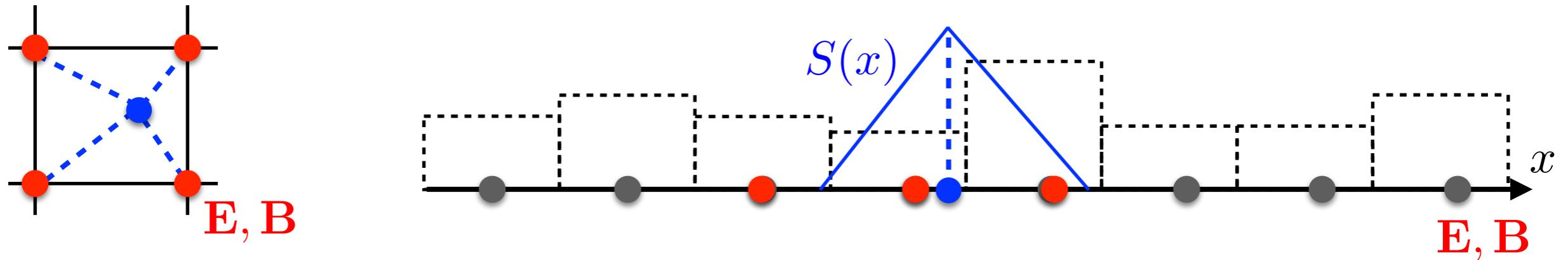
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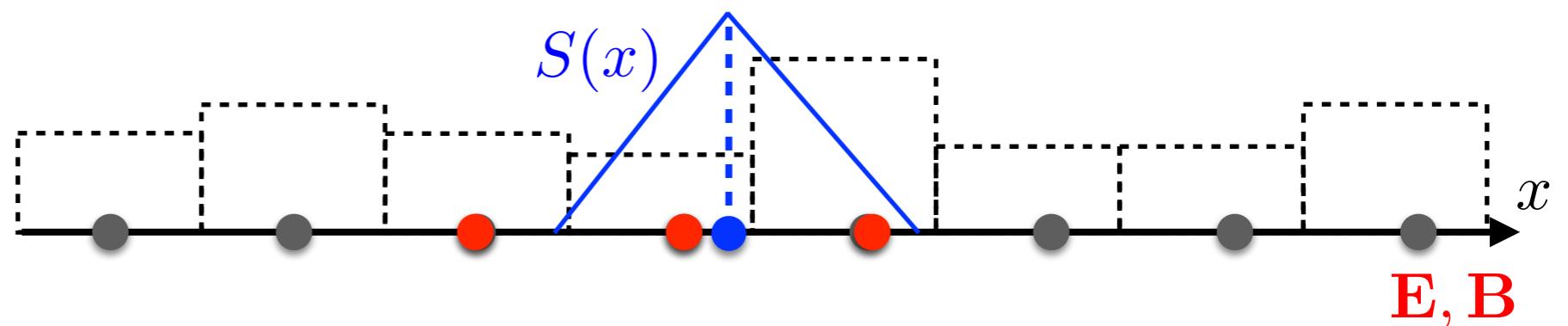
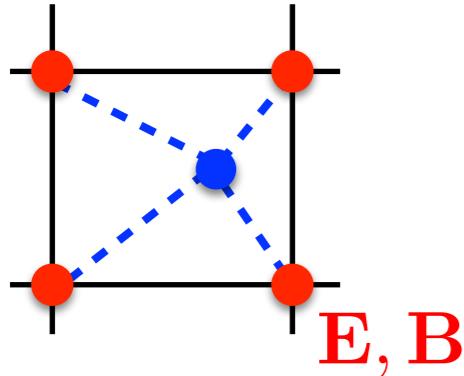
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Field gathering: interpolation at particle position

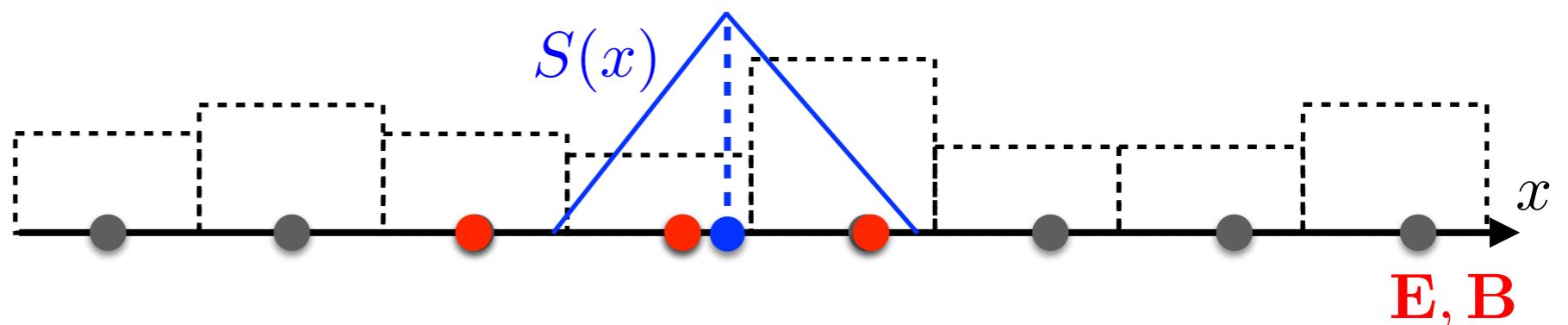
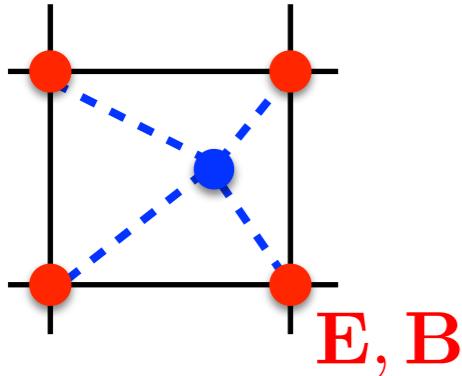


$$(\mathbf{E}, \mathbf{B})_p \equiv \int d\mathbf{x} (\mathbf{E}, \mathbf{B})(\mathbf{x}) S(\mathbf{x} - \mathbf{x}_p)$$

$$\begin{aligned} \hat{s}^{(0)}(x) &= \Delta x \delta(x), \\ \hat{s}^{(1)}(x) &= \begin{cases} 1 & \text{if } |x| \leq \frac{1}{2} \Delta x, \\ 0 & \text{otherwise,} \end{cases} \\ \hat{s}^{(2)}(x) &= \begin{cases} \left(1 - \left|\frac{x}{\Delta x}\right|\right) & \text{if } |x| \leq \Delta x, \\ 0 & \text{otherwise,} \end{cases} \\ \hat{s}^{(3)}(x) &= \begin{cases} \frac{3}{4} \left[1 - \frac{4}{3} \left(\frac{x}{\Delta x}\right)^2\right] & \text{if } |x| \leq \frac{1}{2} \Delta x, \\ \frac{9}{8} \left(1 - \frac{2}{3} \left|\frac{x}{\Delta x}\right|\right)^2 & \text{if } \frac{1}{2} \Delta x < |x| \leq \frac{3}{2} \Delta x, \\ 0 & \text{otherwise,} \end{cases} \\ \hat{s}^{(4)}(x) &= \begin{cases} \frac{2}{3} \left[1 - \frac{3}{2} \left(\frac{x}{\Delta x}\right)^2 + \frac{3}{4} \left|\frac{x}{\Delta x}\right|^3\right] & \text{if } |x| \leq \Delta x, \\ \frac{4}{3} \left(1 - \frac{1}{2} \left|\frac{x}{\Delta x}\right|\right)^3 & \text{if } \Delta x < |x| \leq 2 \Delta x, \\ 0 & \text{otherwise.} \end{cases} \end{aligned}$$

Step 1

Field gathering: interpolation at particle position



$$\hat{s}^{(0)}(x) = \Delta x \delta(x),$$

$$\hat{s}^{(1)}(x) = \begin{cases} 1 & \text{if } |x| \leq \frac{1}{2} \Delta x, \\ 0 & \text{otherwise,} \end{cases}$$

$$\hat{s}^{(2)}(x) = \begin{cases} \left(1 - \left|\frac{x}{\Delta x}\right|\right) & \text{if } |x| \leq \Delta x, \\ 0 & \text{otherwise,} \end{cases}$$

$$\hat{s}^{(3)}(x) = \begin{cases} \frac{3}{4} \left[1 - \frac{4}{3} \left(\frac{x}{\Delta x}\right)^2 \right] & \text{if } |x| \leq \frac{1}{2} \Delta x, \\ \frac{9}{8} \left(1 - \frac{2}{3} \left|\frac{x}{\Delta x}\right|\right)^2 & \text{if } \frac{1}{2} \Delta x < |x| \leq \frac{3}{2} \Delta x, \\ 0 & \text{otherwise,} \end{cases}$$

$$\hat{s}^{(4)}(x) = \begin{cases} \frac{2}{3} \left[1 - \frac{3}{2} \left(\frac{x}{\Delta x}\right)^2 + \frac{3}{4} \left|\frac{x}{\Delta x}\right|^3 \right] & \text{if } |x| \leq \Delta x, \\ \frac{4}{3} \left(1 - \frac{1}{2} \left|\frac{x}{\Delta x}\right|\right)^3 & \text{if } \Delta x < |x| \leq 2 \Delta x, \\ 0 & \text{otherwise.} \end{cases}$$

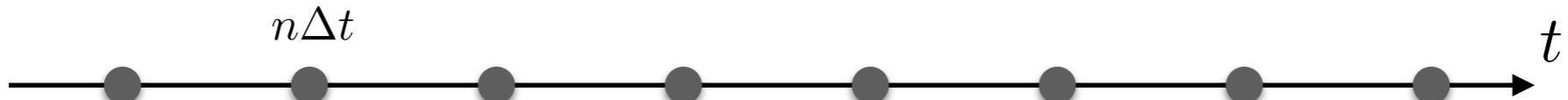
$$(\mathbf{E}, \mathbf{B})_p \equiv \int d\mathbf{x} (\mathbf{E}, \mathbf{B})(\mathbf{x}) S(\mathbf{x} - \mathbf{x}_p)$$

Step 2

The Boris leap-frog *pusher* is a very popular method to advance particles

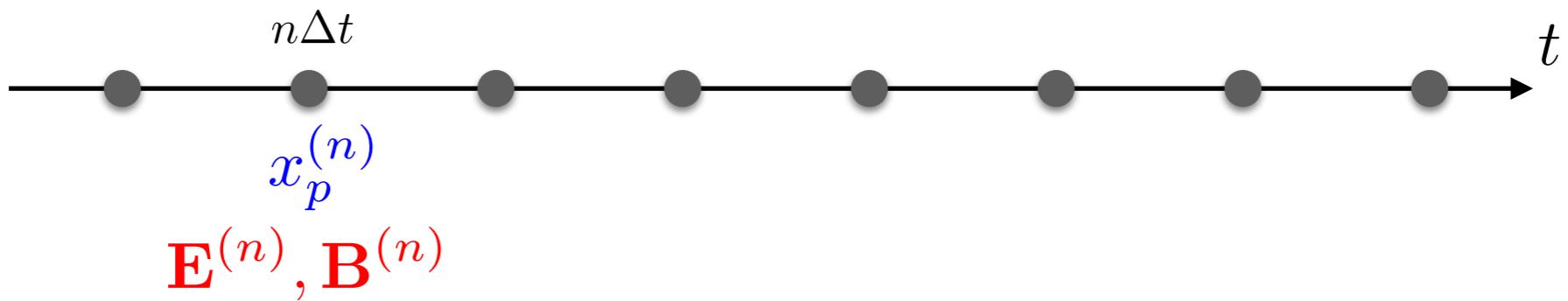
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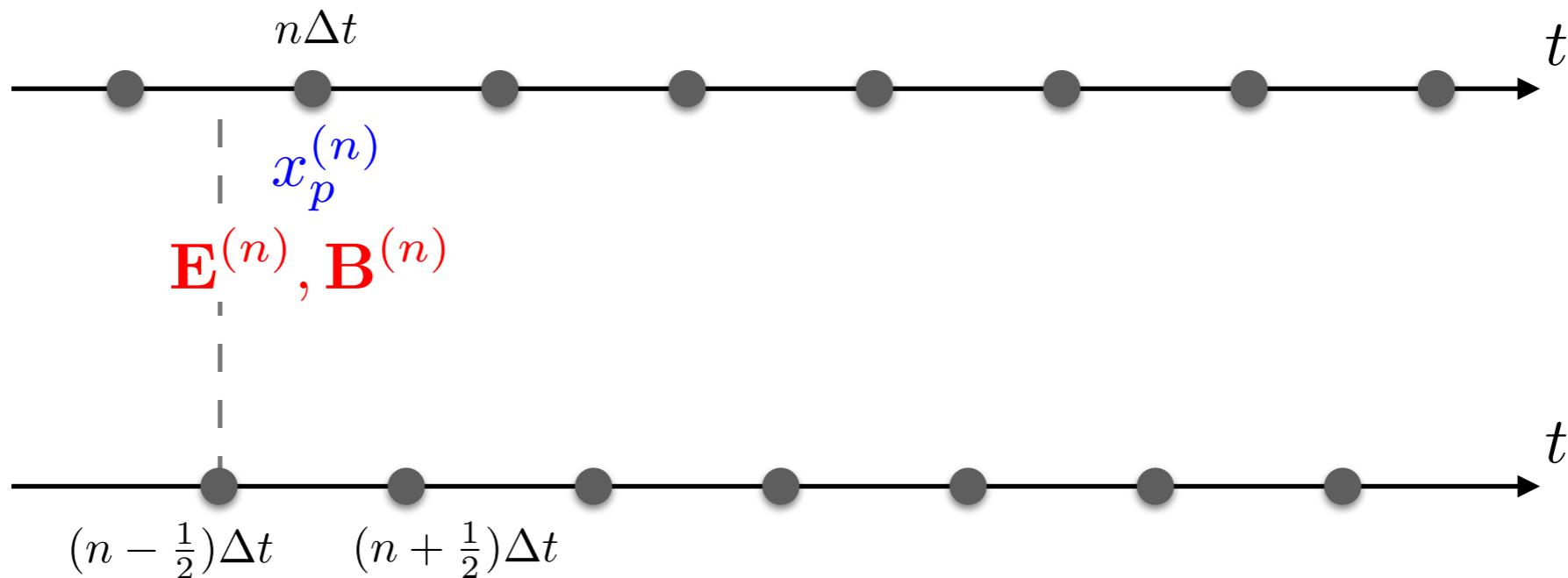
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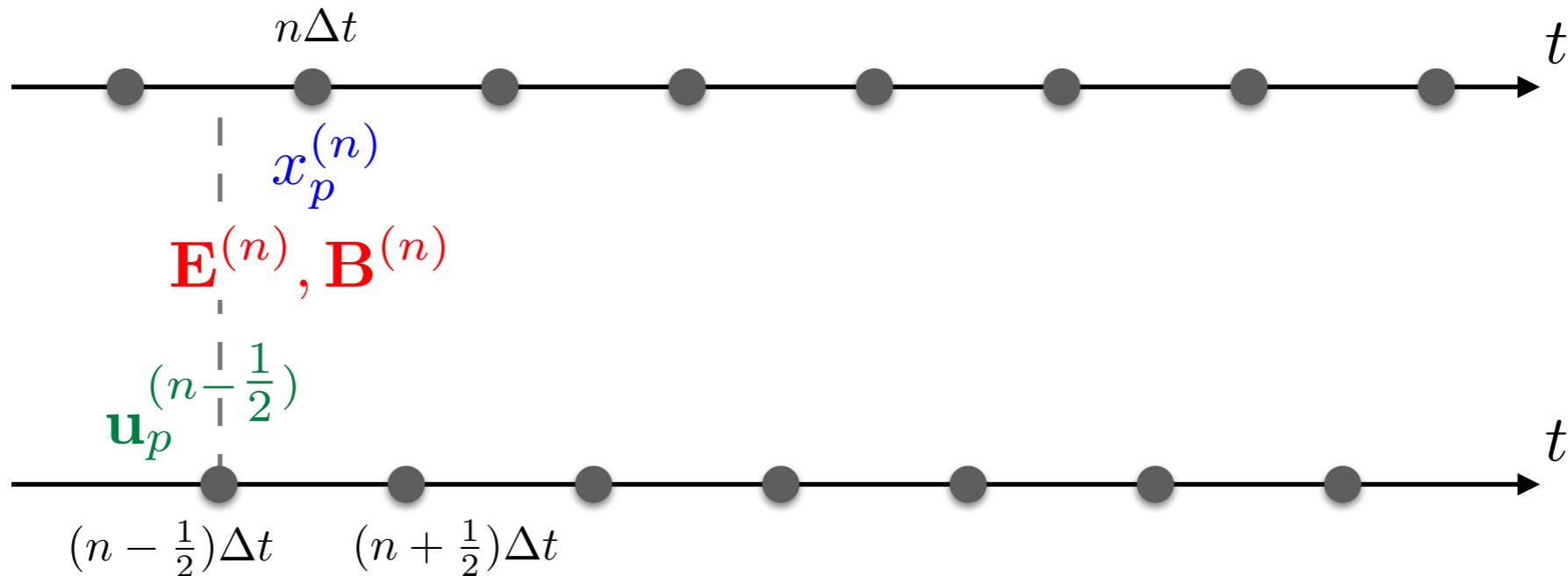
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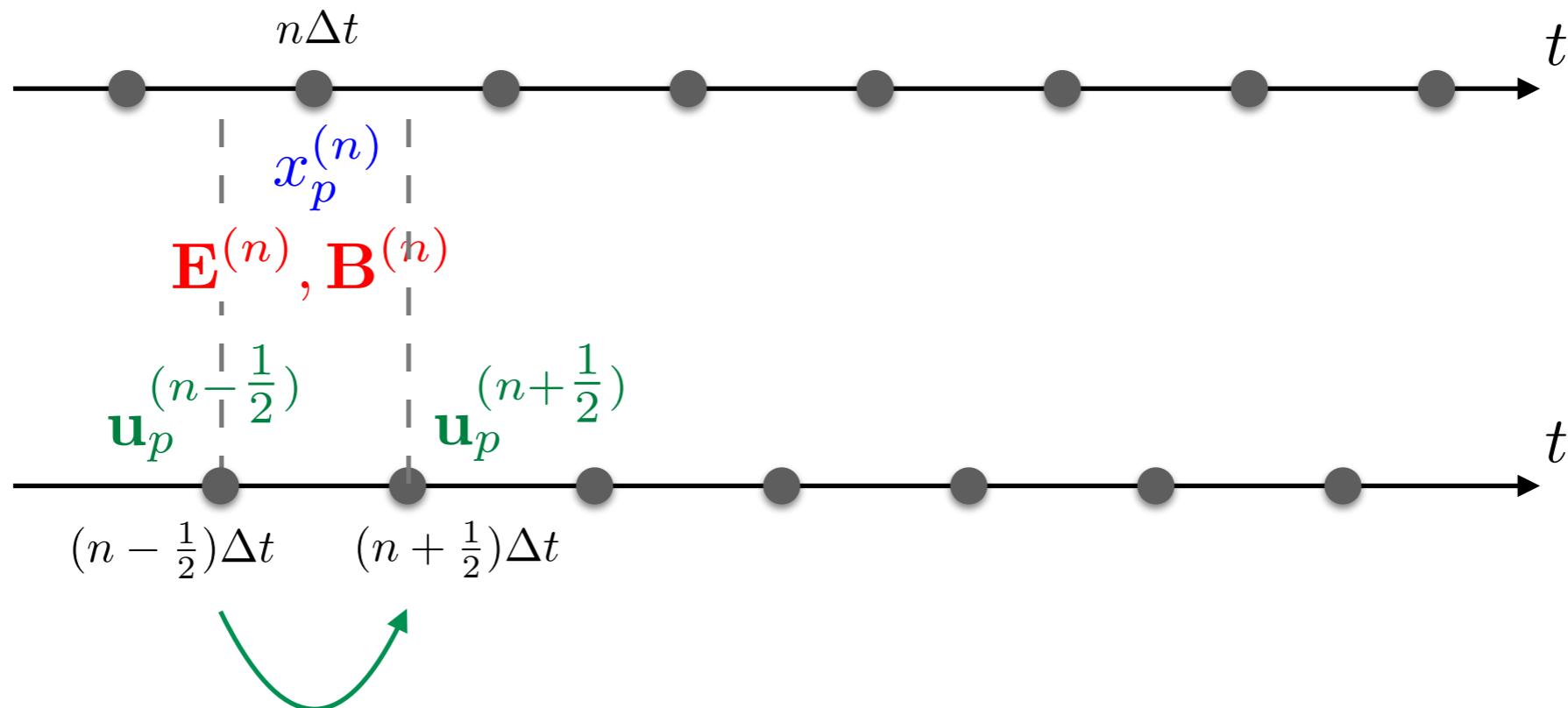
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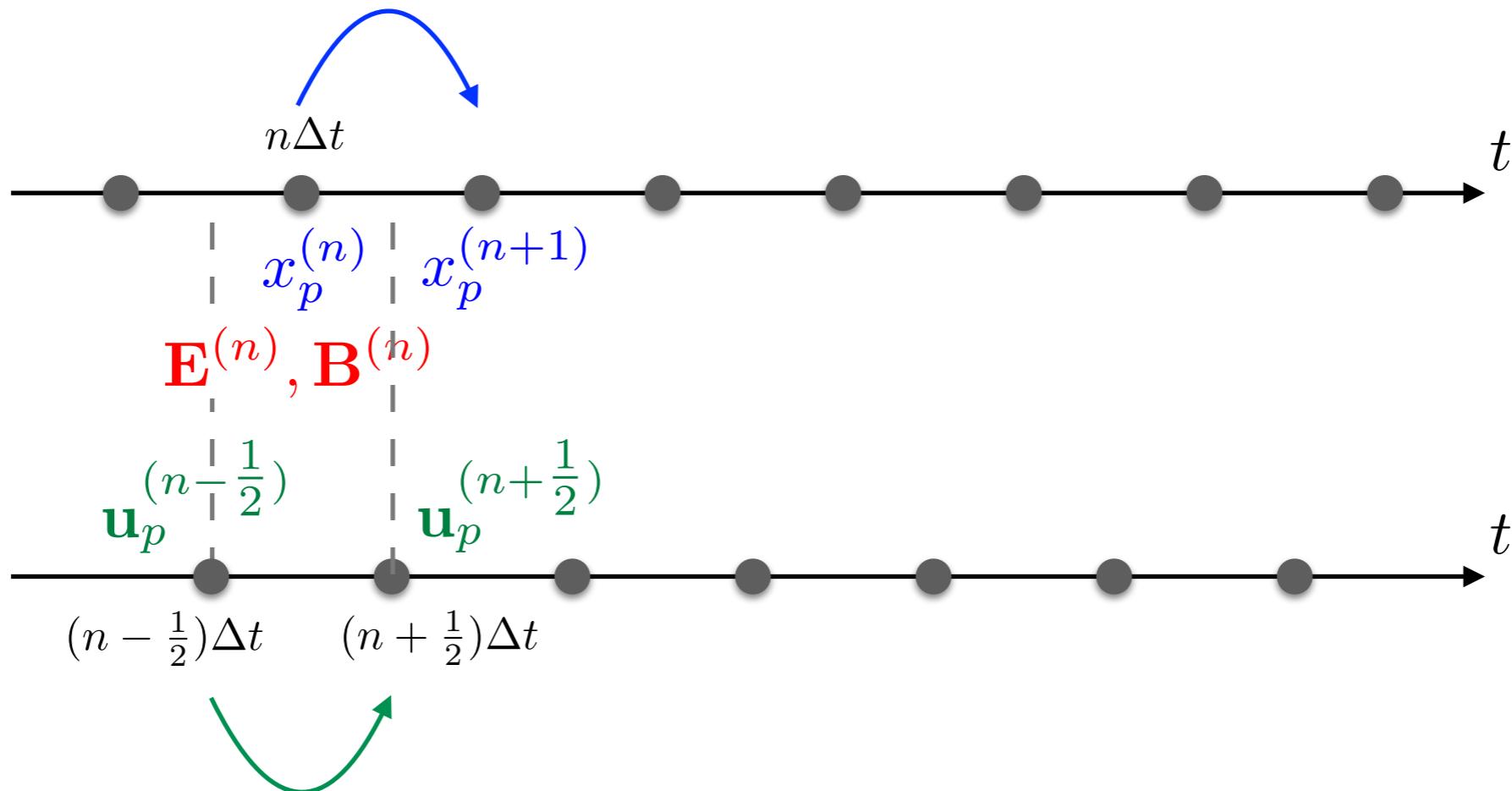
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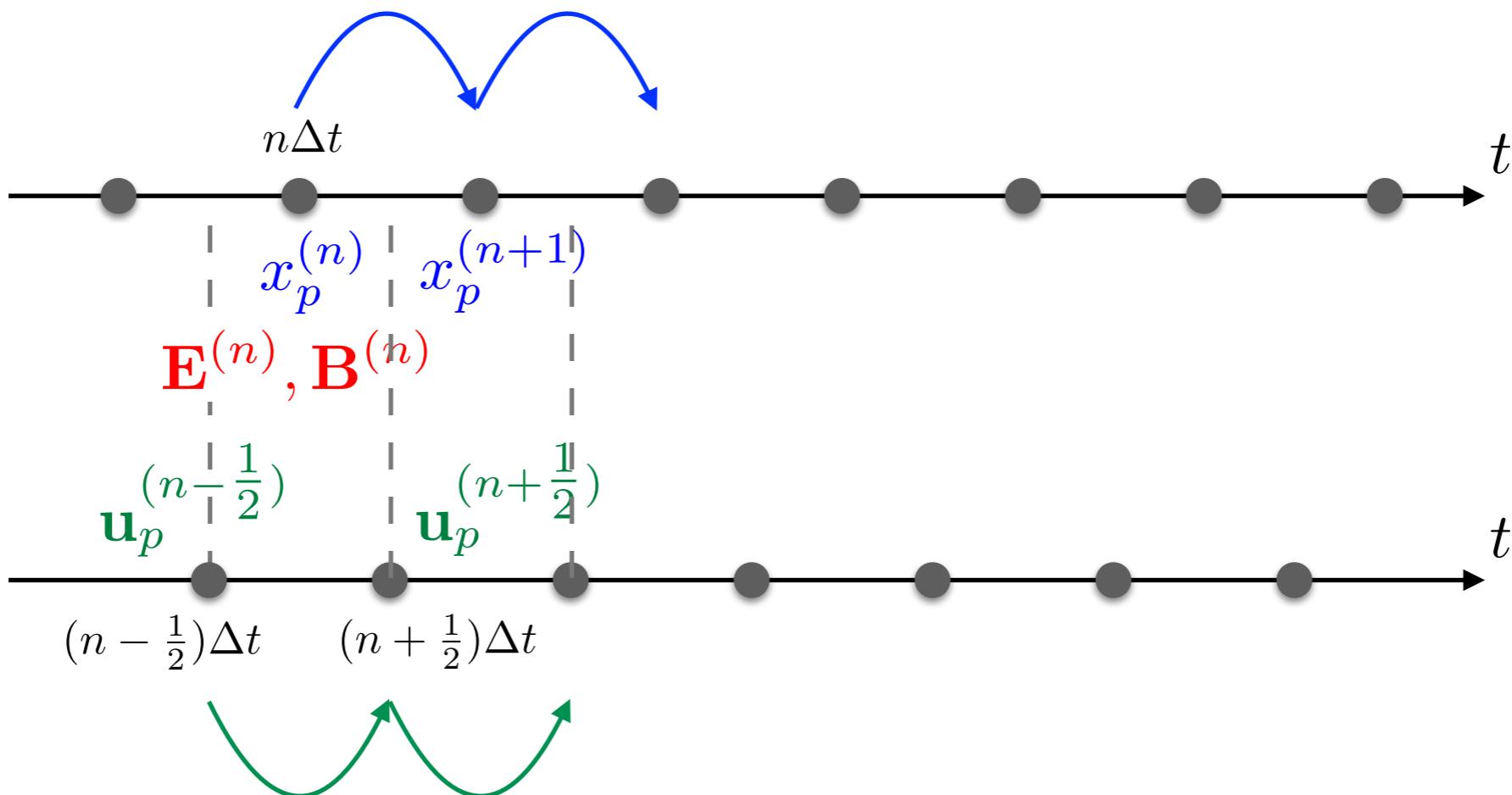
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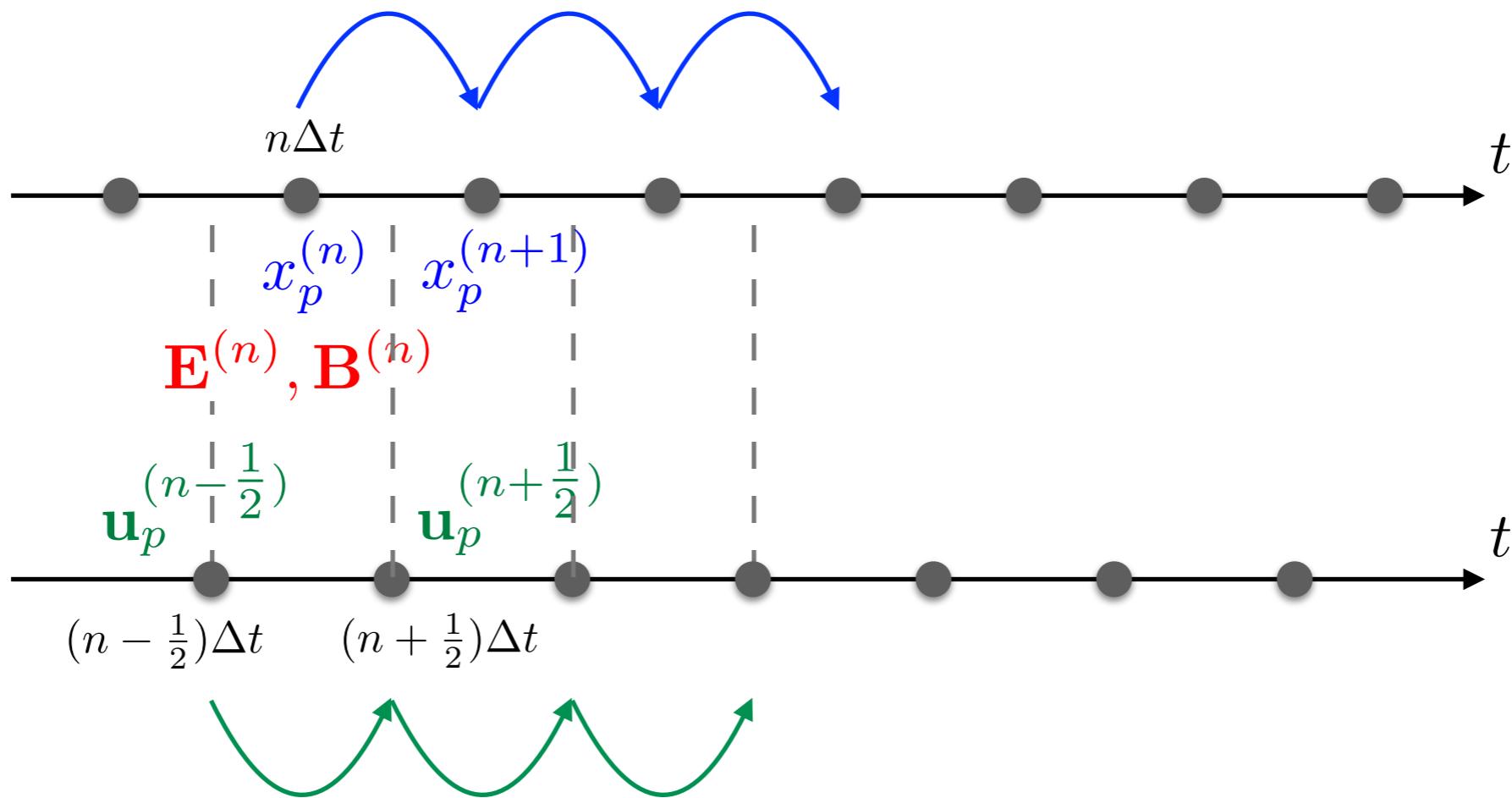
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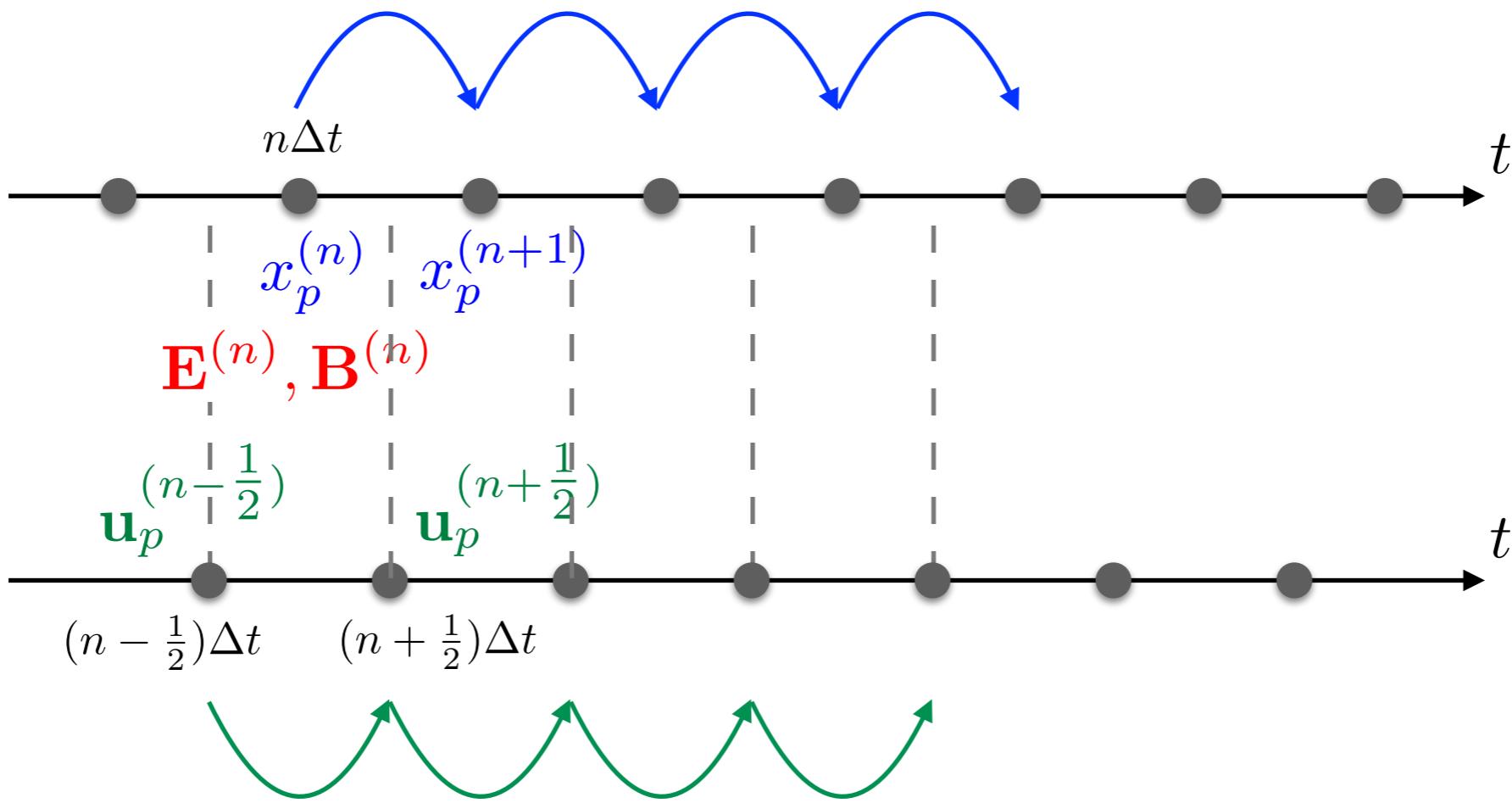
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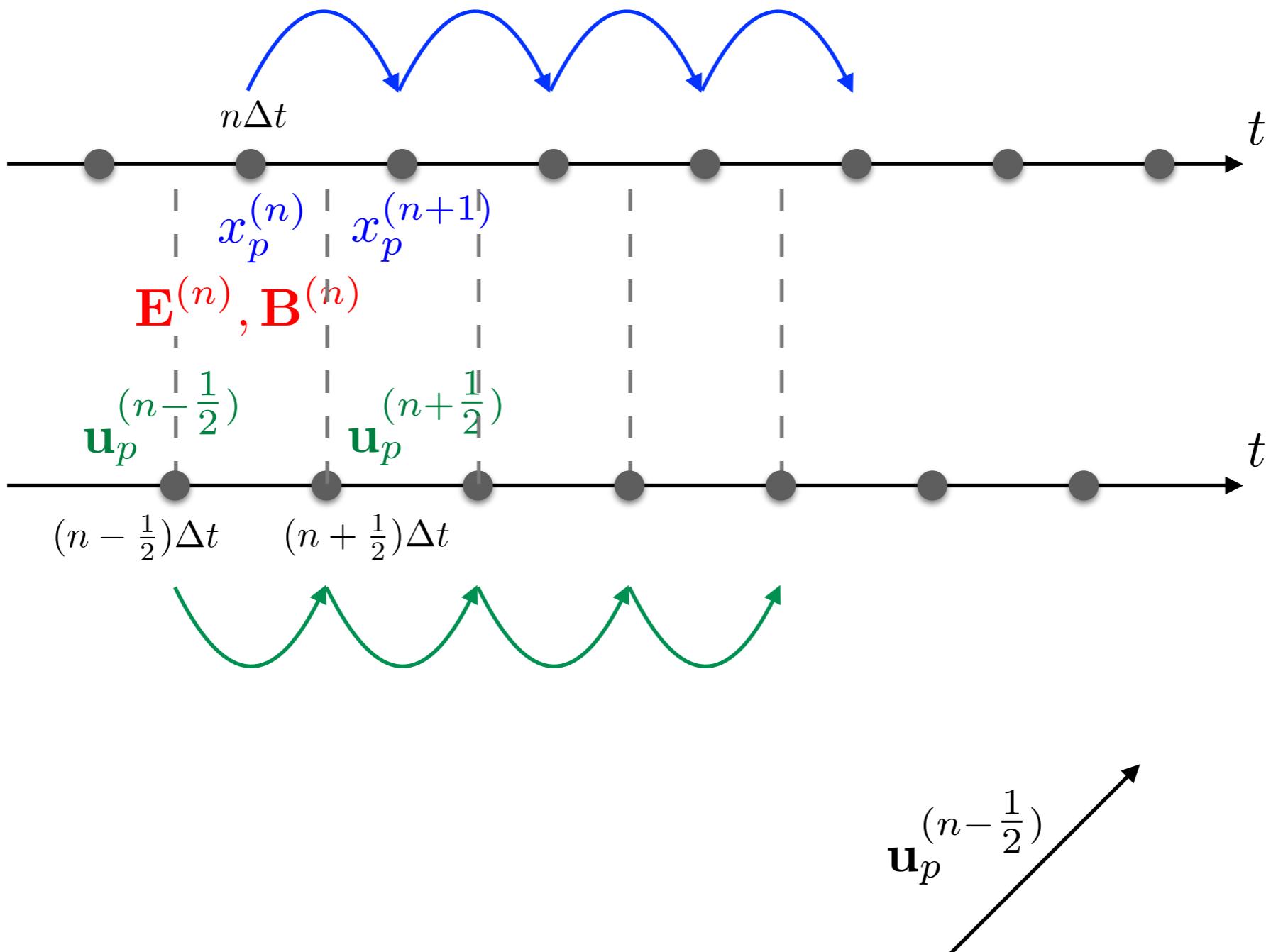
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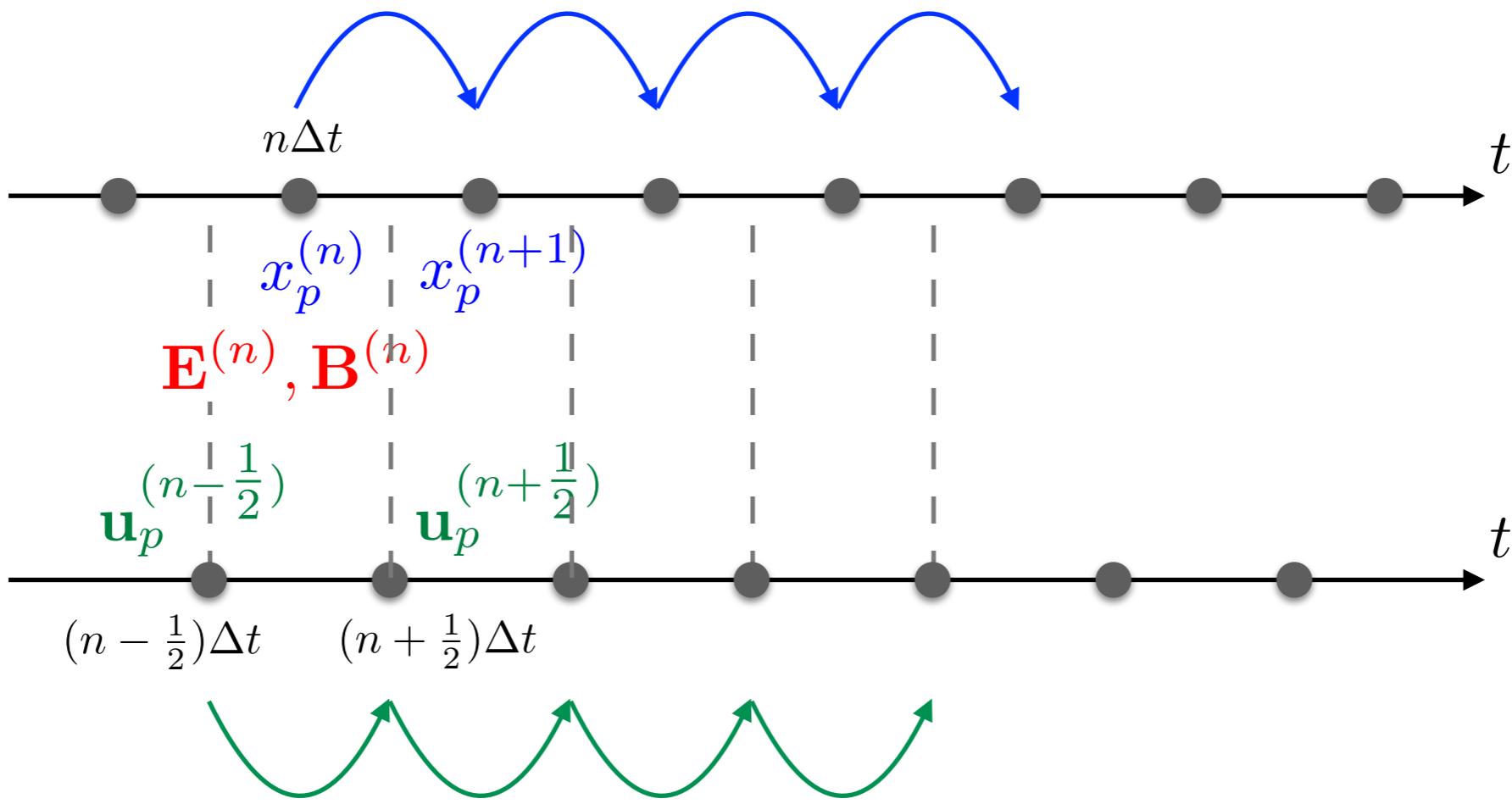
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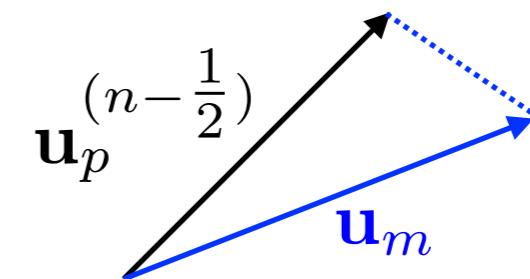


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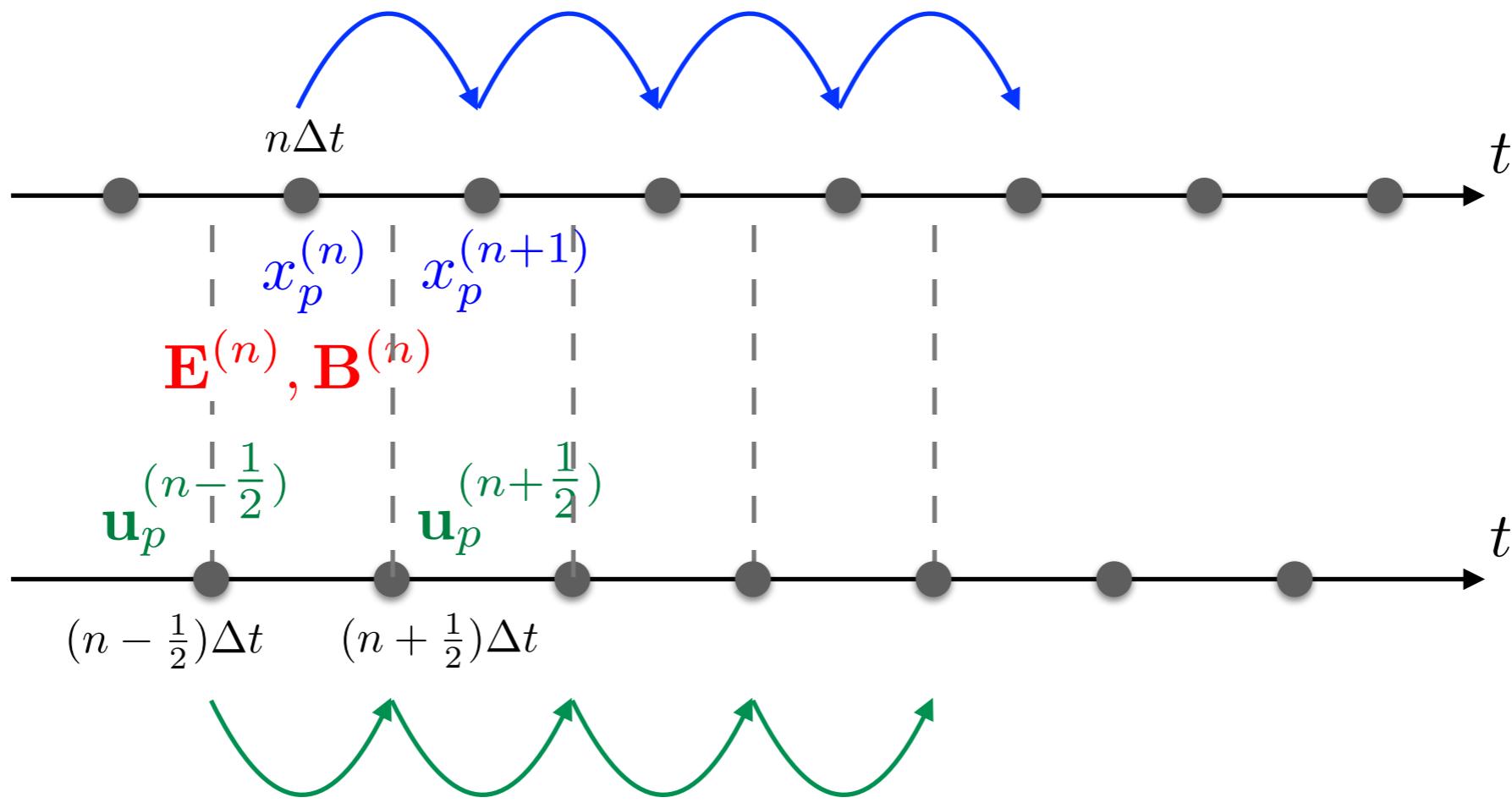


$$\mathbf{u}_m = \mathbf{u}_p^{(n-\frac{1}{2})} + \frac{q_s}{m_s} \frac{\Delta t}{2} \mathbf{E}_p$$



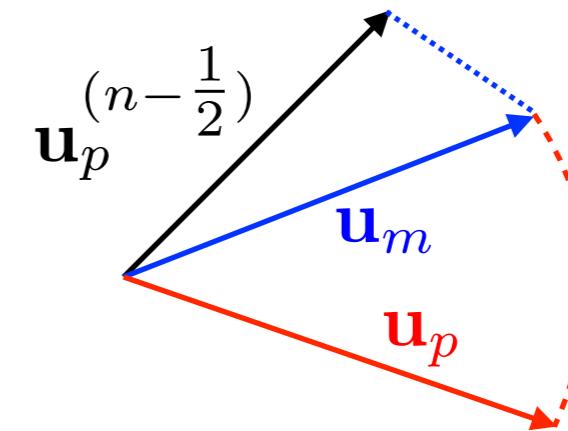
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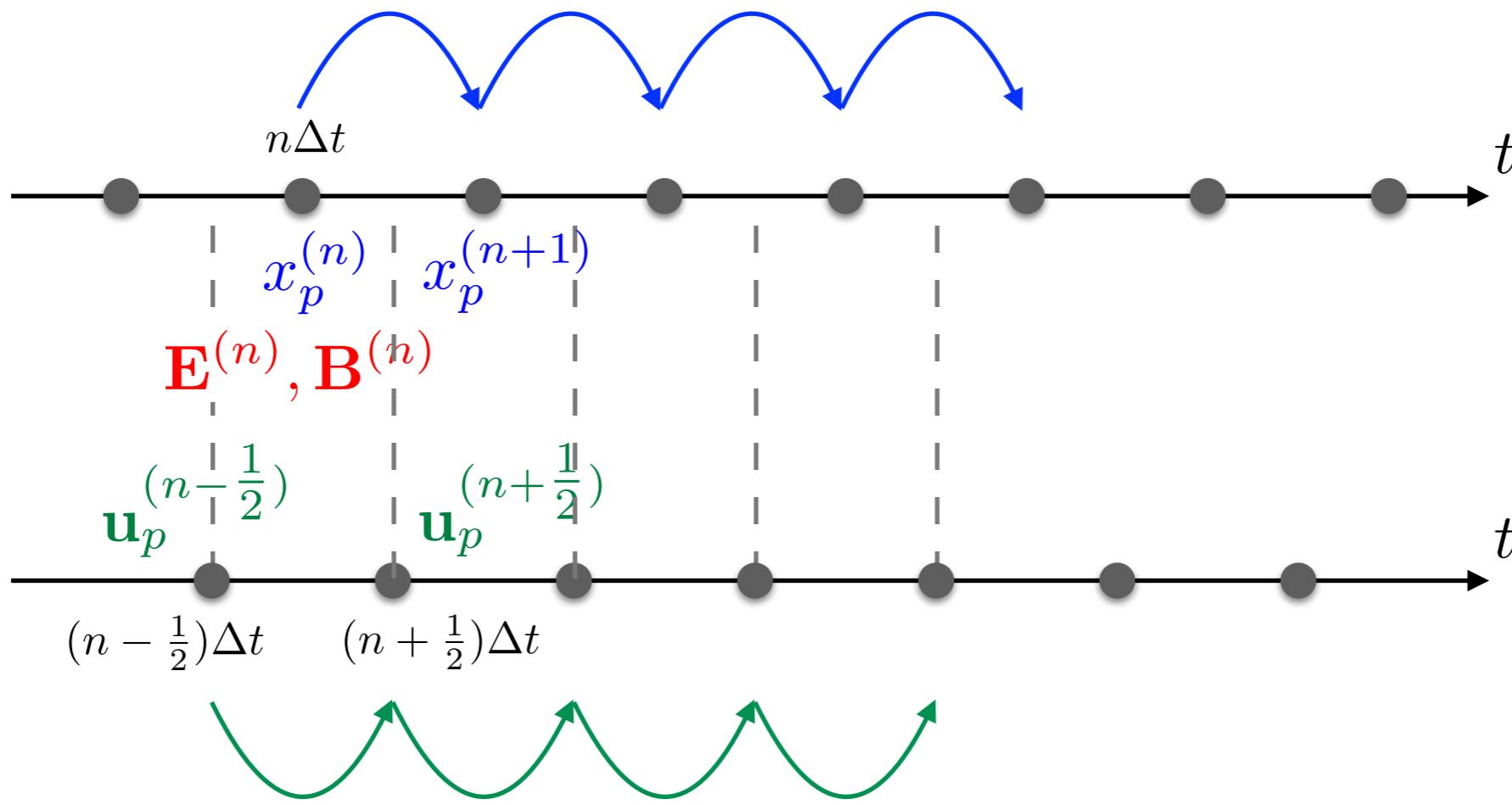
$$\mathbf{u}_m = \mathbf{u}_p^{(n-\frac{1}{2})} + \frac{q_s}{m_s} \frac{\Delta t}{2} \mathbf{E}_p$$

$$\mathbf{u}_p = \mathbf{u}_p^{(n-\frac{1}{2})} + \frac{q_s}{m_s} \Delta t \mathcal{M}(\mathbf{B}_p) \mathbf{u}_m$$



Step 2

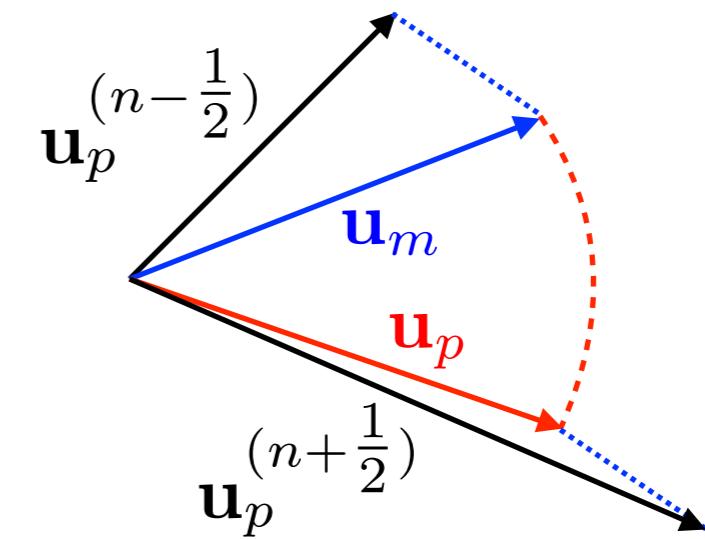
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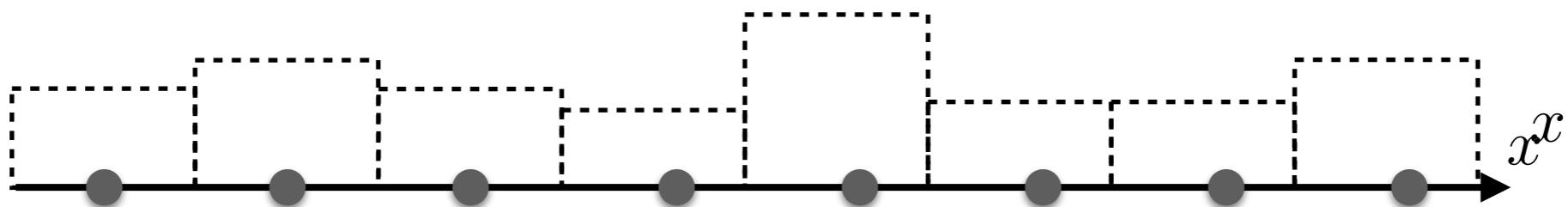
$$\mathbf{u}_p = \mathbf{u}_p^{(n - \frac{1}{2})} + \frac{q_s}{m_s} \Delta t \mathcal{M}(\mathbf{B}_p) \mathbf{u}_m$$

$$\mathbf{u}_p^{(n + \frac{1}{2})} = \mathbf{u}_p + \frac{q_s}{m_s} \frac{\Delta t}{2} \mathbf{E}_p$$



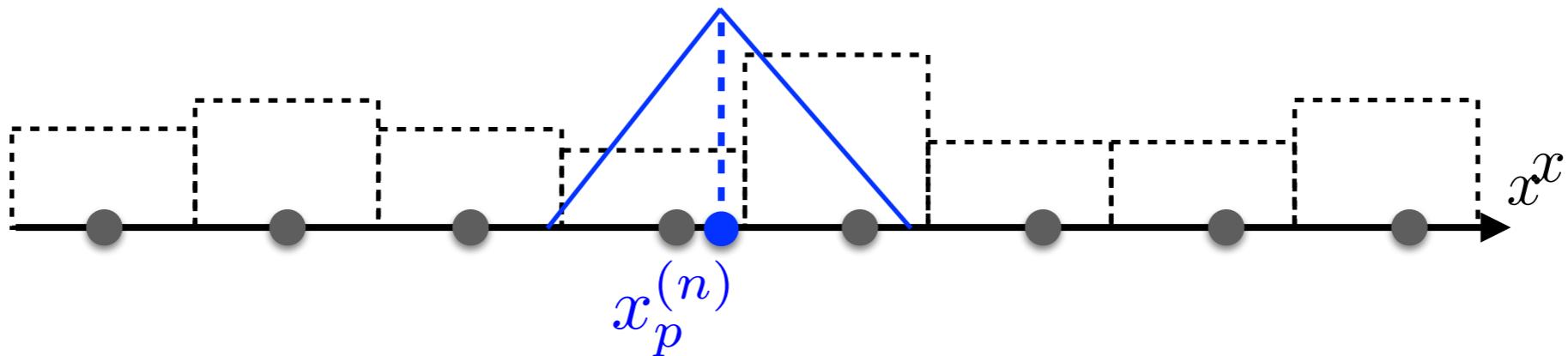
Step 3

Charge-conserving current deposition scheme are available among which Esirkepov's is ‘most’ popular



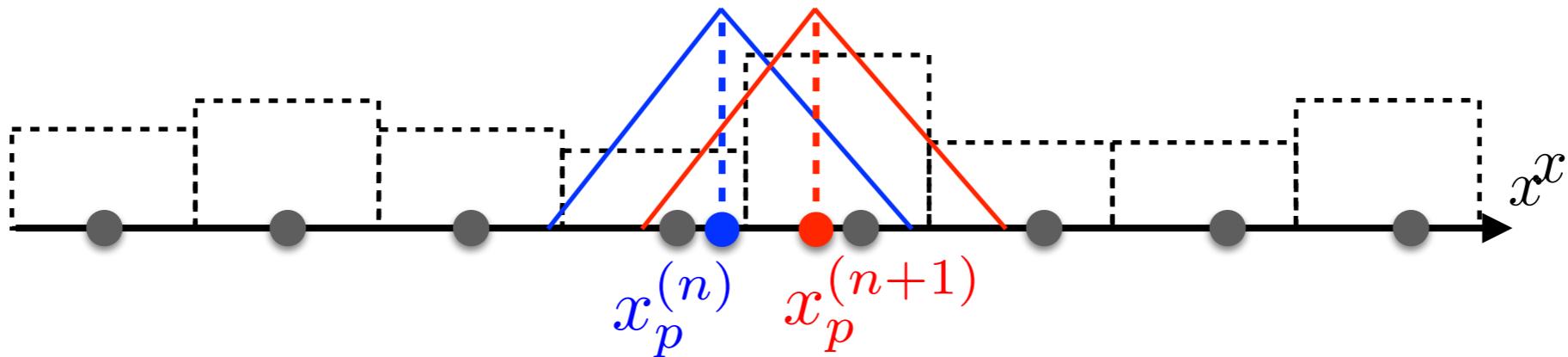
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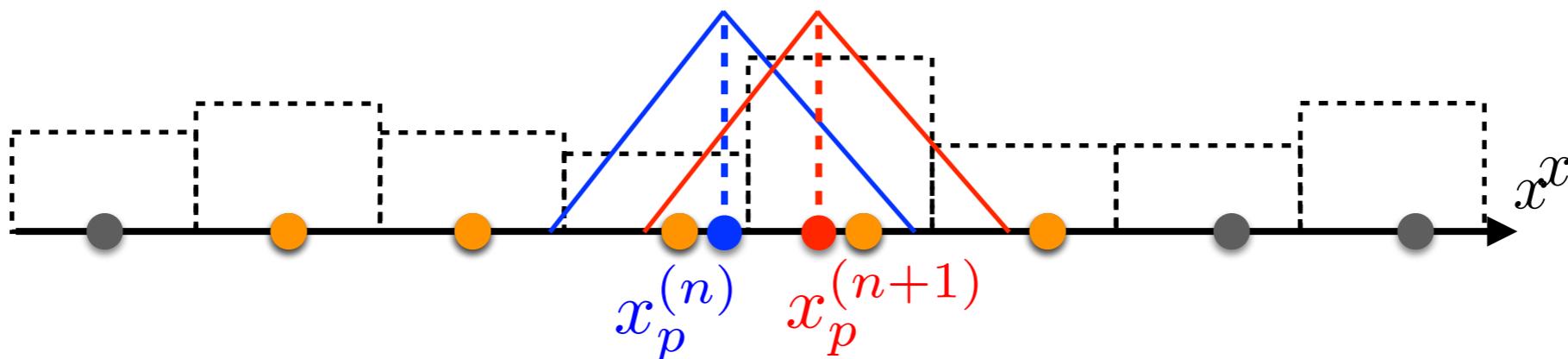
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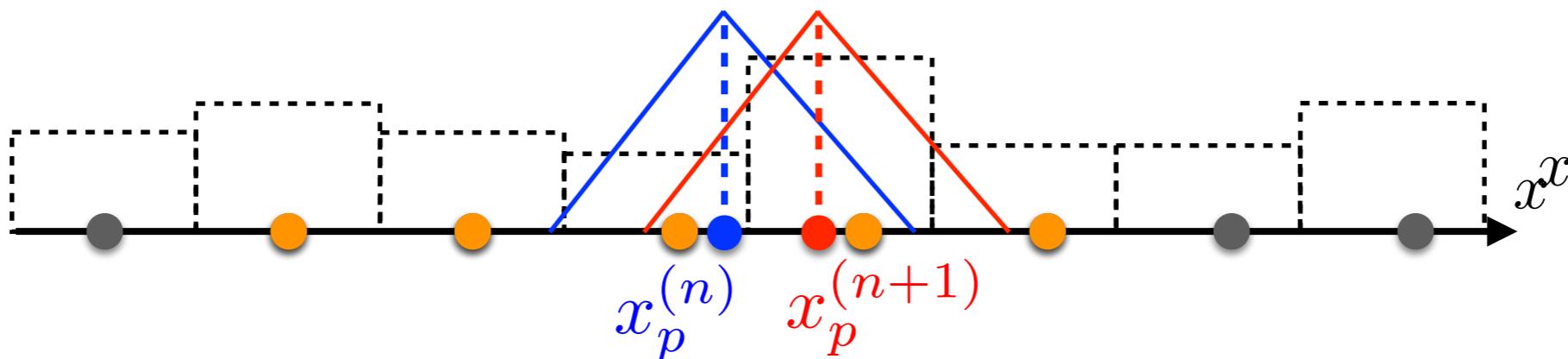
In 1D, current deposition is easily done directly from charge conservation:

$$\partial_x J_x = -\partial_t \rho$$

while other component are ‘directly’ projected onto the grid (see interpolation)

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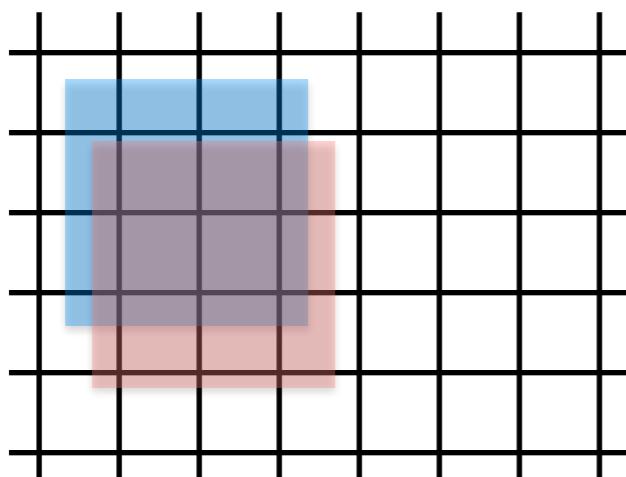


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In 2D & 3D, Esirkepov’s method allows to conserve charge (within machine precision)

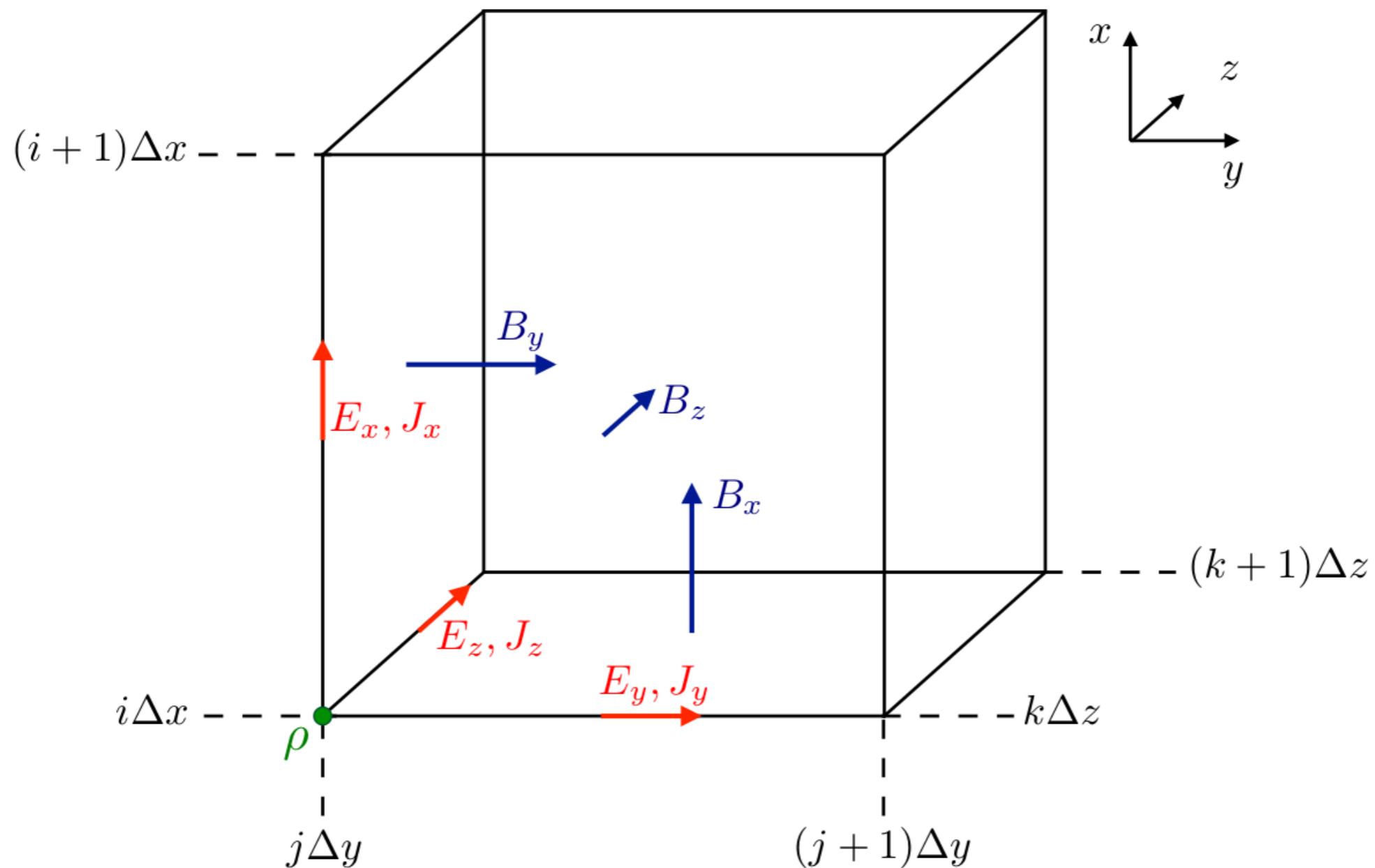


$$(J_{x,p})_{i+\frac{1}{2},j}^{(n+\frac{1}{2})} = (J_{x,p})_{i-\frac{1}{2},j}^{(n+\frac{1}{2})} + q_s w_p \frac{\Delta x}{\Delta t} (W_x)_{i+\frac{1}{2},j}^{(n+\frac{1}{2})}$$

$$(J_{y,p})_{i,j+\frac{1}{2}}^{(n+\frac{1}{2})} = (J_{y,p})_{i,j-\frac{1}{2}}^{(n+\frac{1}{2})} + q_s w_p \frac{\Delta y}{\Delta t} (W_y)_{j,i+\frac{1}{2}}^{(n+\frac{1}{2})}$$

Step 4

The Finite-Difference Time-Domain (FDTD) method is a popular method for solving Maxwell's Equations



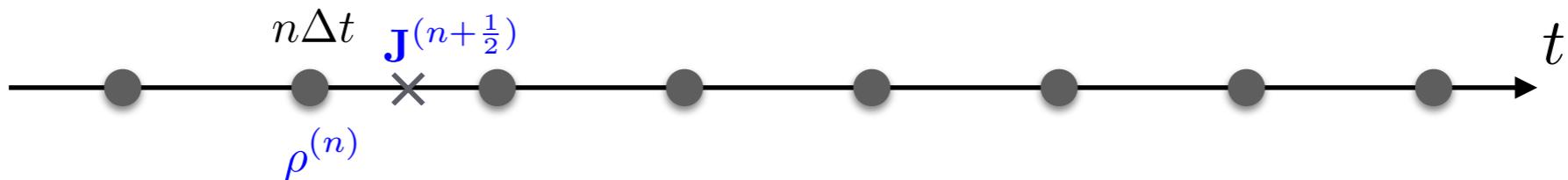
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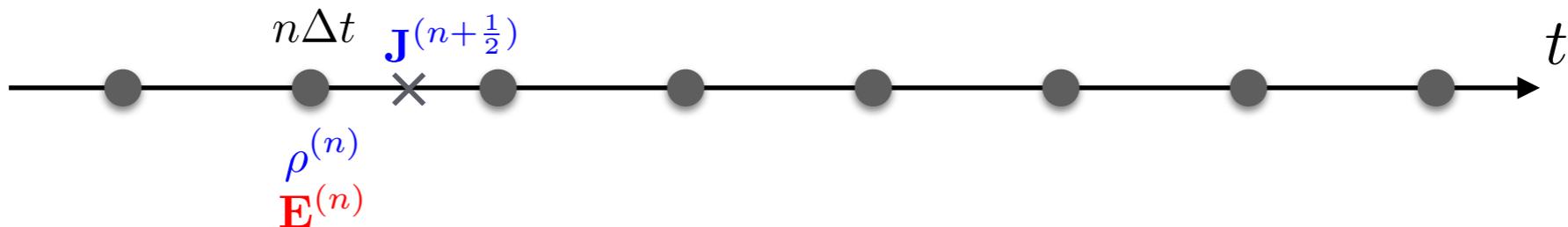
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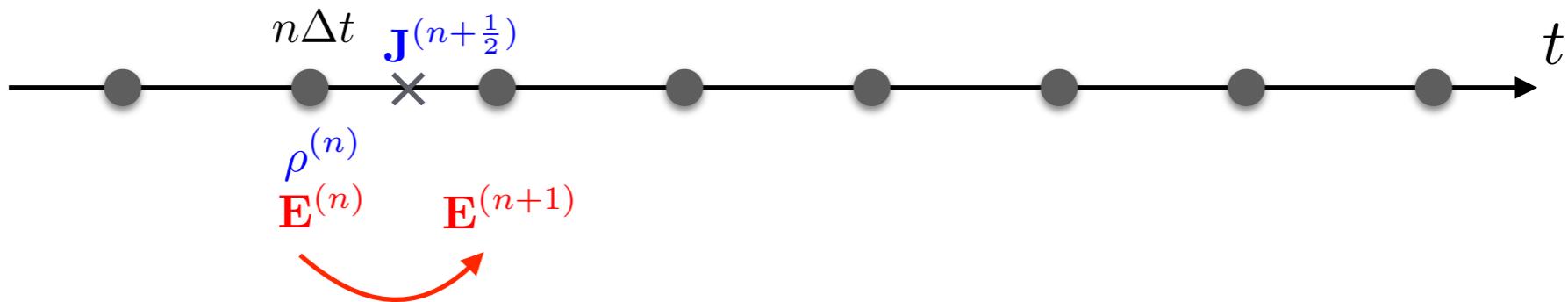
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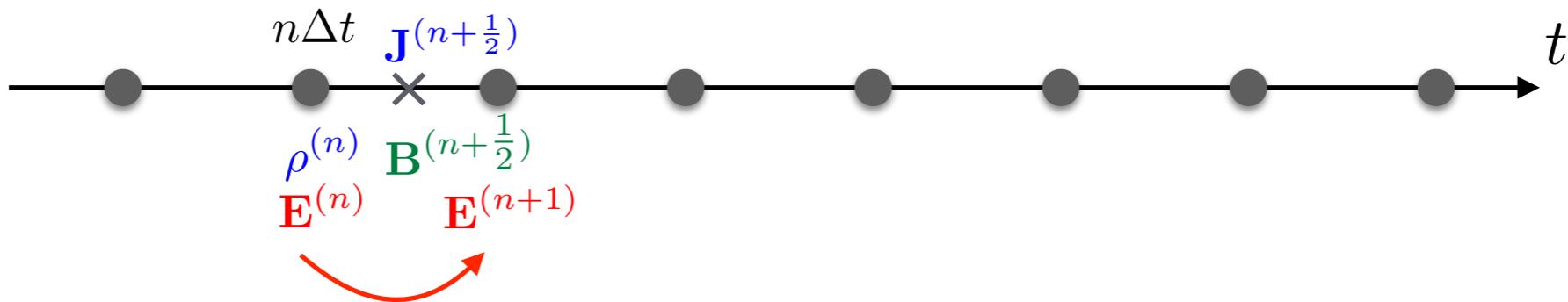
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Solving Ampère's equation: $\partial_t E_y = -J_y - \partial_x B_z$

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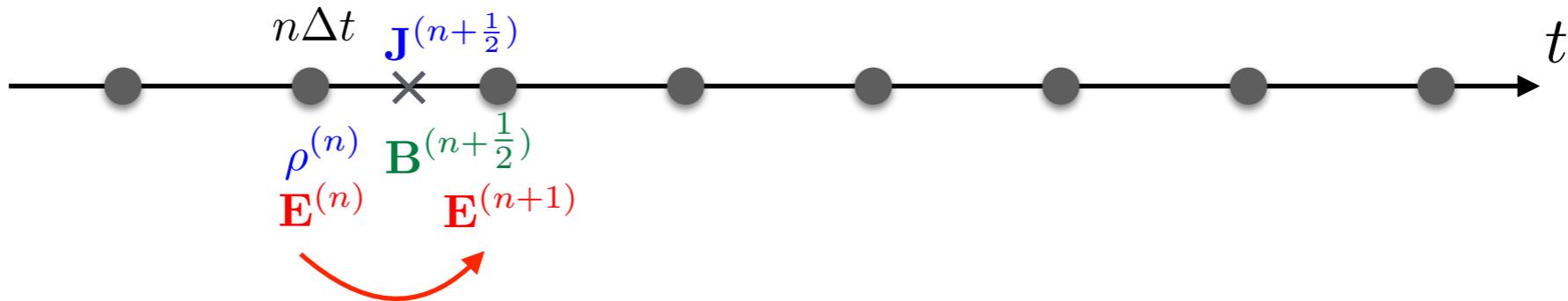


Solving Ampère's equation: $\partial_t E_y = -J_y - \partial_x B_z$

time-centering
$$\frac{(E_y)^{(n+1)} - (E_y)^{(n)}}{\Delta t} = -J_y^{(n+\frac{1}{2})} - (\partial_x B_z)^{(n+\frac{1}{2})}$$

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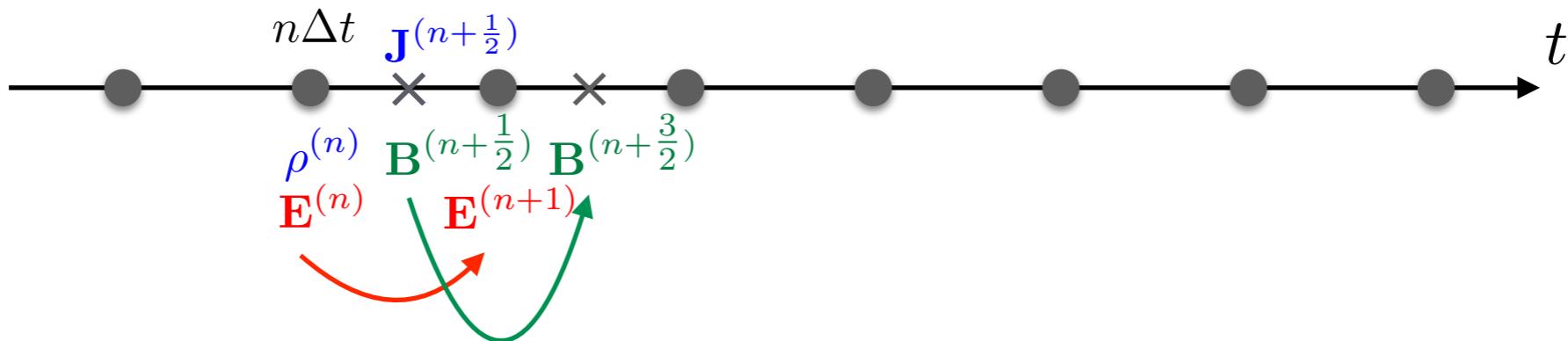
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space-centering
$$\frac{(E_y)_i^{(n+1)} - (E_y)_i^{(n)}}{\Delta t} = -(J_y)_i^{(n+\frac{1}{2})} - \frac{(B_z)_{i+\frac{1}{2}}^{(n+\frac{1}{2})} - (B_z)_{i-\frac{1}{2}}^{(n-\frac{1}{2})}}{\Delta x}$$

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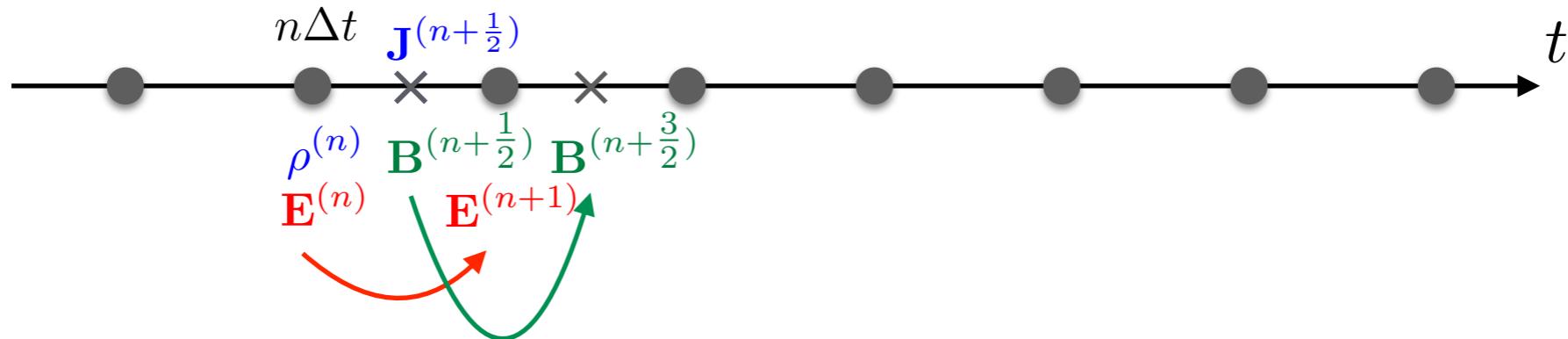
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Solving Faraday's equation: $\partial_t B_z = \partial_x E_y$

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Solving Ampère's equation: $\partial_t E_y = -J_y - \partial_x B_z$

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$$\text{space/time-centering} \quad \frac{(B_z)_{i+\frac{1}{2}}^{(n+\frac{3}{2})} - (B_z)_{i+\frac{1}{2}}^{(n+\frac{1}{2})}}{\Delta t} = \frac{(E_y)_{i+1}^{(n+1)} - (E_y)_i^{(n+1)}}{\Delta x}$$

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Numerical analysis of the FDTD solvers gives you access to the numerical dispersion relation & CFL condition

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The *numerical* electromagnetic wave equation in a vacuum

$$\partial_t^N \mathbf{E} = +\nabla^N \times \mathbf{B}$$

$$\partial_t^N \mathbf{B} = -\nabla^N \times \mathbf{E}$$

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with:

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Using the standard technique to derive the wave equation leads to:

$$\partial_{tt}^N \mathbf{E} + \sum_\mu \partial_{\mu\mu}^N \mathbf{E} = 0$$

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Looking for *numerical* solution in the form:

$$(E_y)_{i,j+\frac{1}{2},k}^{(n)} = E_{y0} \exp \left\{ i \left[ik_x \Delta x + (j + \frac{1}{2}) k_y \Delta y + k_z \Delta z - n \omega \Delta t \right] \right\}$$

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After some algebra, one finds the *numerical dispersion relation*:

$$\frac{\sin^2(\omega\Delta t/2)}{\Delta t^2} = \sum_{\mu} \frac{\sin^2(k_{\mu}\Delta\mu/2)}{\Delta\mu^2}$$

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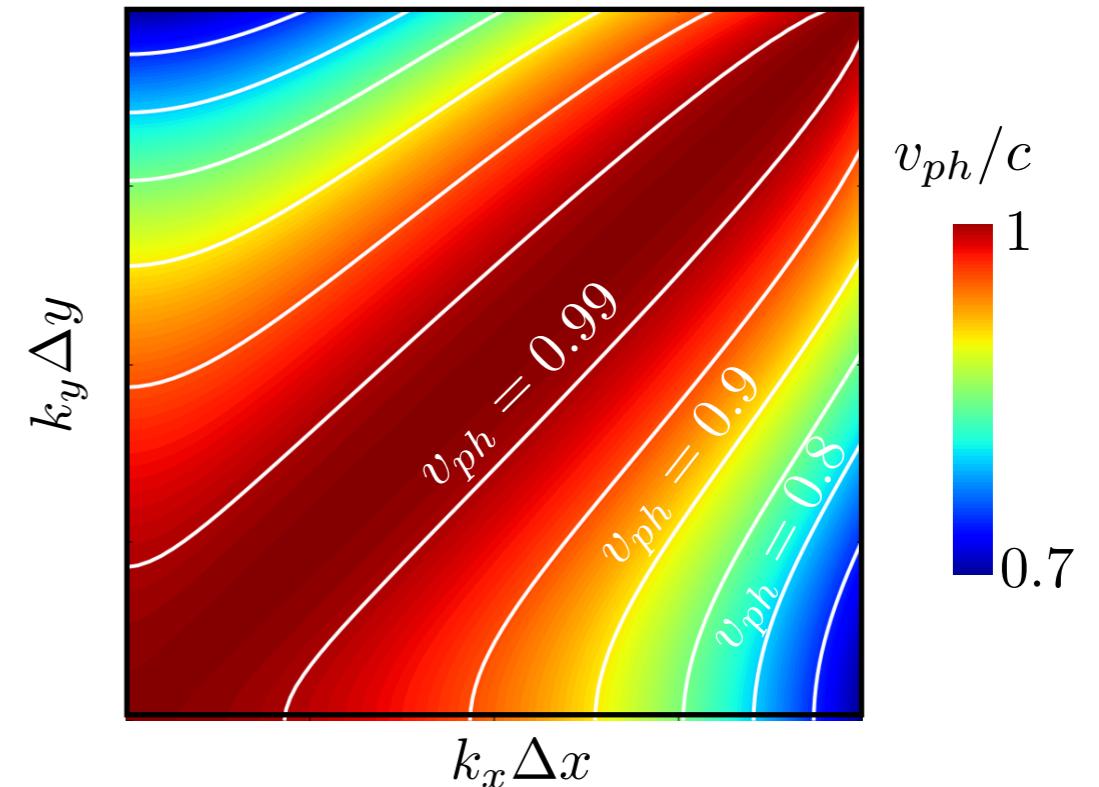
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The FDTD solver is subject to *numerical dispersion* as the numerical light wave velocity is found to depend on its wavenumber and orientation.



A quick summary

The PIC approach in a nutshell

Initialization	time step $n = 0$, time $t = 0$
Particle loading	$\forall p$, define $(\mathbf{x}_p)^{n=0}$, $(\mathbf{u}_p)^{n=-\frac{1}{2}}$
Charge projection on grid	$[\forall p, (\mathbf{x}_p)^{n=0}] \rightarrow \rho^{(n=0)}(\mathbf{x})$
Compute initial fields	<ul style="list-style-type: none"> - solve Poisson on grid: $[\rho^{(n=0)}(\mathbf{x})] \rightarrow \mathbf{E}_{\text{stat}}^{(n=0)}(\mathbf{x})$ - add external fields: $\mathbf{E}^{(n=0)}(\mathbf{x}) = \mathbf{E}_{\text{stat}}^{(n=0)}(\mathbf{x}) + \mathbf{E}_{\text{ext}}^{(n=0)}(\mathbf{x})$ $\mathbf{B}^{(n=\frac{1}{2})}(\mathbf{x}) = \mathbf{B}_{\text{ext}}^{(n=\frac{1}{2})}(\mathbf{x})$
PIC loop: from time step n to $n + 1$, time $t = (n + 1) \Delta t$	

Restart charge & current densities

Save magnetic fields value (used to center magnetic fields)

Interpolate fields at particle positions $\forall p, [\mathbf{x}_p, \mathbf{E}^{(n)}(\mathbf{x}), \mathbf{B}^{(n)}(\mathbf{x})] \rightarrow \mathbf{E}_p^{(n)}, \mathbf{B}_p^{(n)}$

Push particles

- compute new velocity $\forall p, \mathbf{p}_p^{(n-\frac{1}{2})} \left[\mathbf{E}_p^{(n)}, \mathbf{B}_p^{(n)} \right] \mathbf{p}_p^{(n+\frac{1}{2})}$
- compute new position $\forall p, \mathbf{x}_p^{(n)} \left[\mathbf{p}_p^{(n+\frac{1}{2})} \right] \mathbf{x}_p^{(n+1)}$

Project current onto the grid using a charge-conserving scheme

$$\left[\forall p \quad \mathbf{x}_p^{(n)}, \mathbf{x}_p^{(n+1)}, \mathbf{p}_p^{(n+\frac{1}{2})} \right] \rightarrow \mathbf{J}^{(n+\frac{1}{2})}(\mathbf{x})$$

Solve Maxwell's equations

- solve Maxwell-Faraday: $\mathbf{E}^{(n)}(\mathbf{x}) \left[\mathbf{J}^{(n+\frac{1}{2})(\mathbf{x})} \right] \mathbf{E}^{(n+1)}(\mathbf{x})$
- solve Maxwell-Ampère: $\mathbf{B}^{(n+\frac{1}{2})}(\mathbf{x}) \left[\mathbf{E}^{(n+1)}(\mathbf{x}) \right] \mathbf{B}^{(n+\frac{3}{2})}(\mathbf{x})$
- center magnetic fields: $\mathbf{B}^{(n+1)}(\mathbf{x}) = \frac{1}{2} \left(\mathbf{B}^{(n+\frac{1}{2})}(\mathbf{x}) + \mathbf{B}^{(n+\frac{3}{2})}(\mathbf{x}) \right)$

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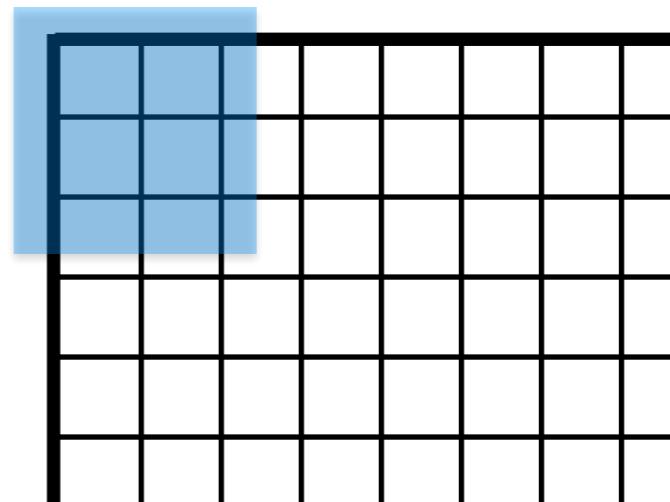
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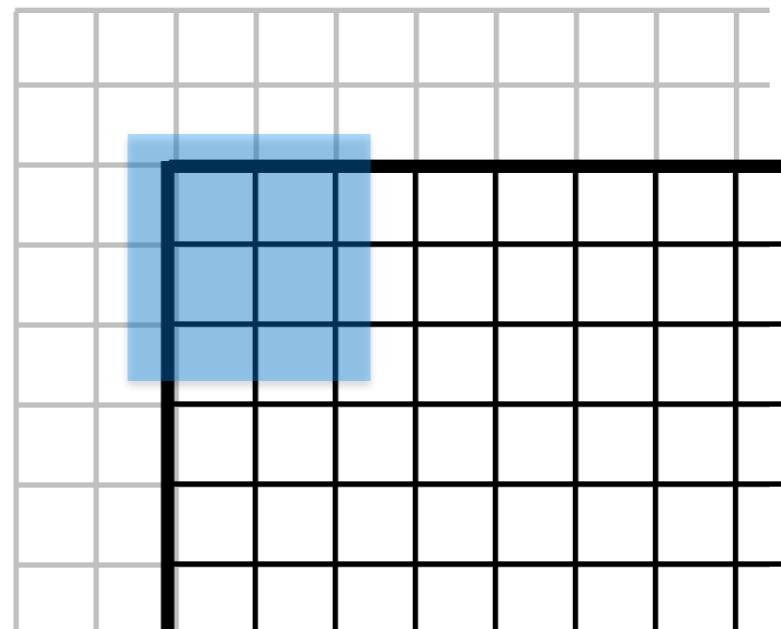
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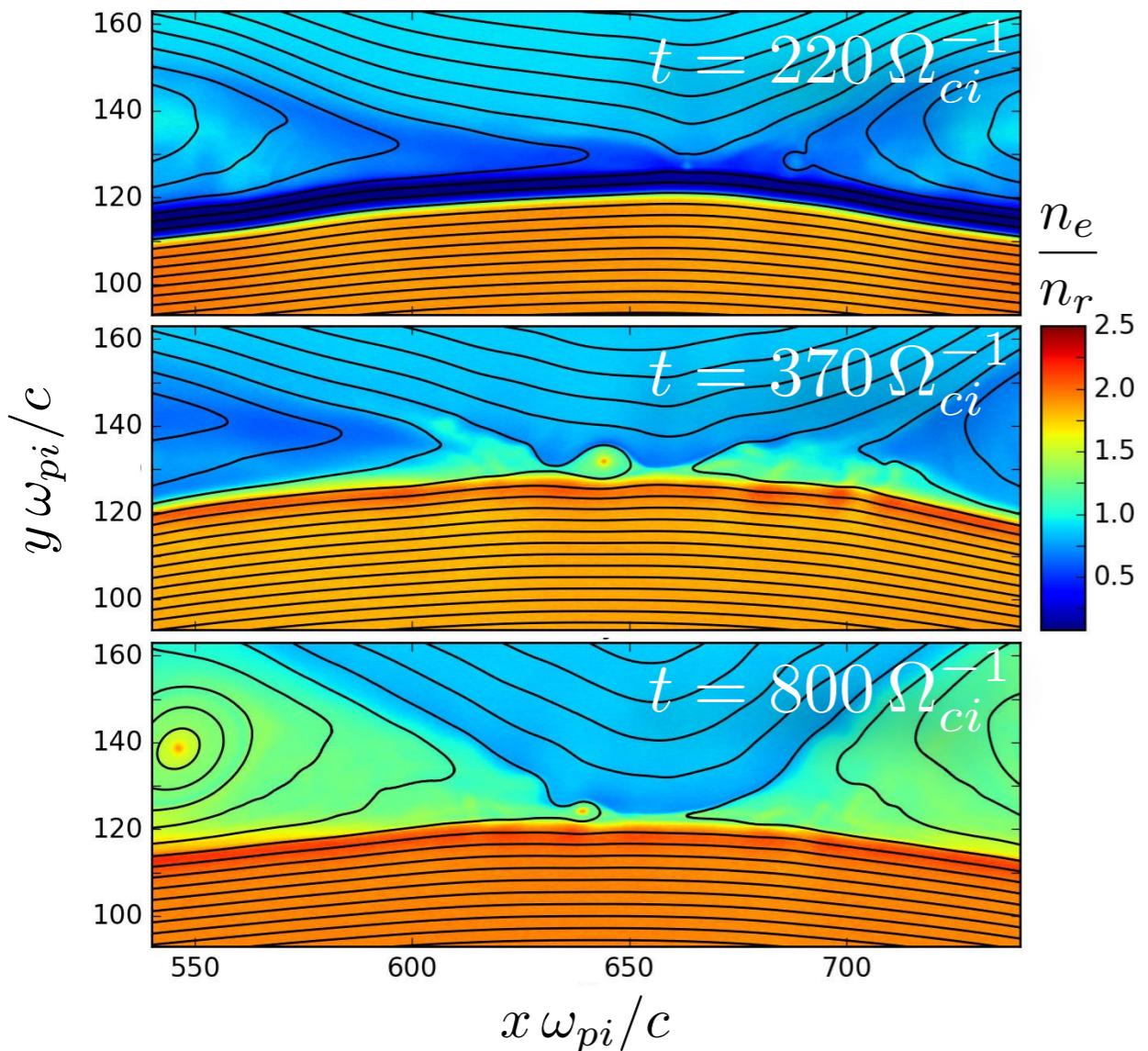
Outlines

- Numerical approach: **how to build your PIC code**
- **High-performance computing:
getting ready for the super-computers**
- Additional modules: **beyond the *collisionless* plasma**
- Some physics highlights: **what you can do with a PIC code**

Parallelization is mandatory for large-scale PIC simulation

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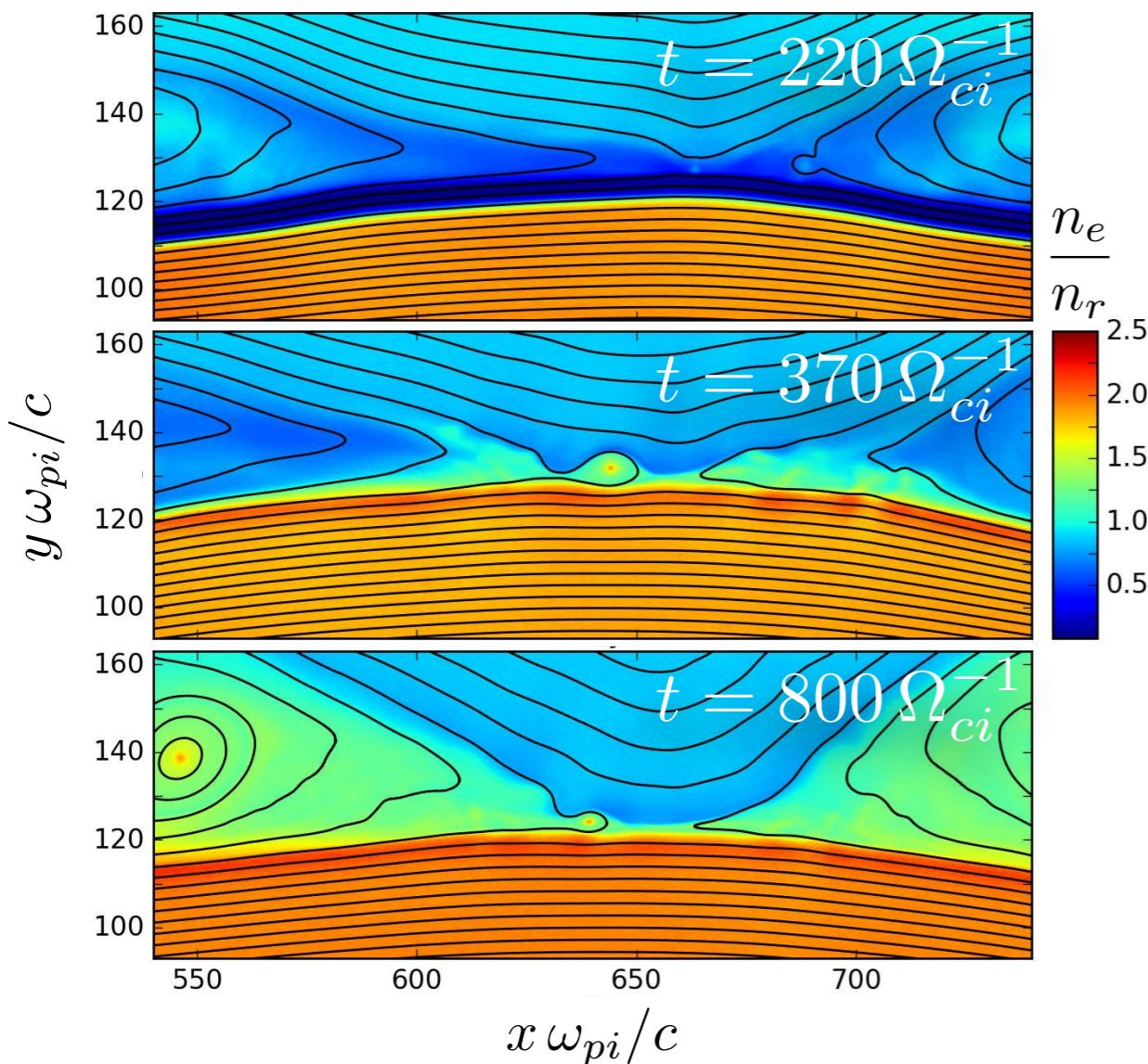
Large scale PIC simulation of magnetic reconnection at the earth magnetopause



Simulation box: $1280 \frac{c}{\omega_{pi}} \times 256 \frac{c}{\omega_{pi}}$
 25600×10240 PIC cells
run up to $t = 800 \Omega_{ci}^{-1}$
 $N_t \sim 9.5 \times 10^5$ timesteps
for a total of 22×10^9 quasi-particles.

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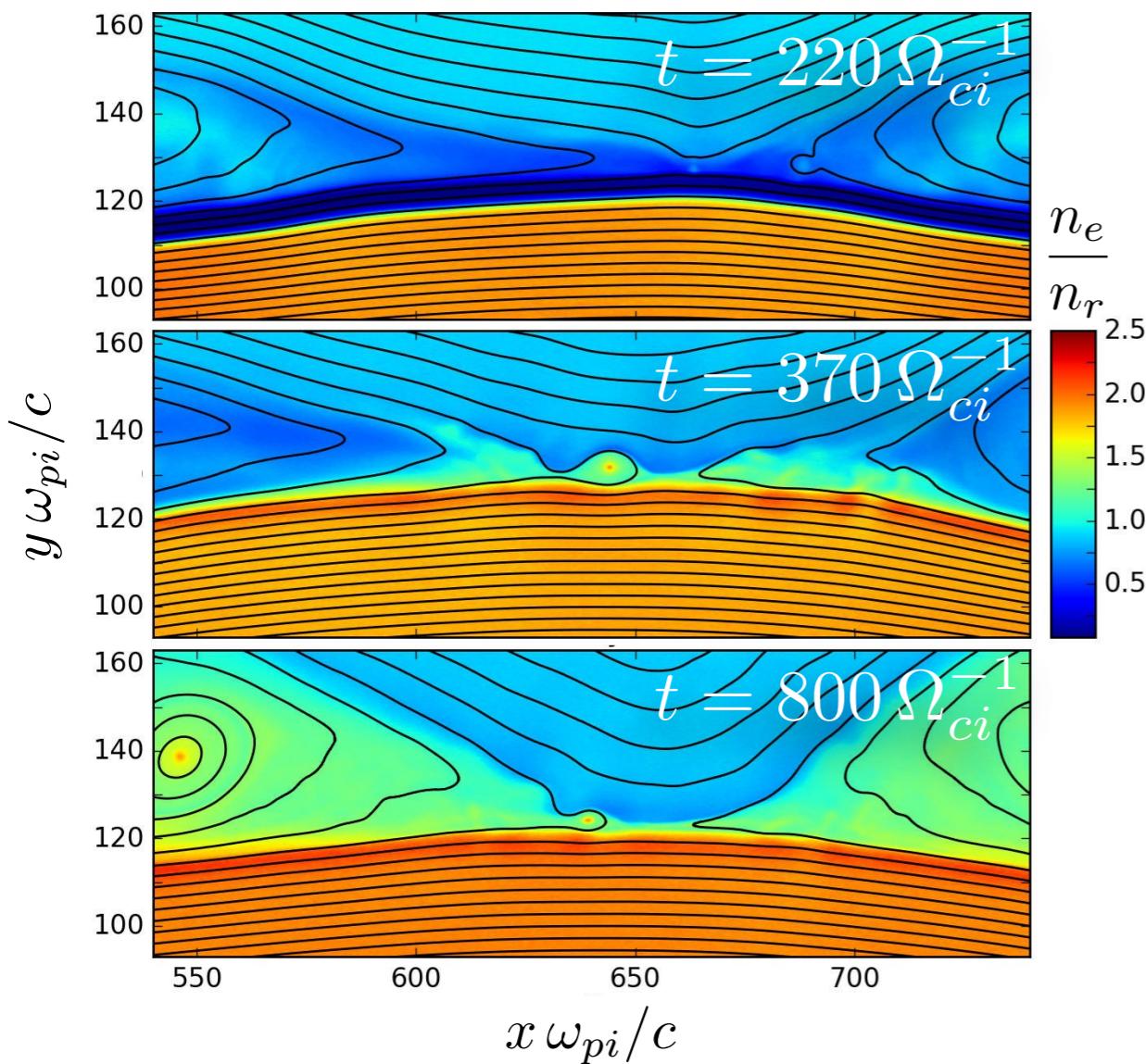


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14 000 000 hours \sim 1600 years!!!

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Solution:
share the work on 16384 CPUs !!!

High-performance computing new paradigms & challenges

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Tianhe-2 34 PF:
17 MW

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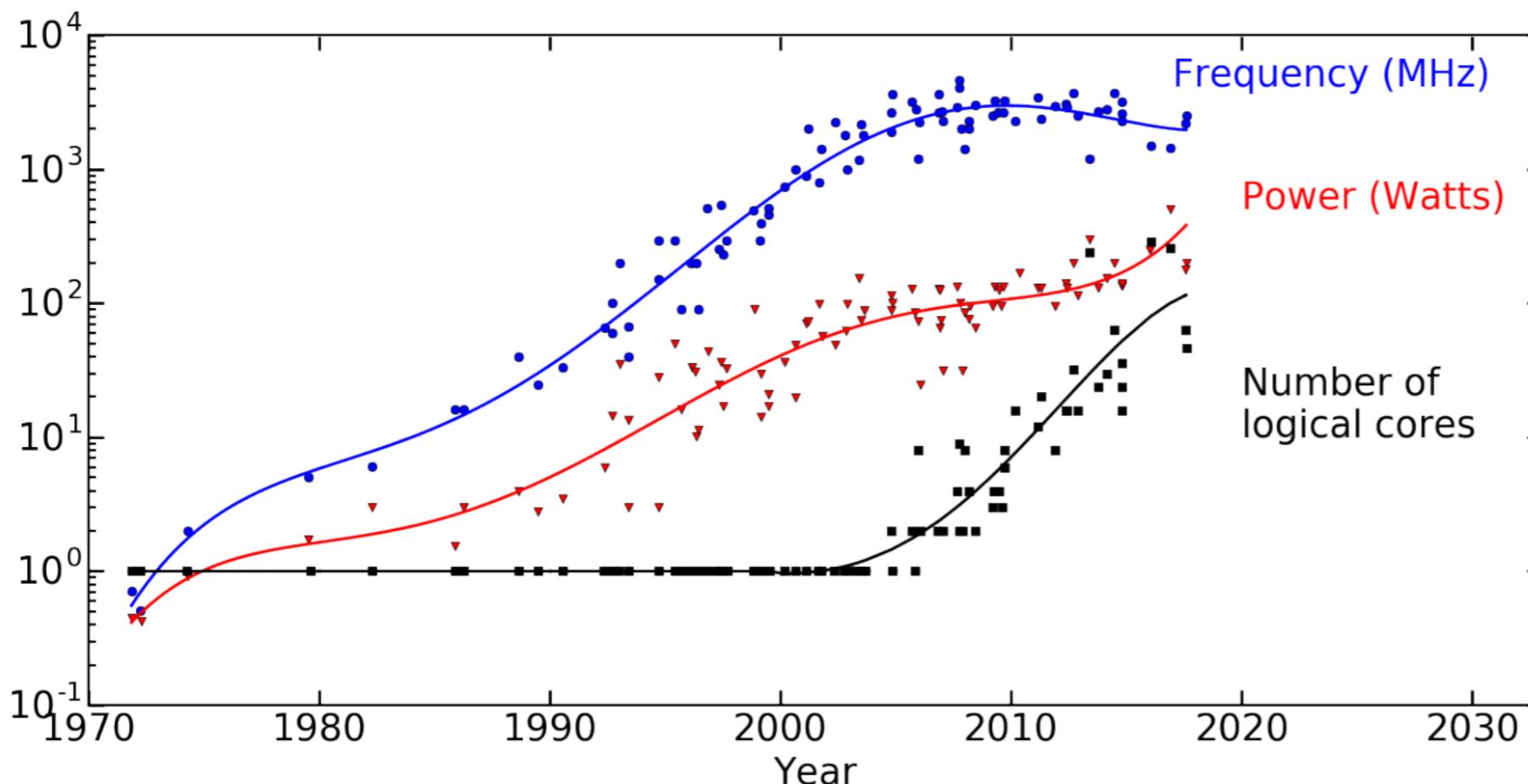
Exascale 1000 PF:

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High-performance computing new paradigms & challenges

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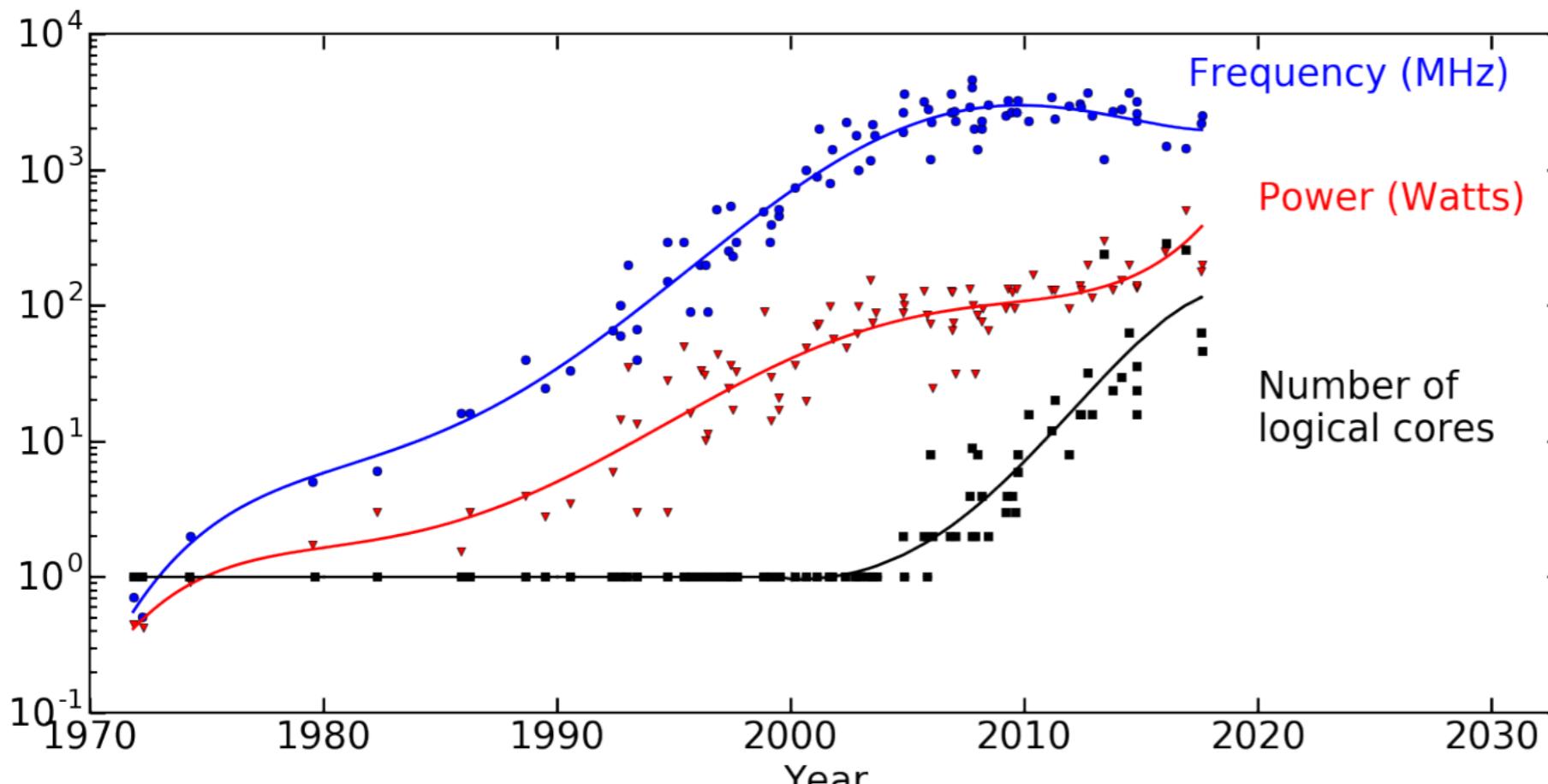


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Parallelism*
massive
hybrid MPI-OpenMP
dynamic (load balance)

Memory
shared vs. distributed
cache use

Vectorization**
SIMD

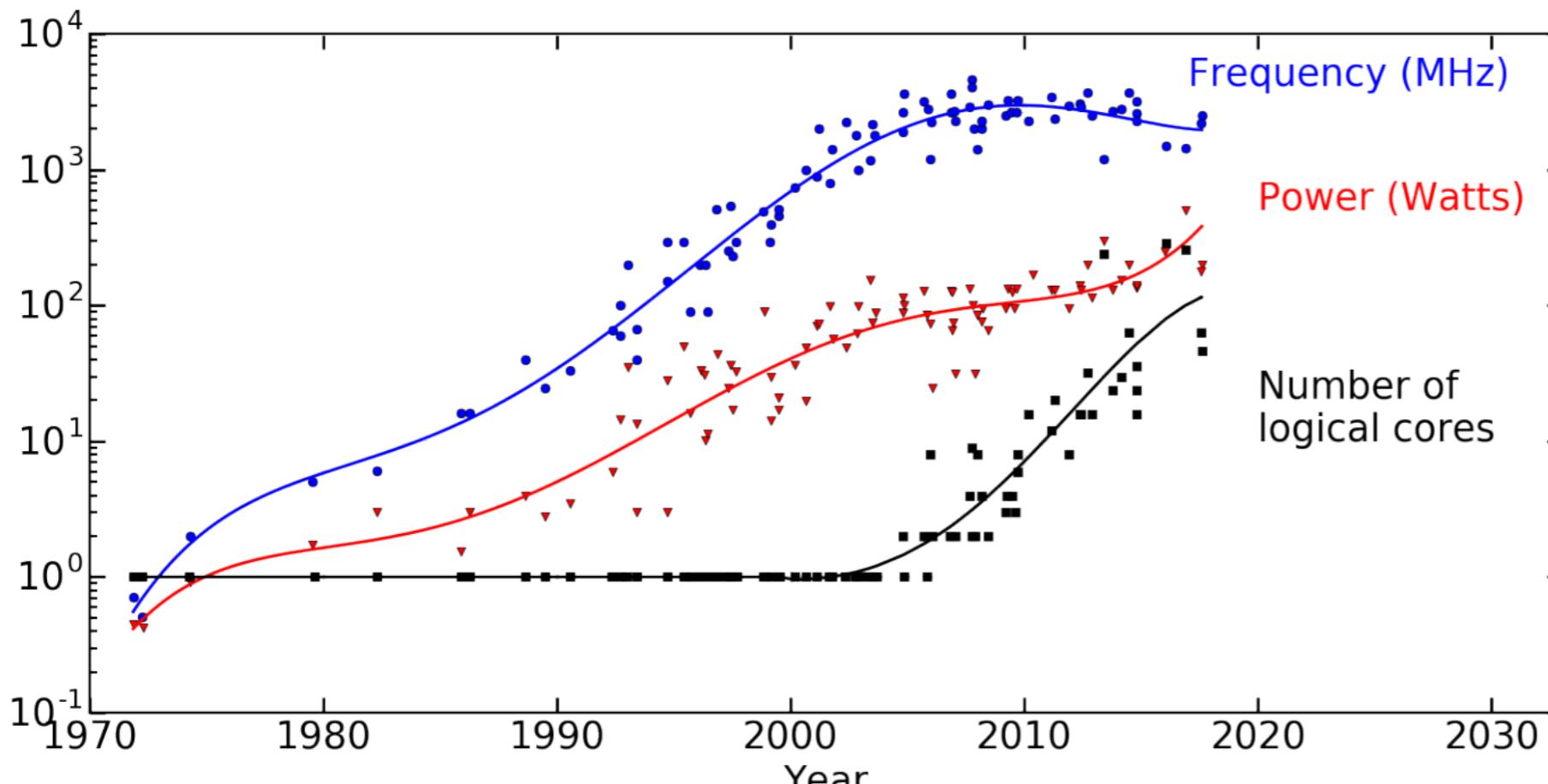
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*Derouillat *et al.*, Comp. Phys. Comm. **222**, 351 (2018)

**Beck *et al.*, arXiv:1810.03949

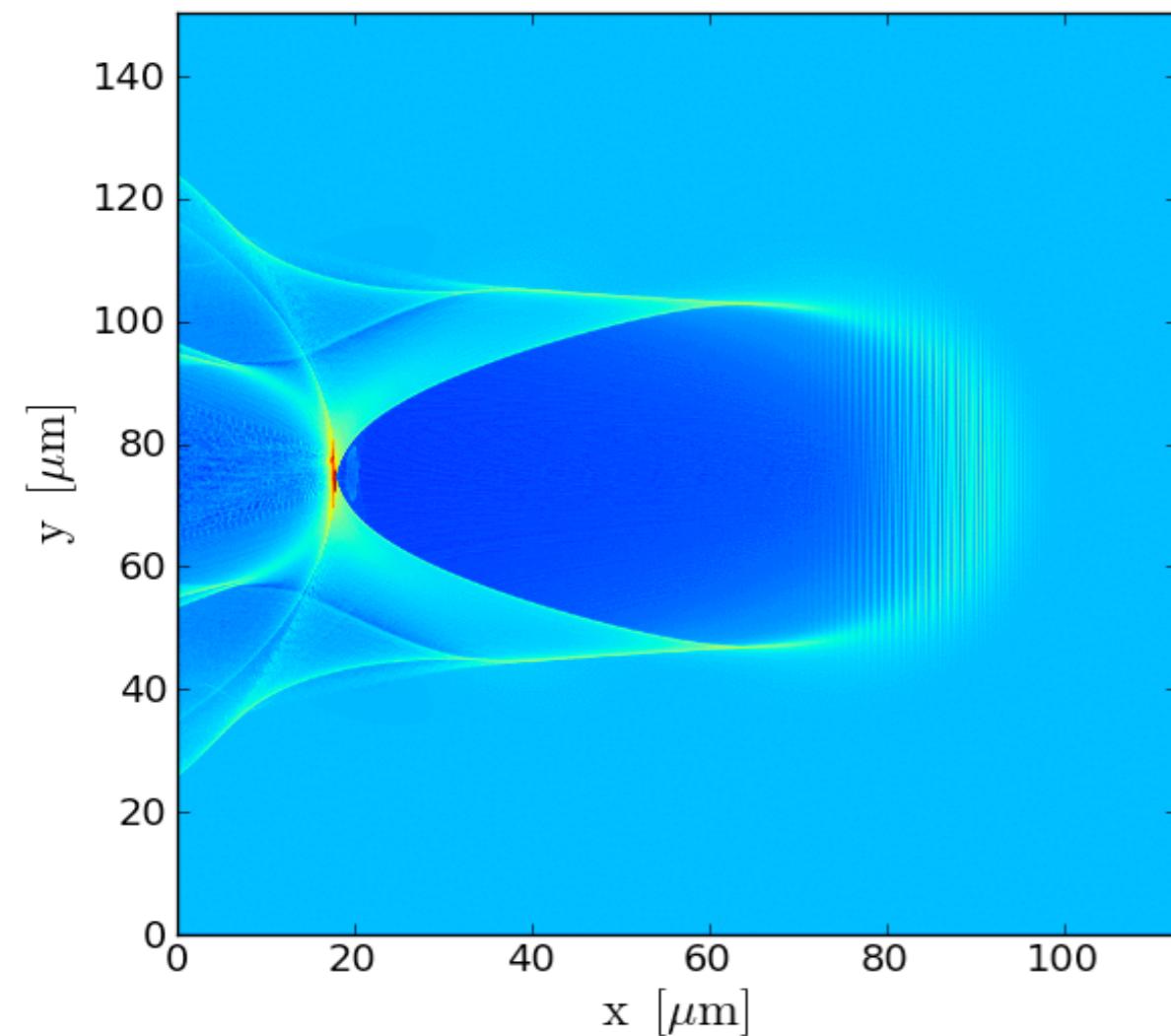
Step 1: Parallelization

PIC codes are well adapted to massive parallelism

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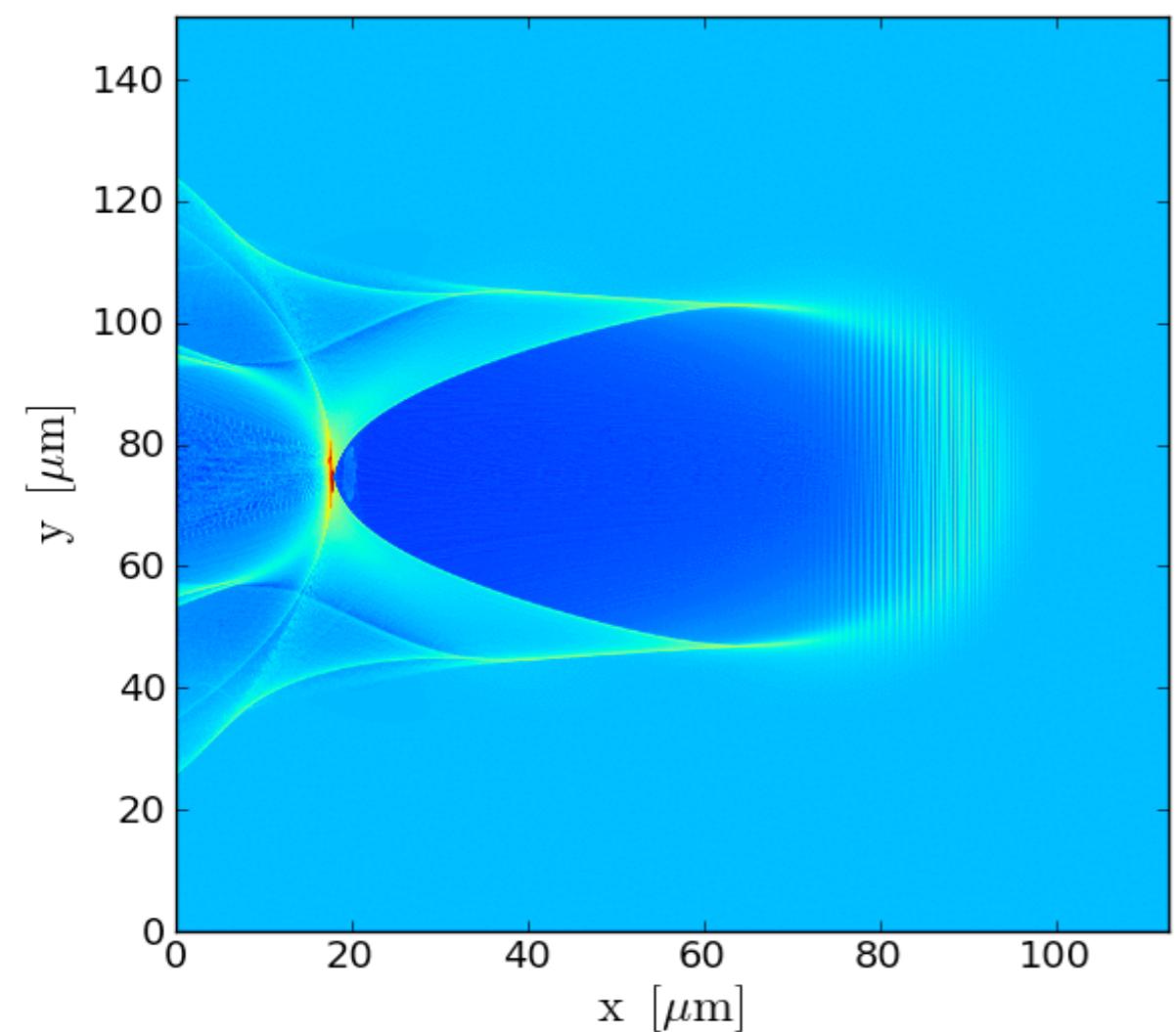
My Simulation (LWFA)



Step 1: Parallelization

PIC codes are well adapted to massive parallelism

My Simulation (LWFA)



My Super-Computer

CE
computing element

CE-0

CE-1

CE-2

CE-3

CE-4

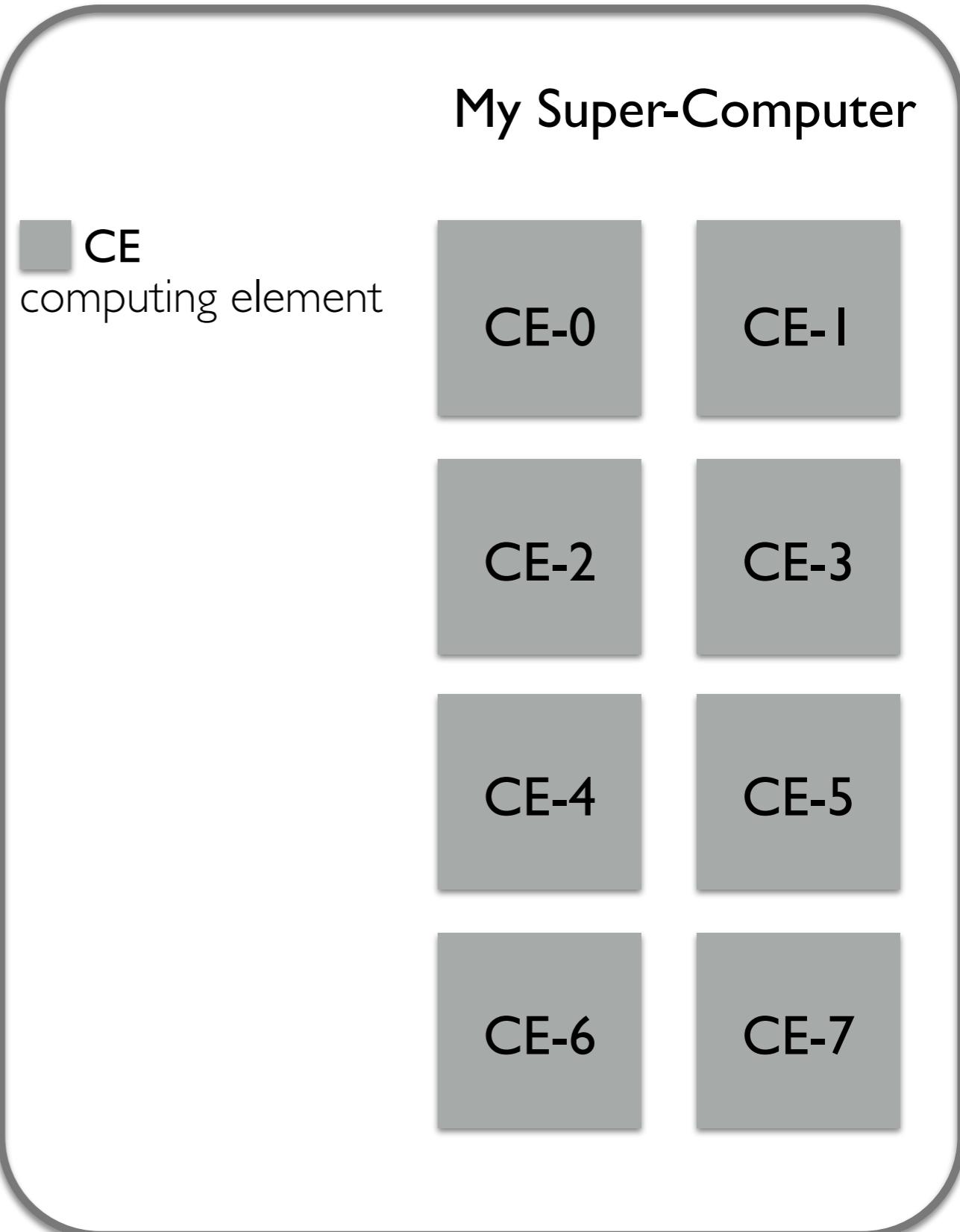
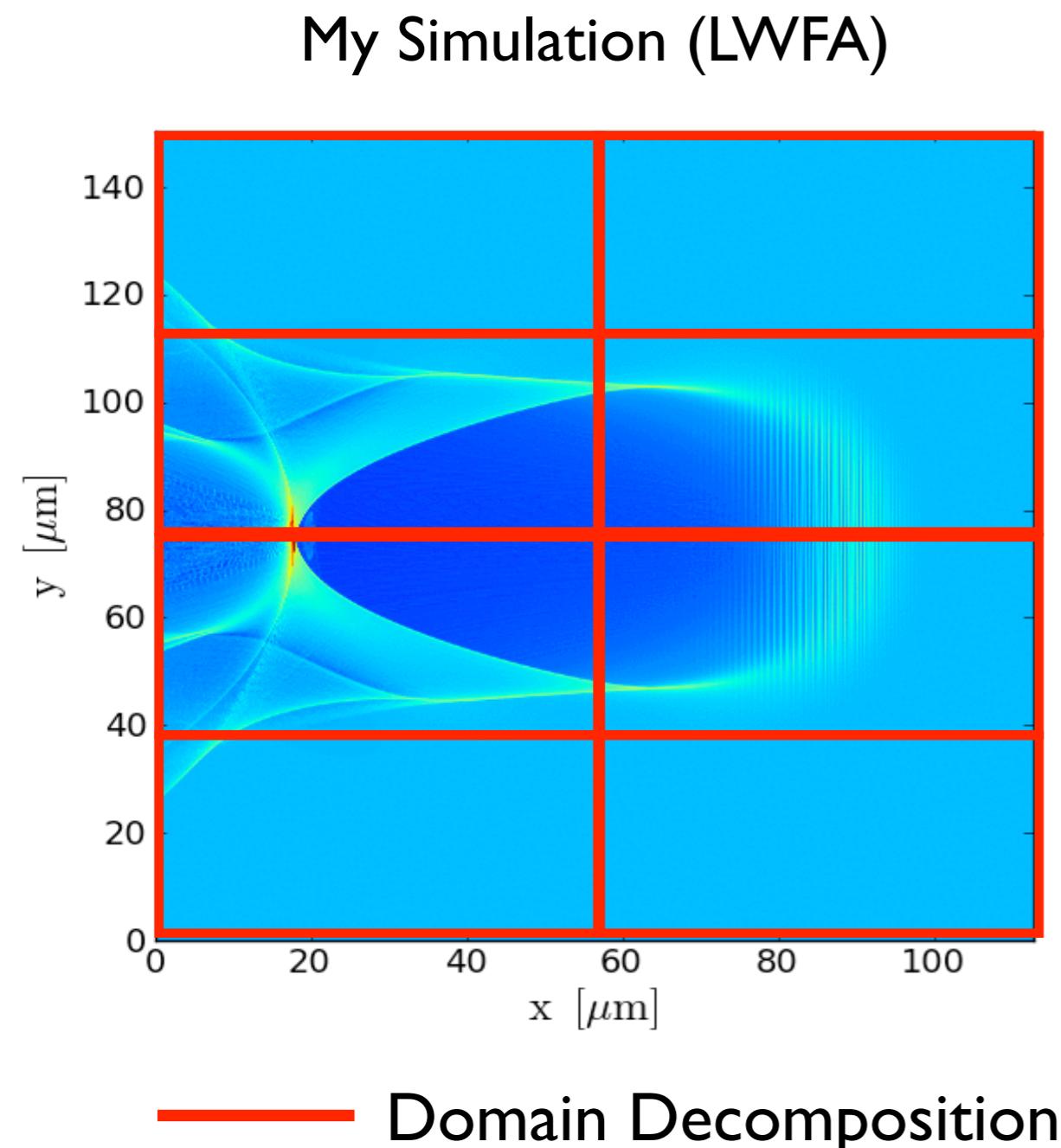
CE-5

CE-6

CE-7

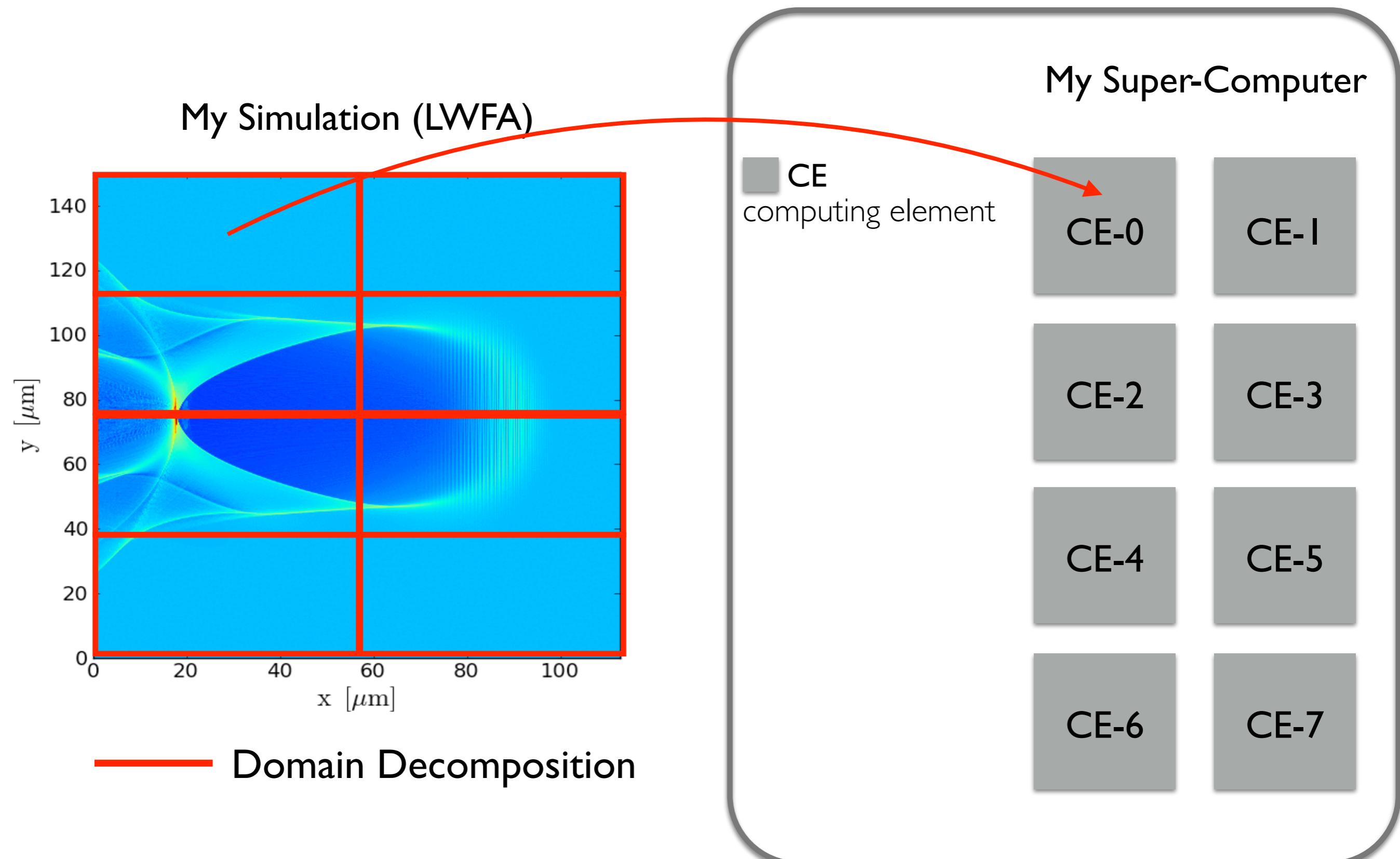
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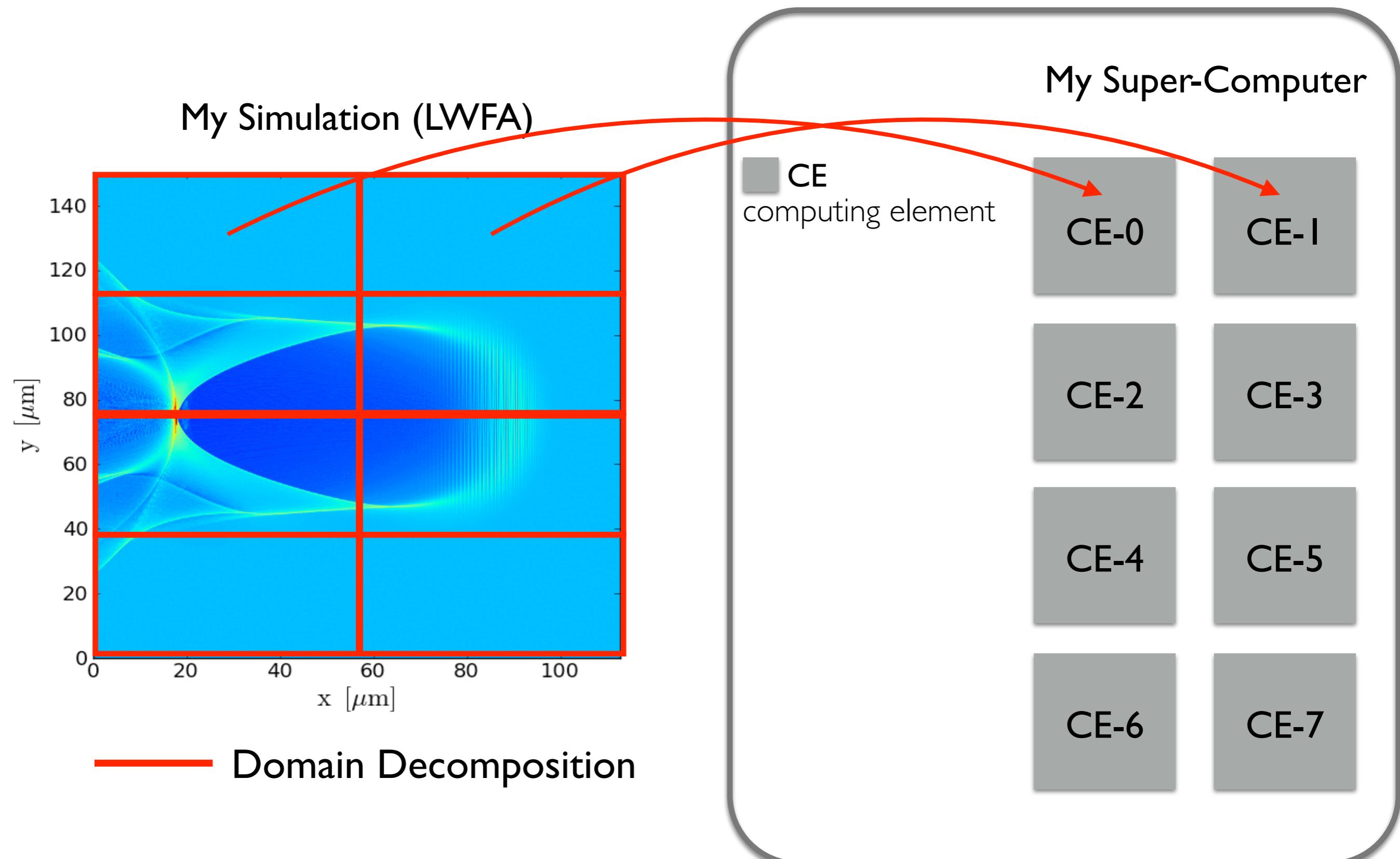
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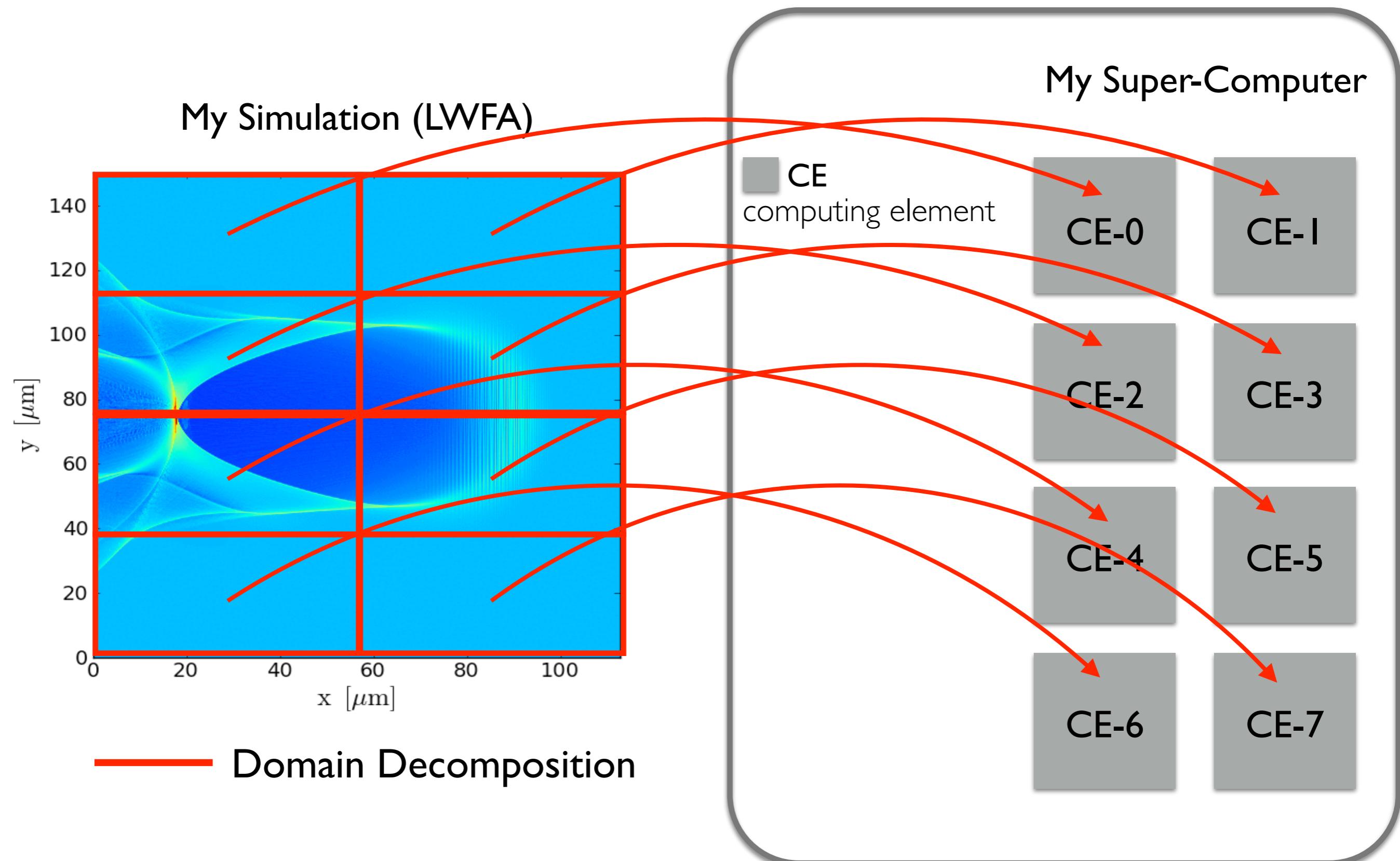
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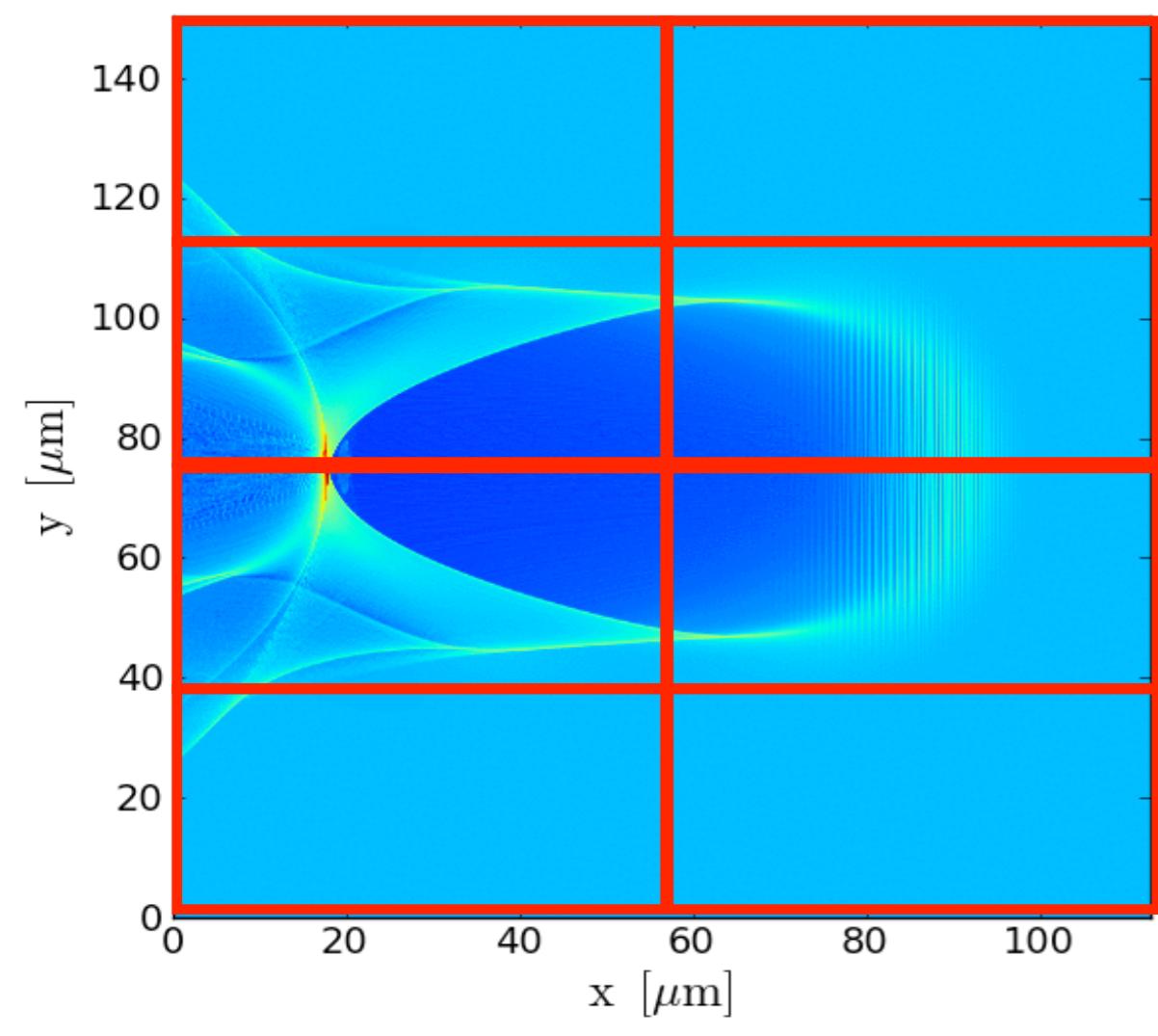
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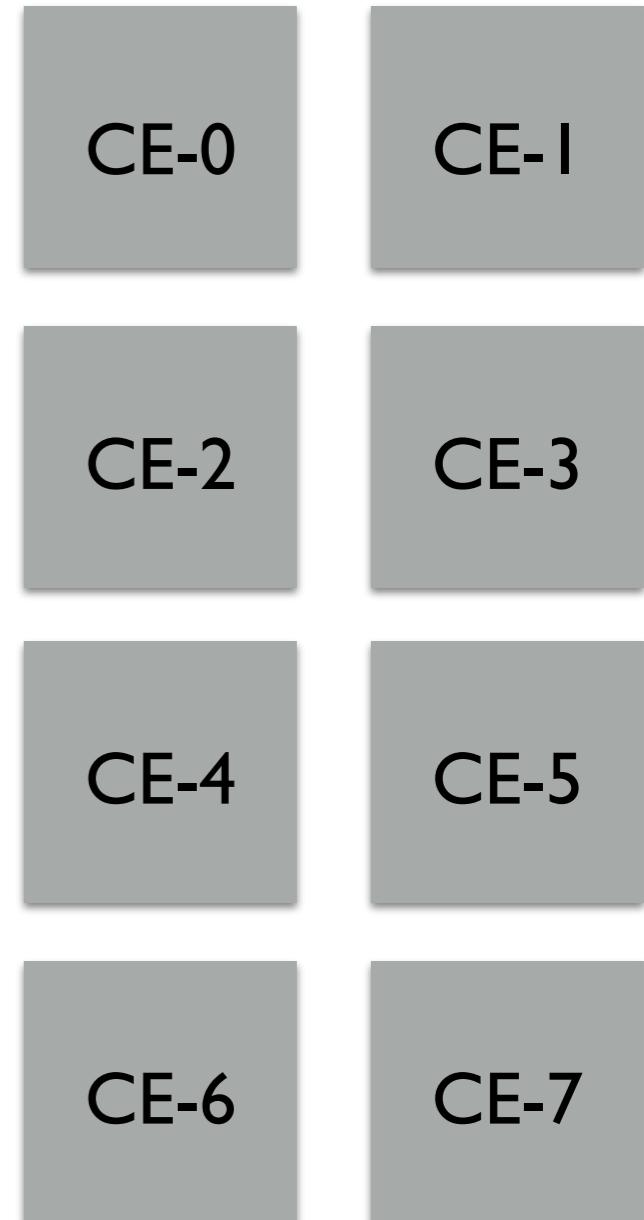
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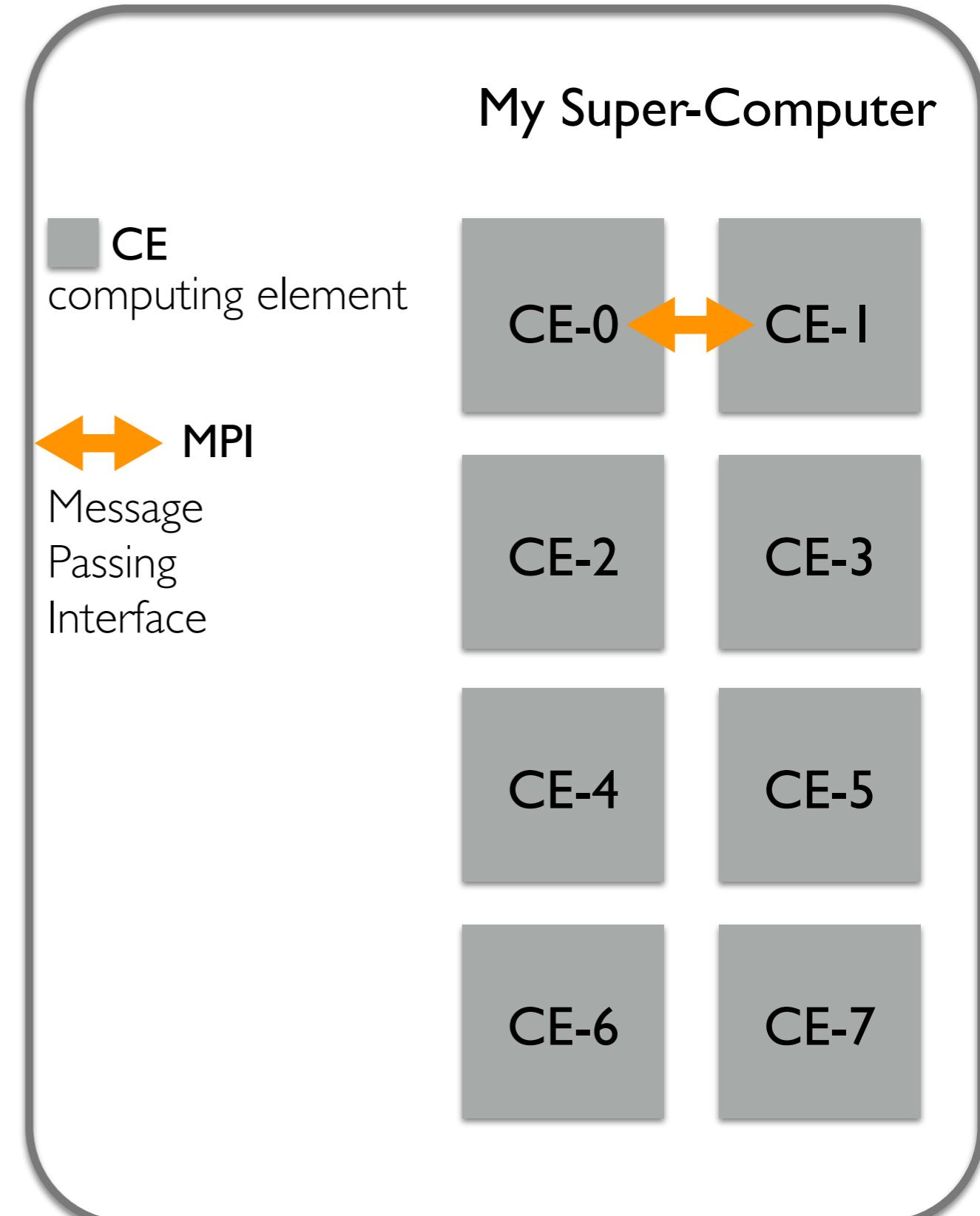
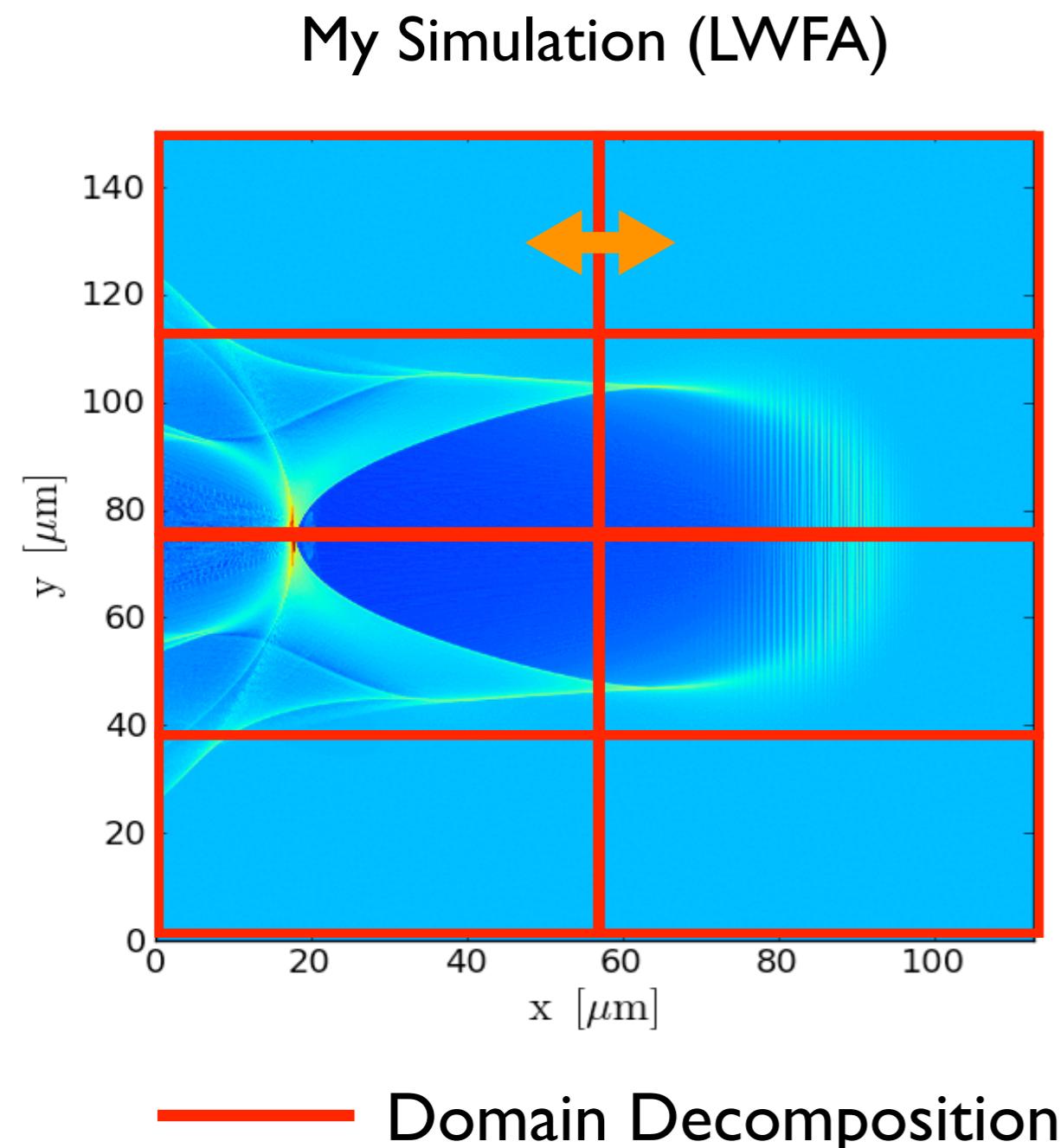
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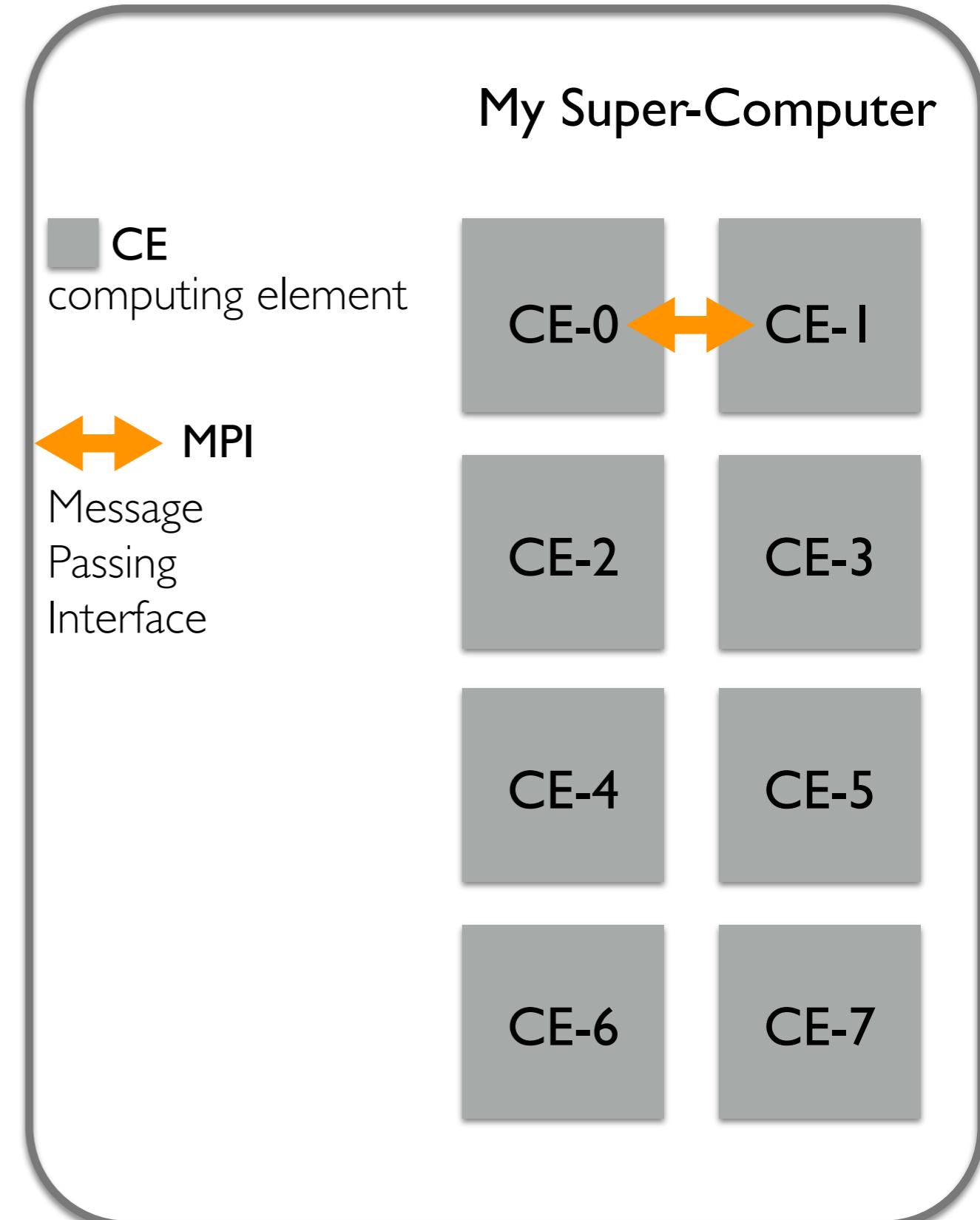
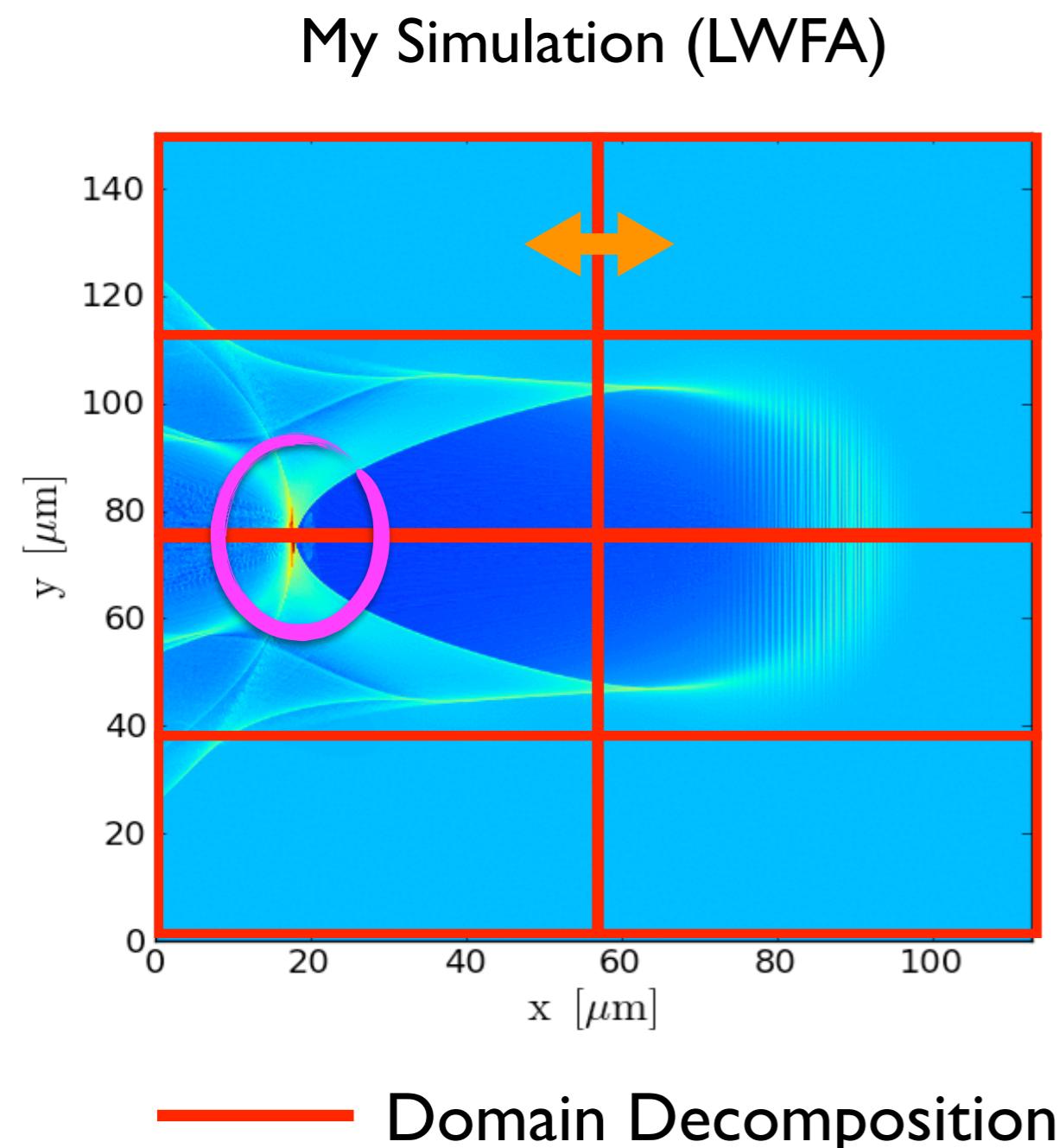
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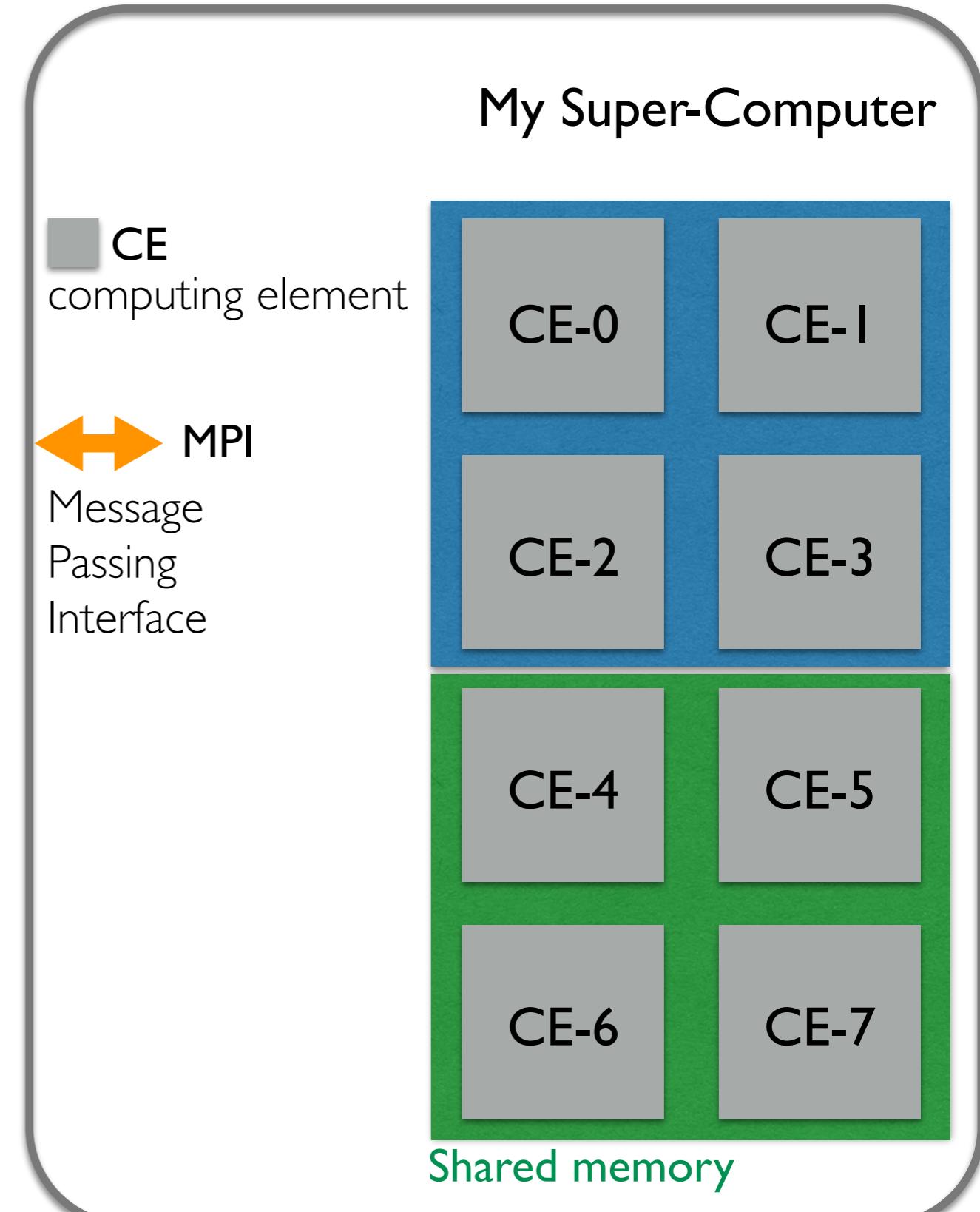
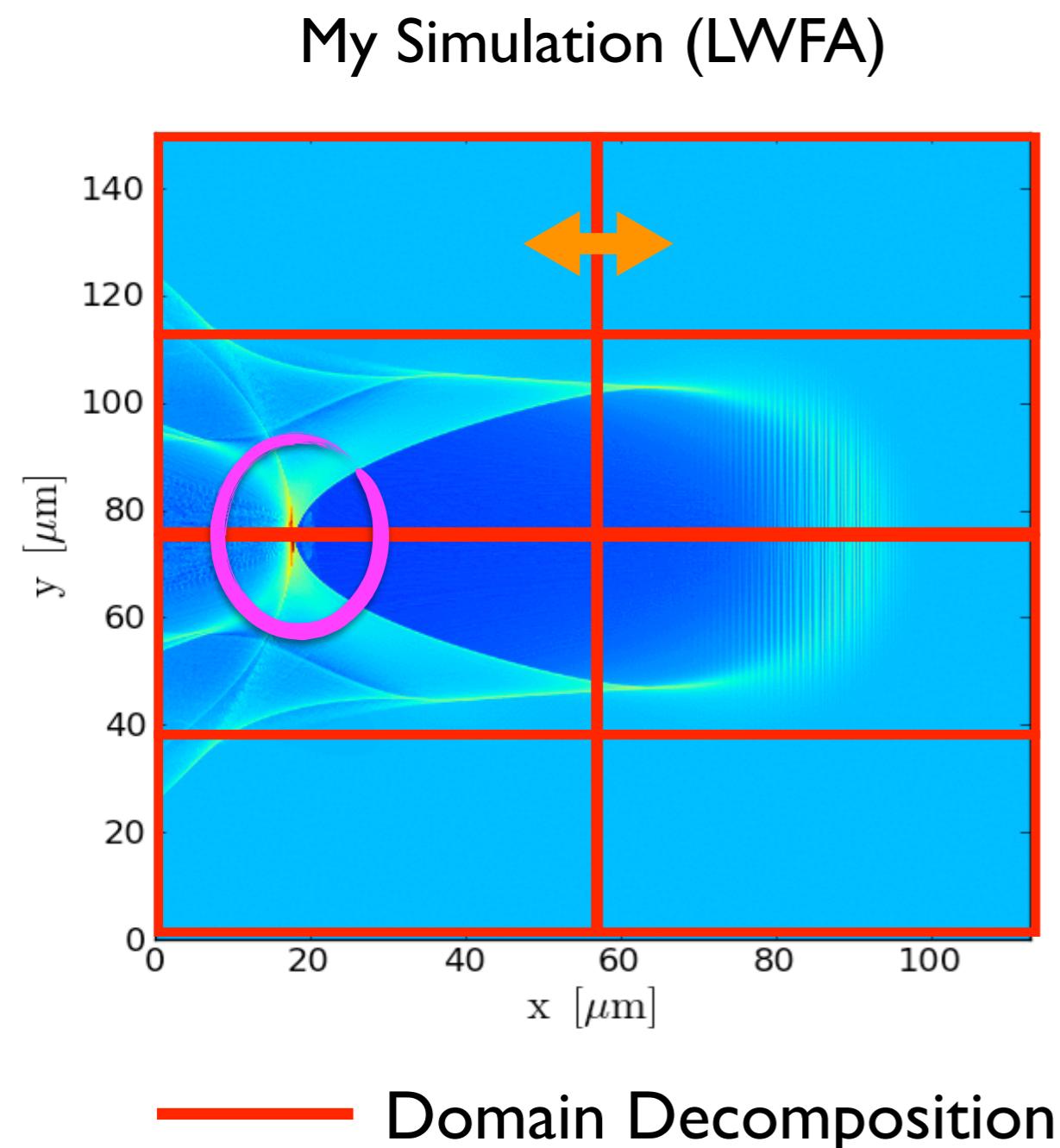
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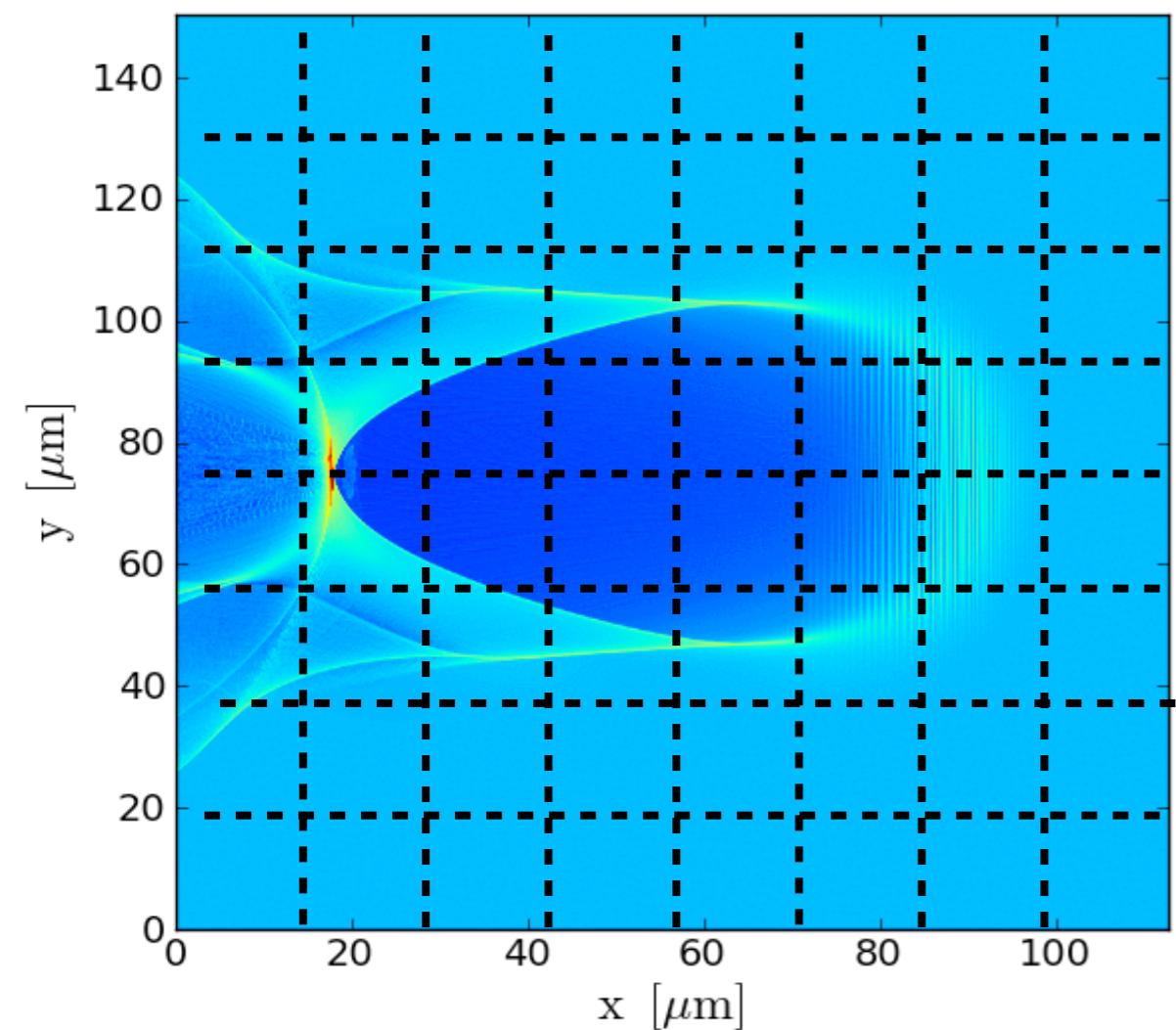
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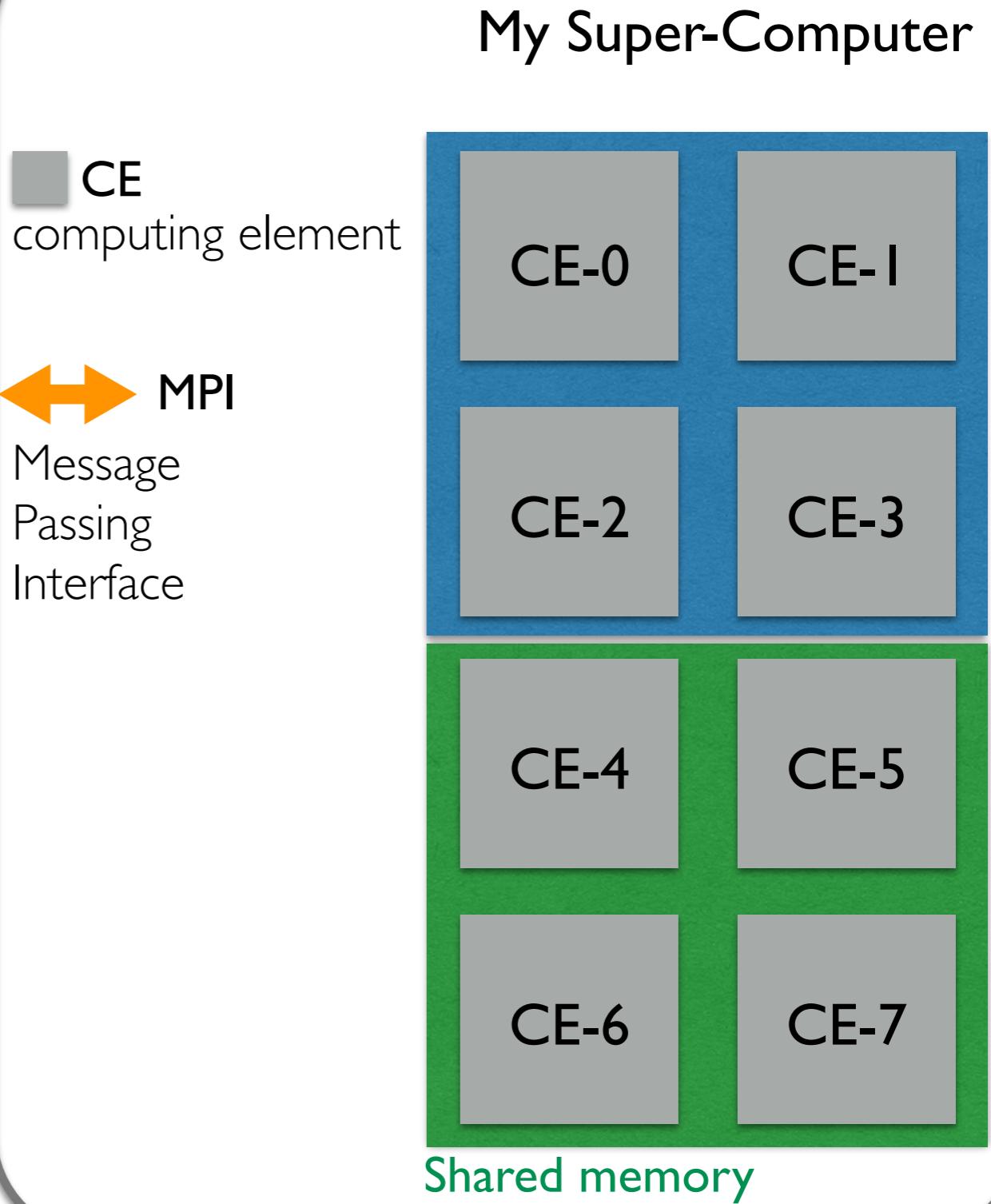
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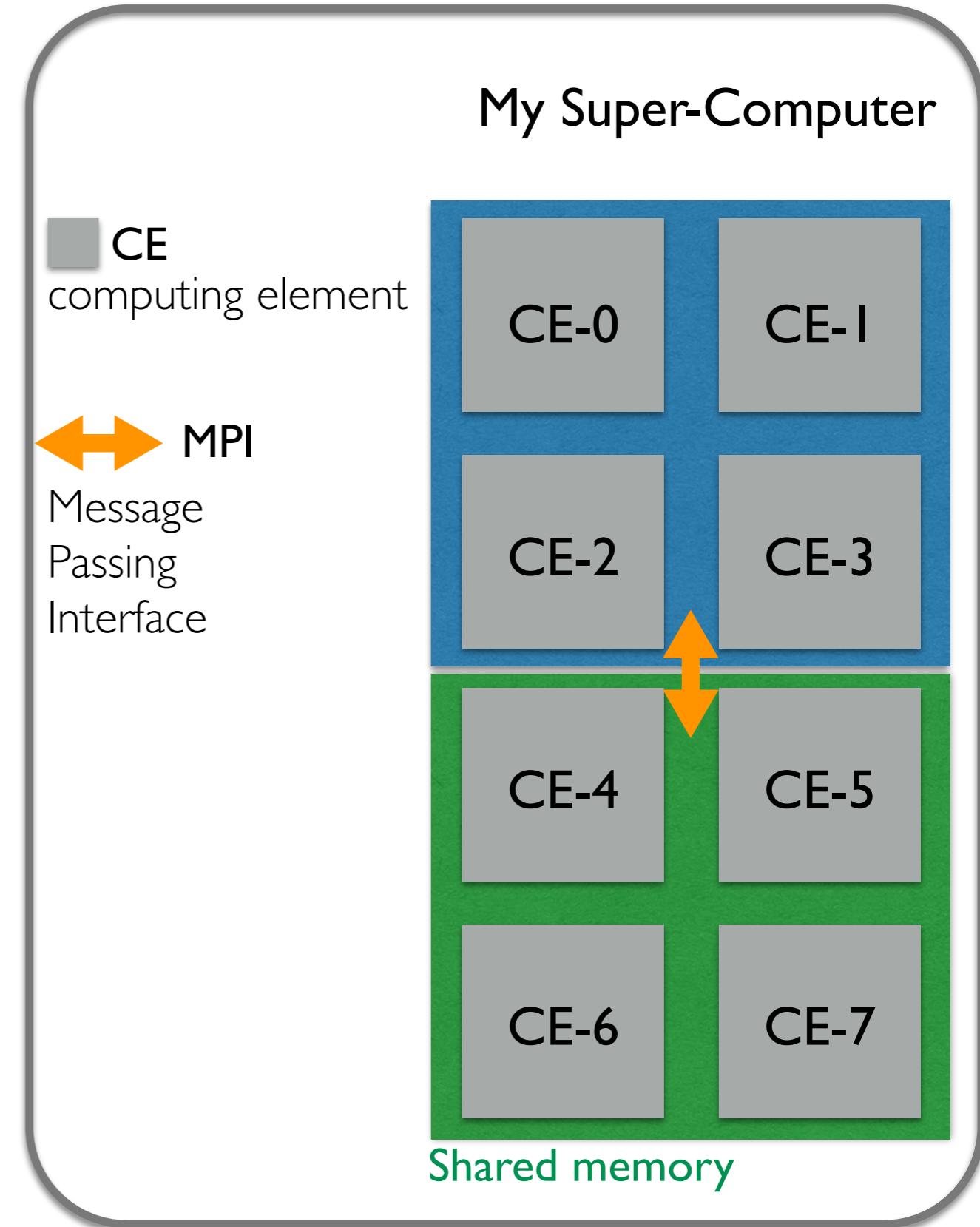
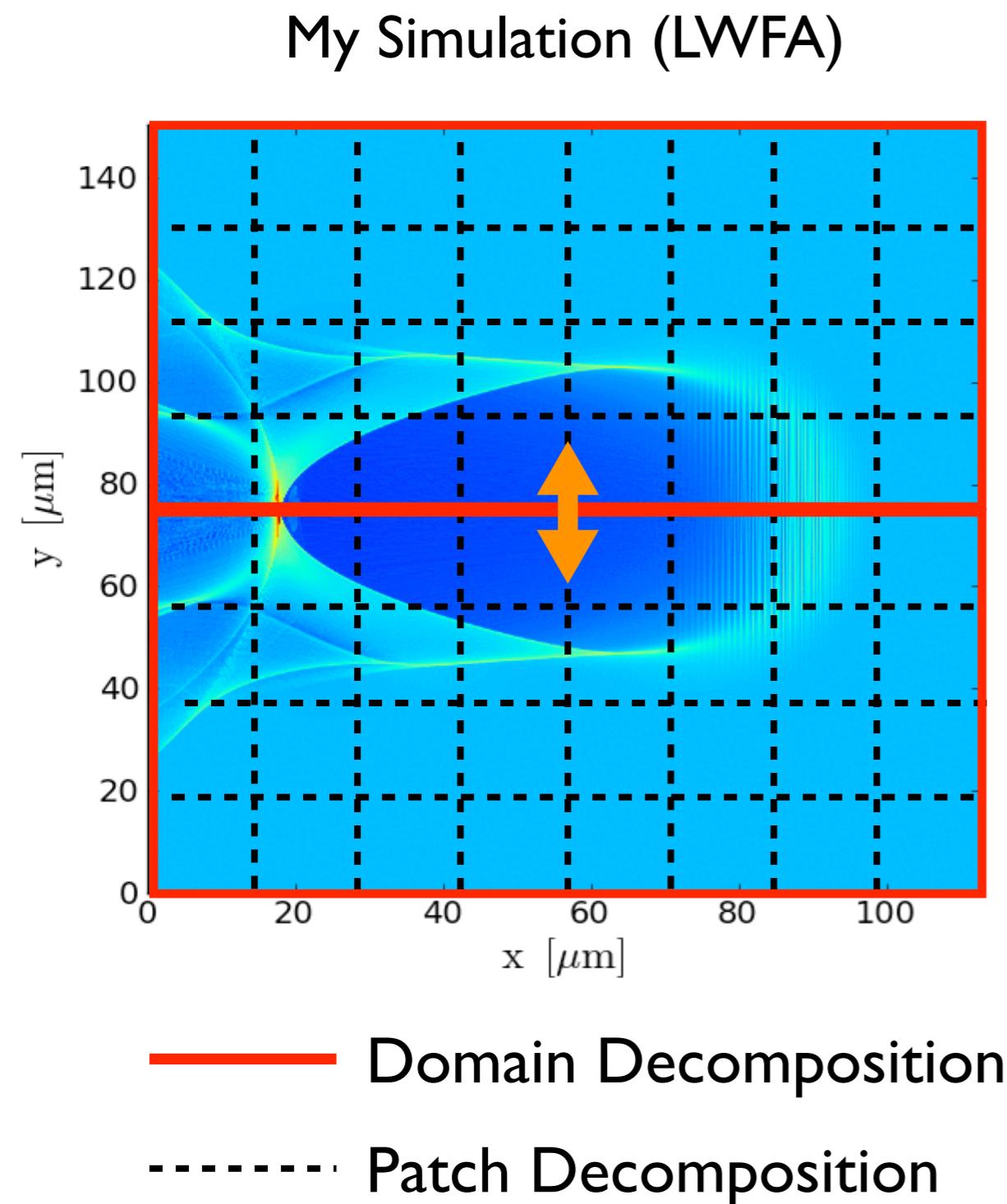


----- Patch Decomposition



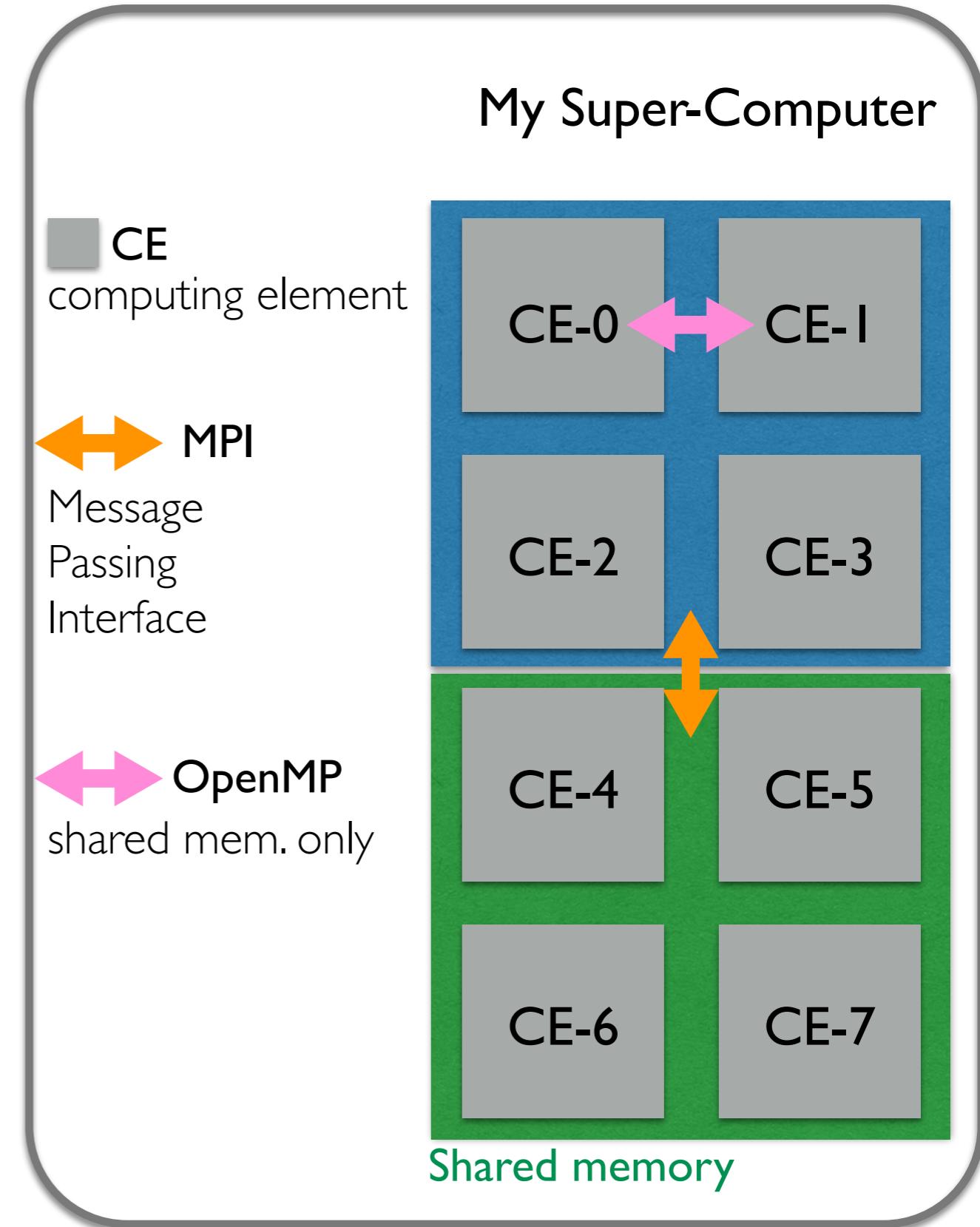
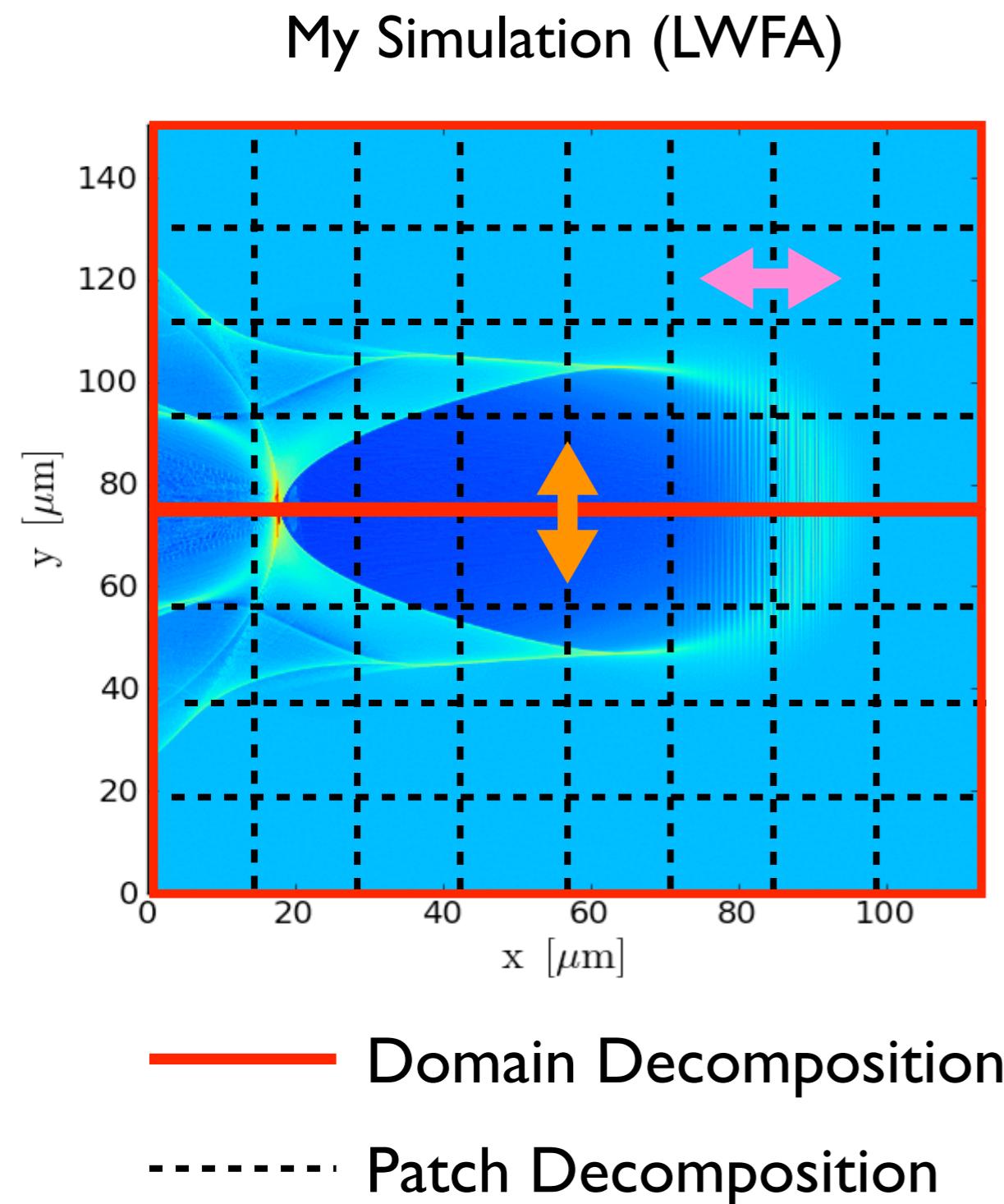
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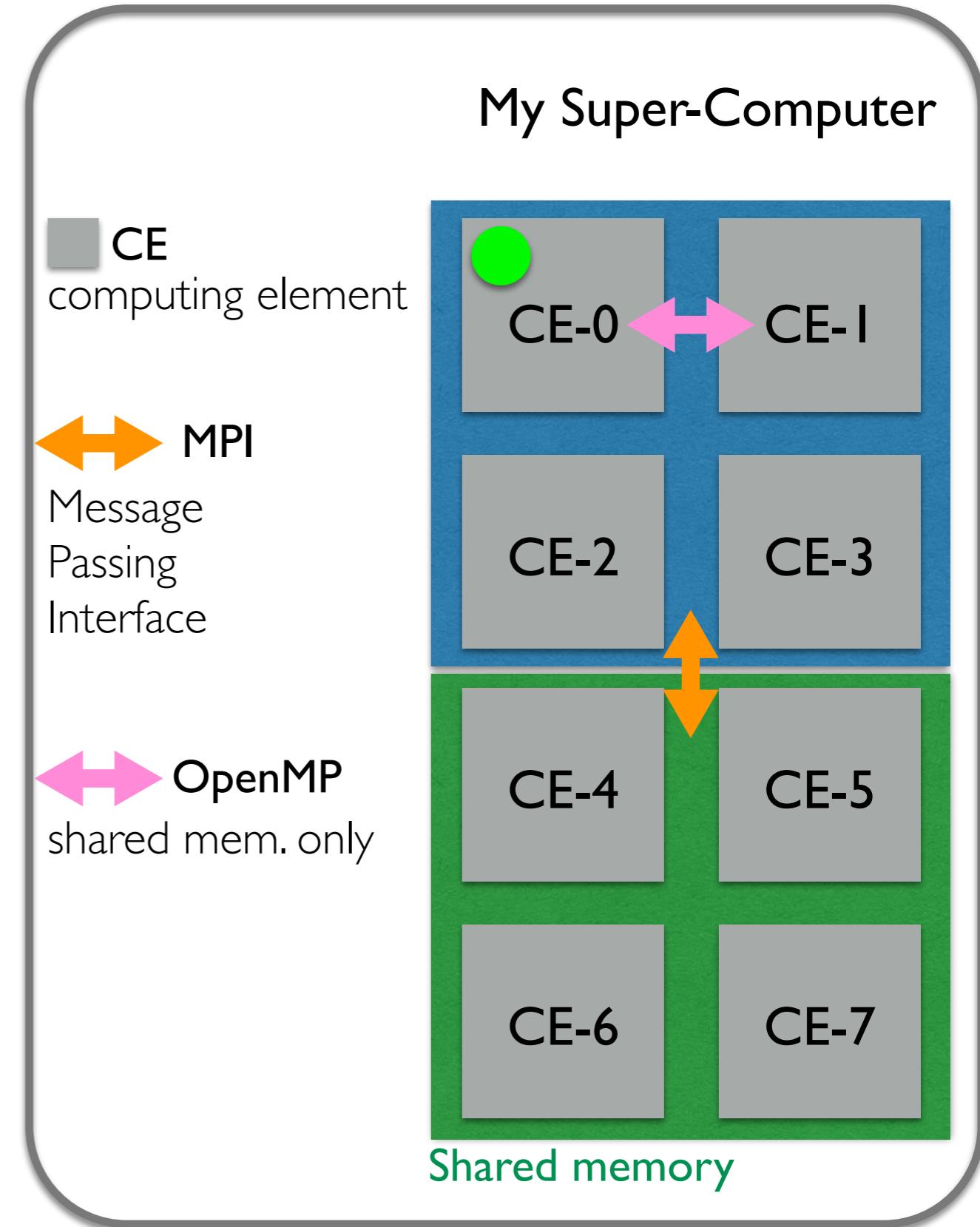
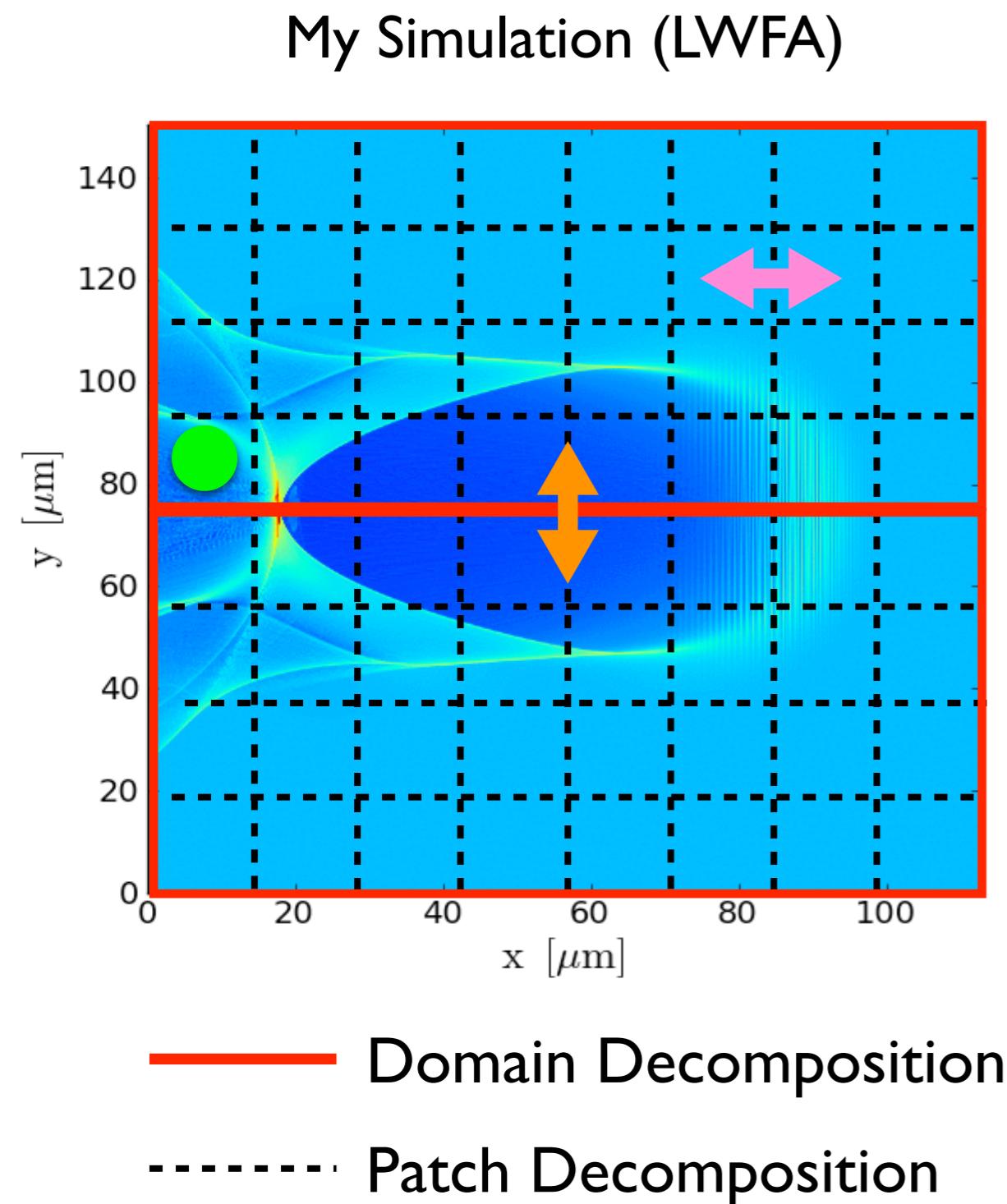
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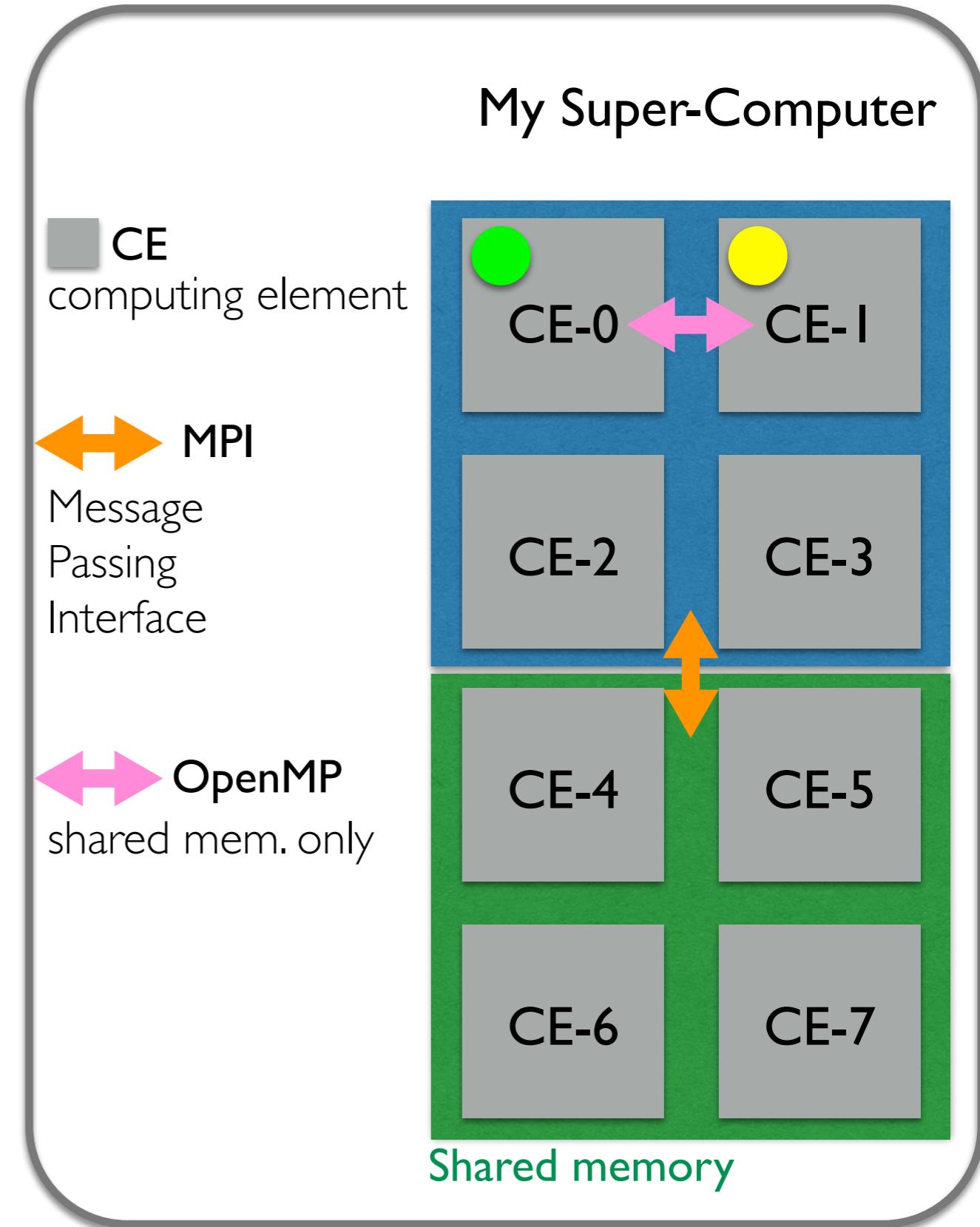
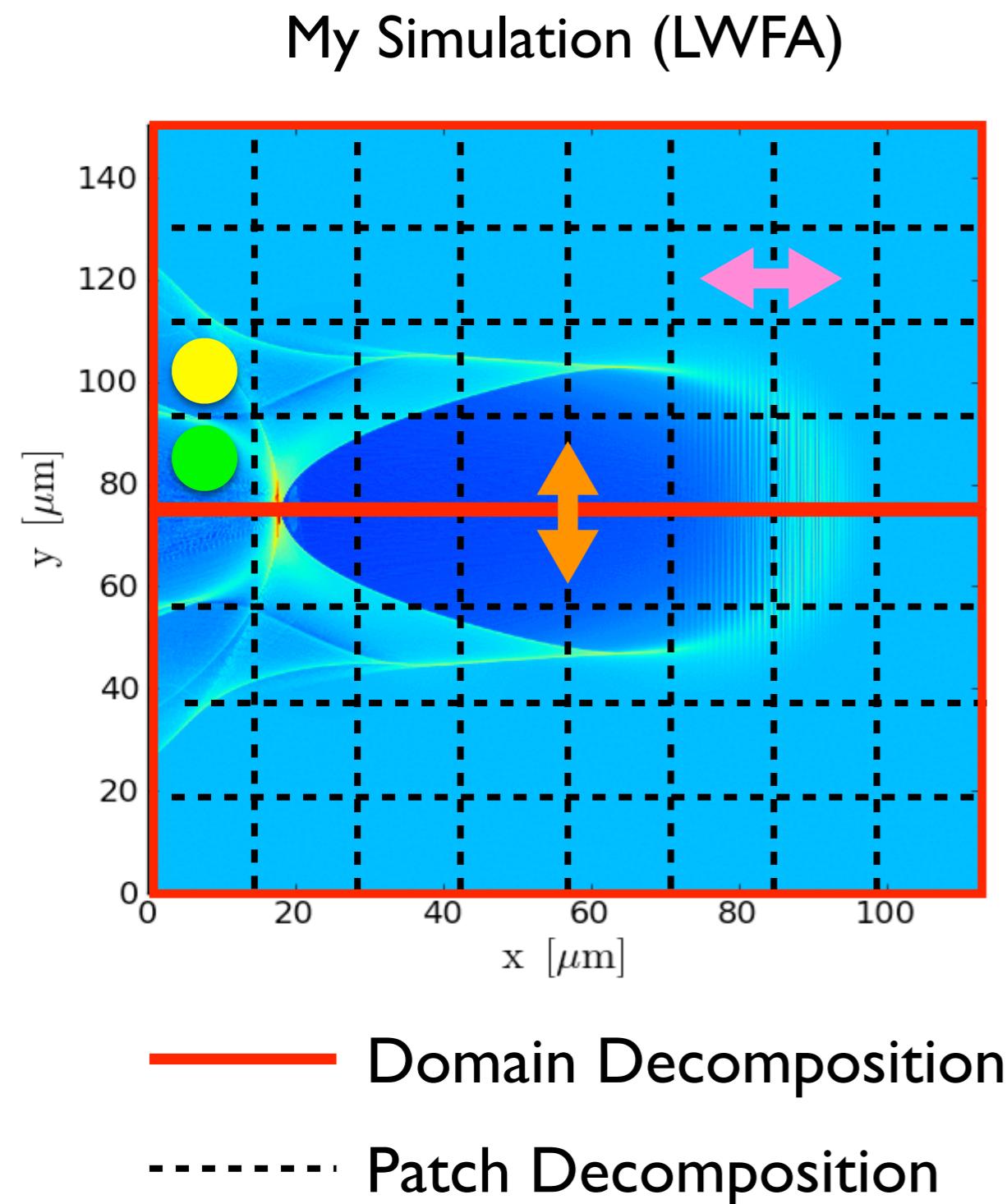
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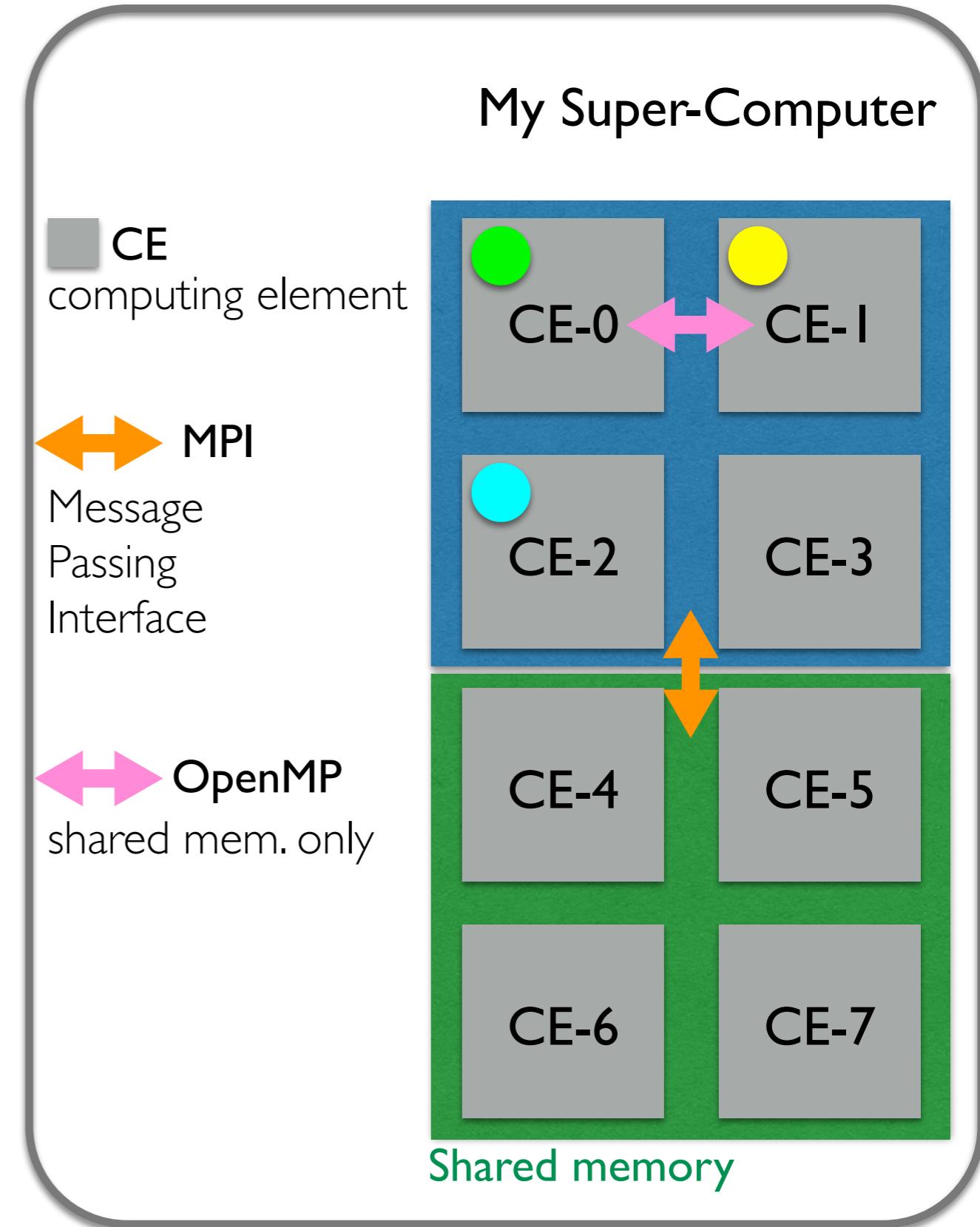
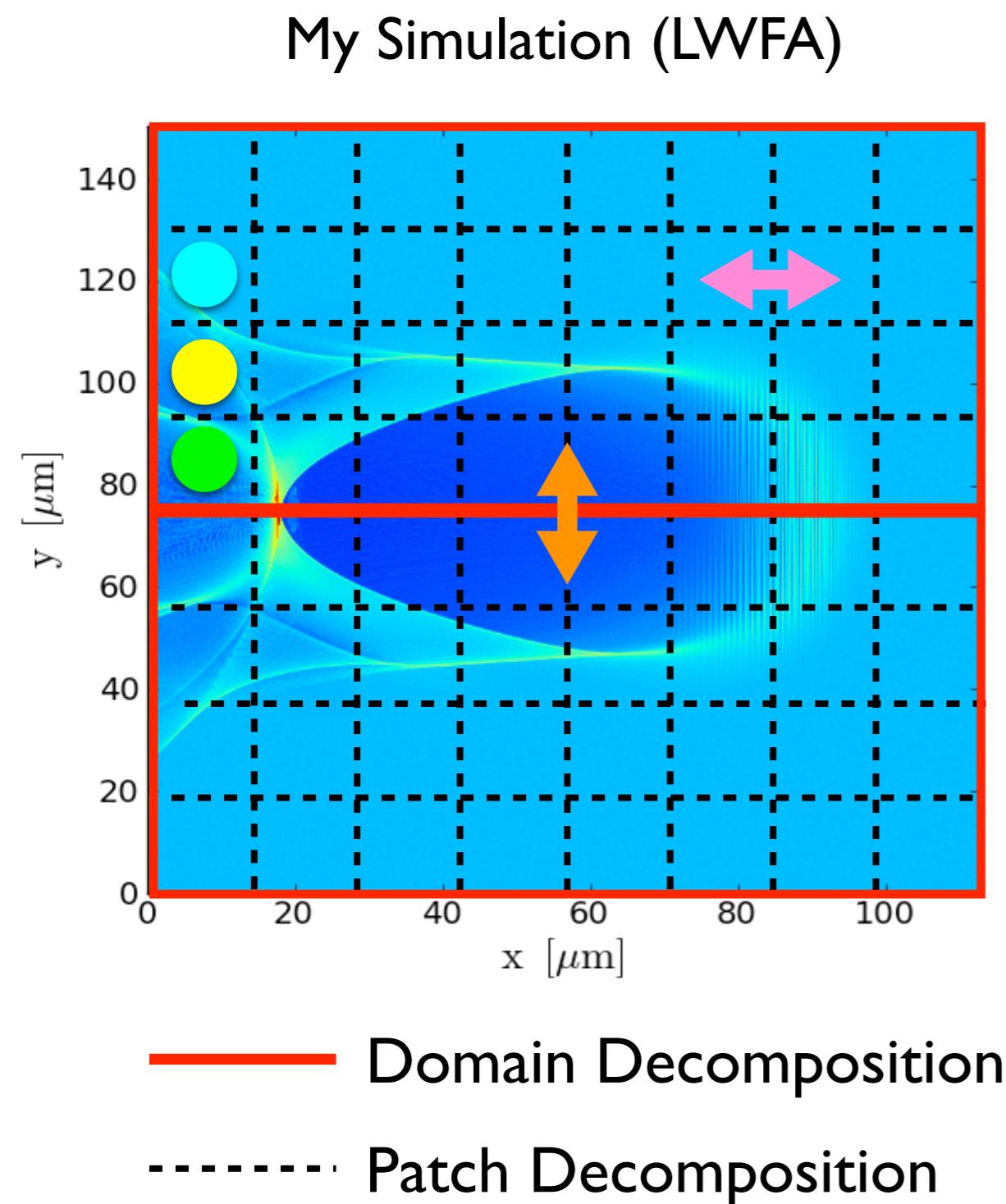
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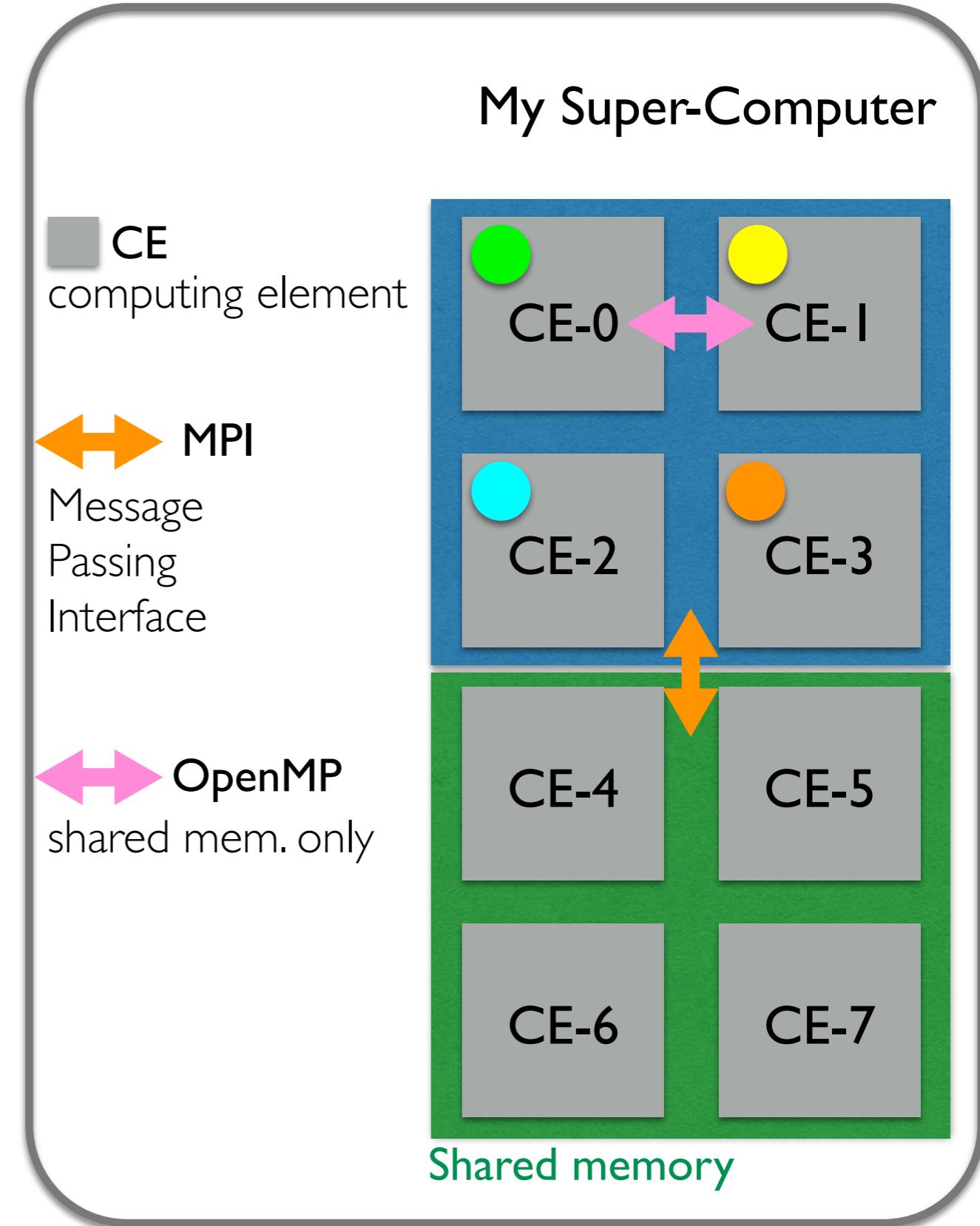
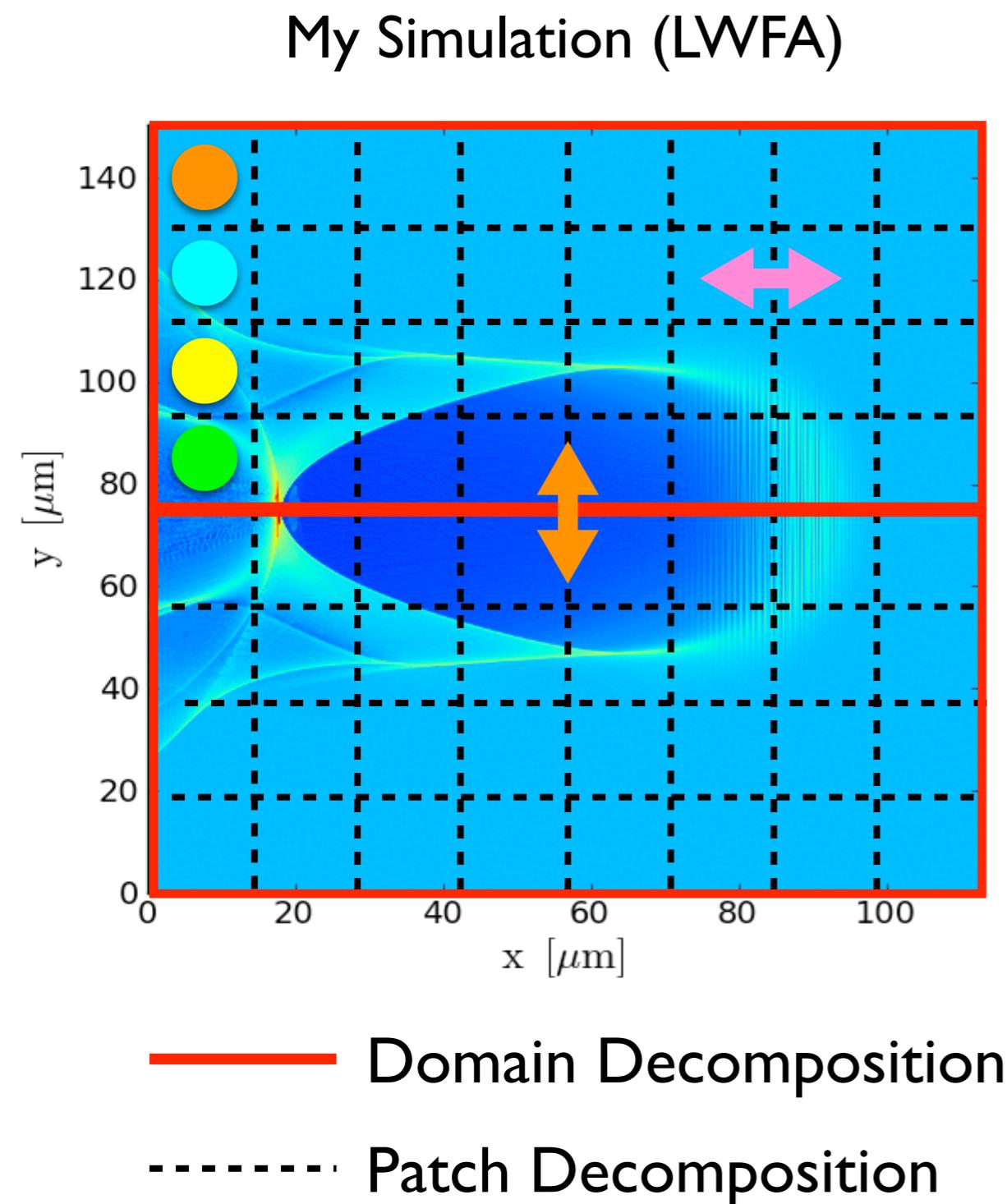
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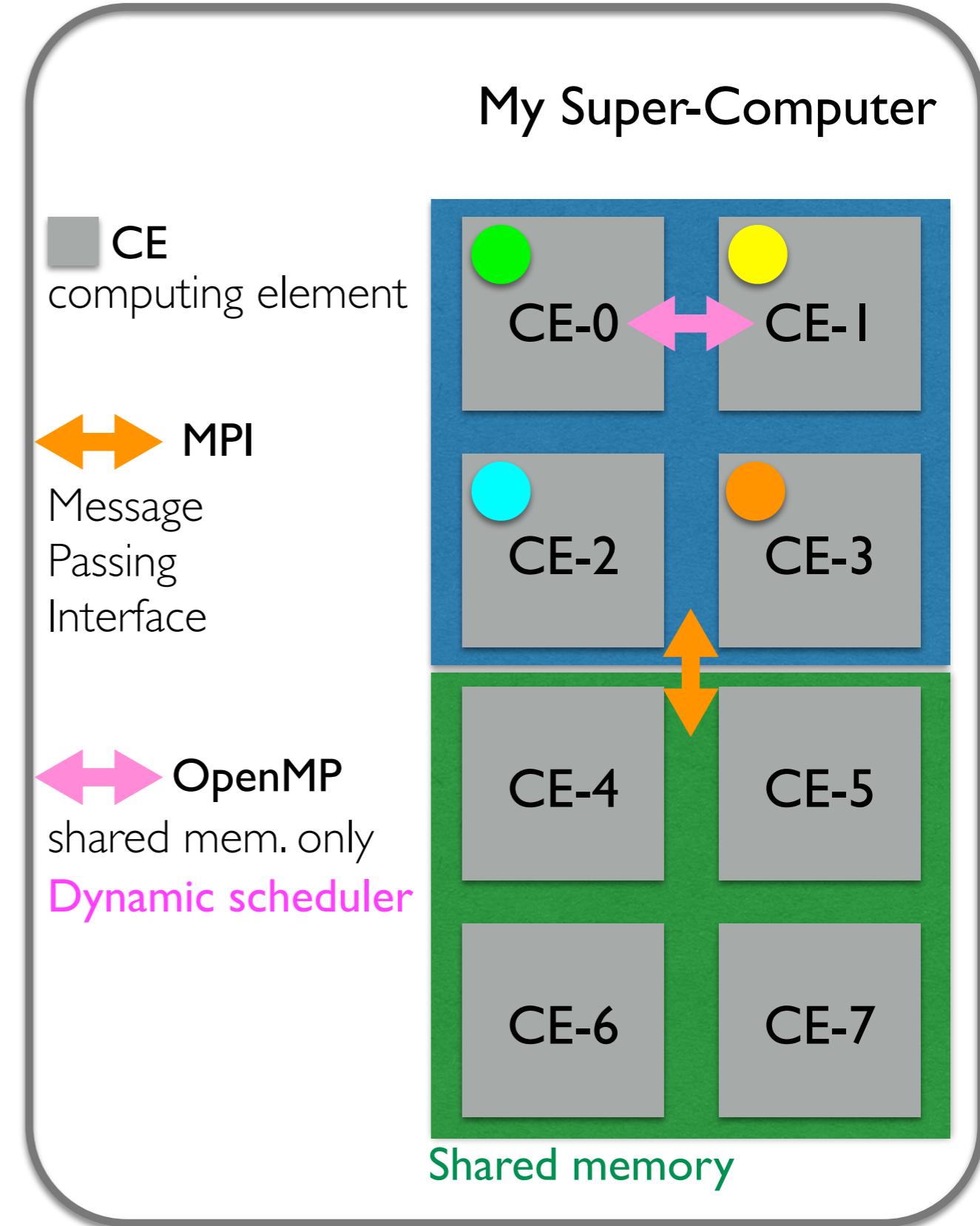
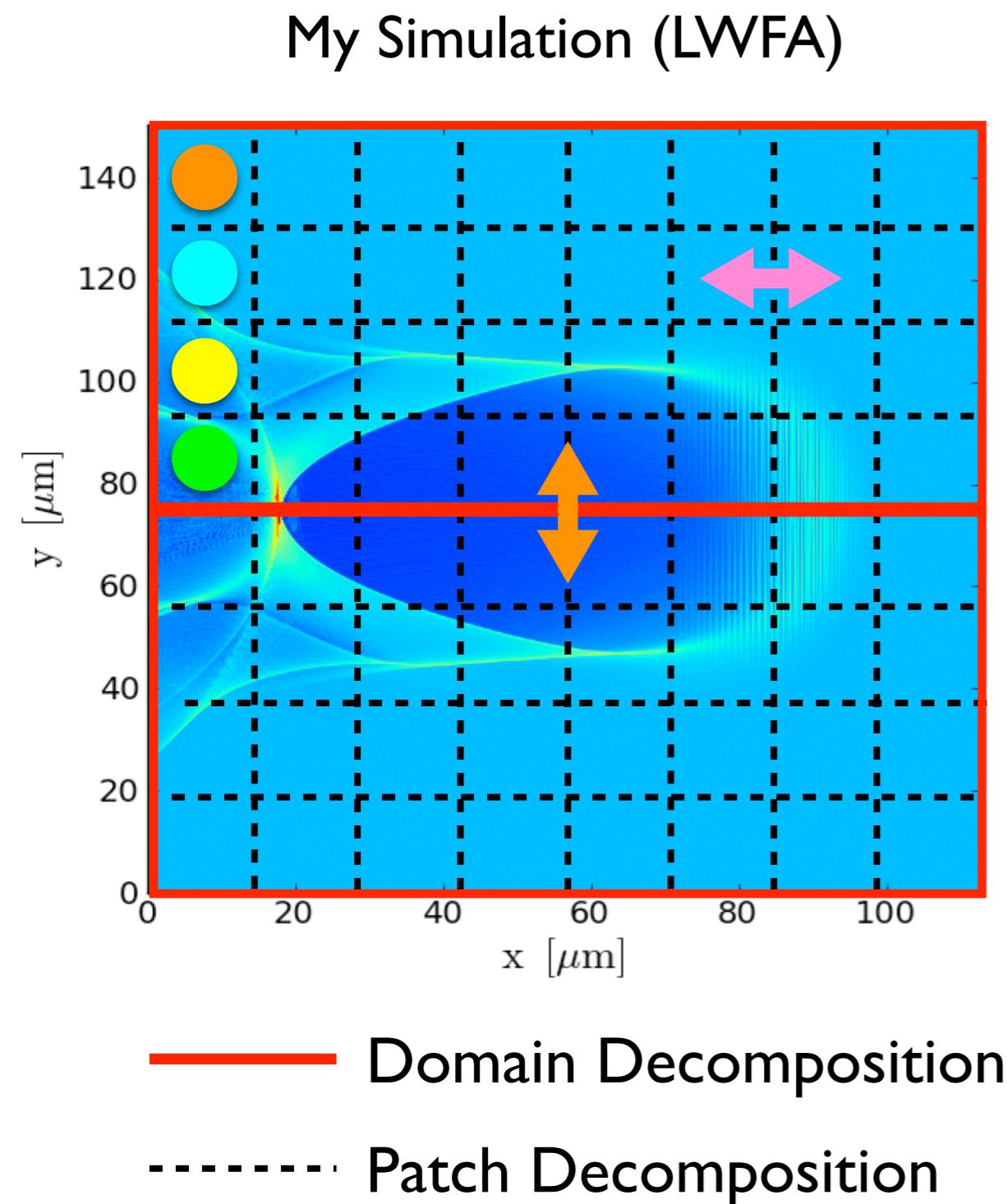
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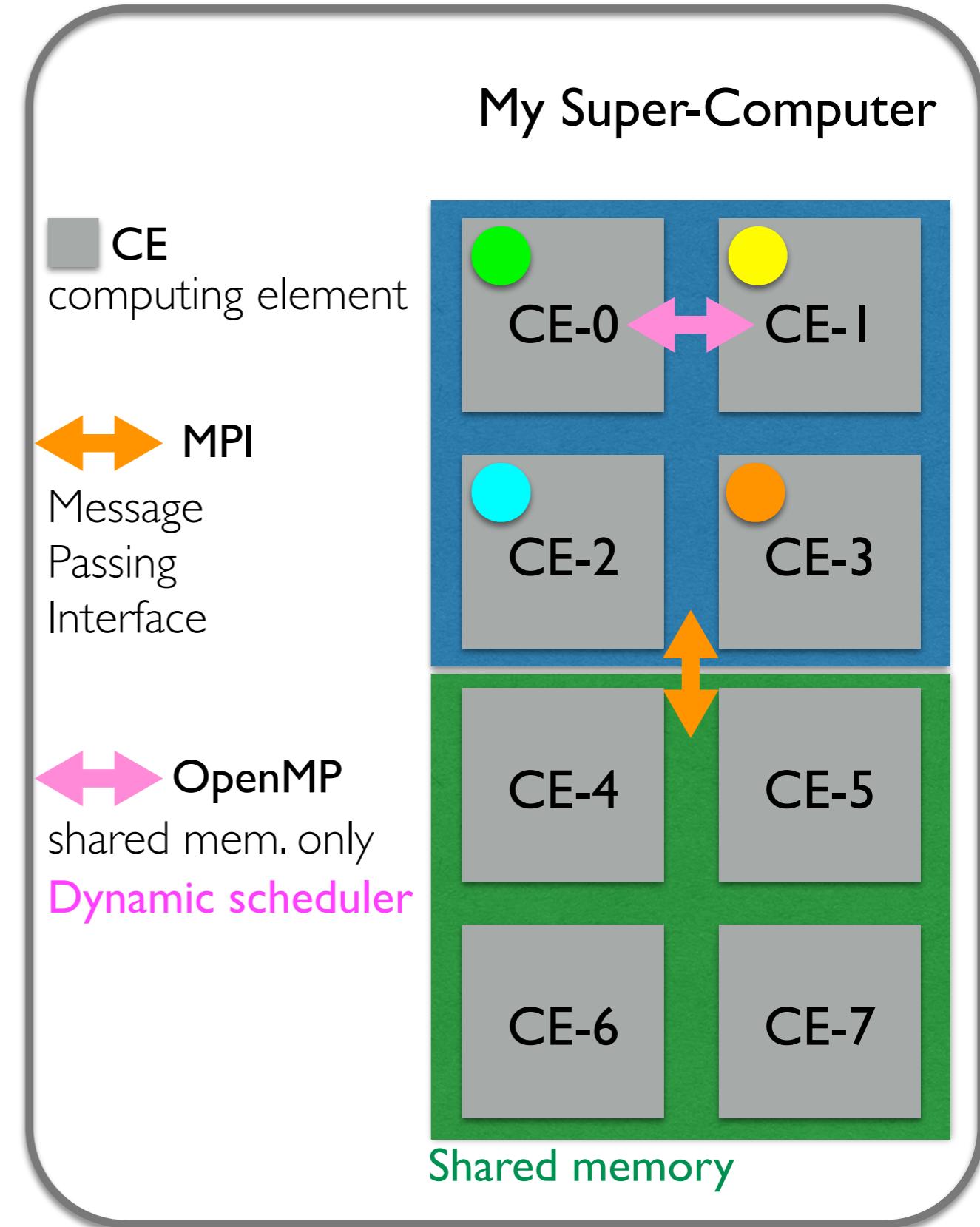
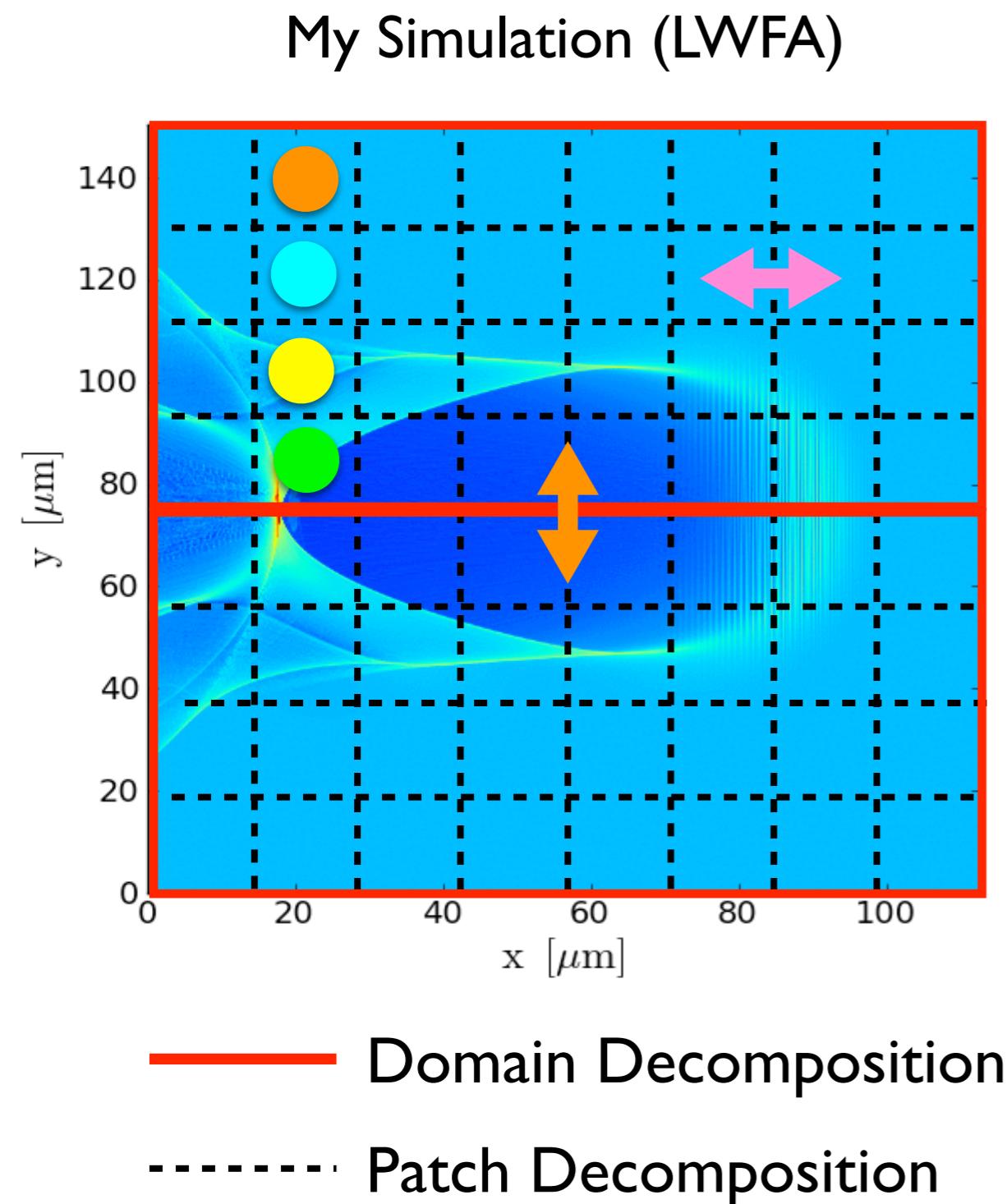
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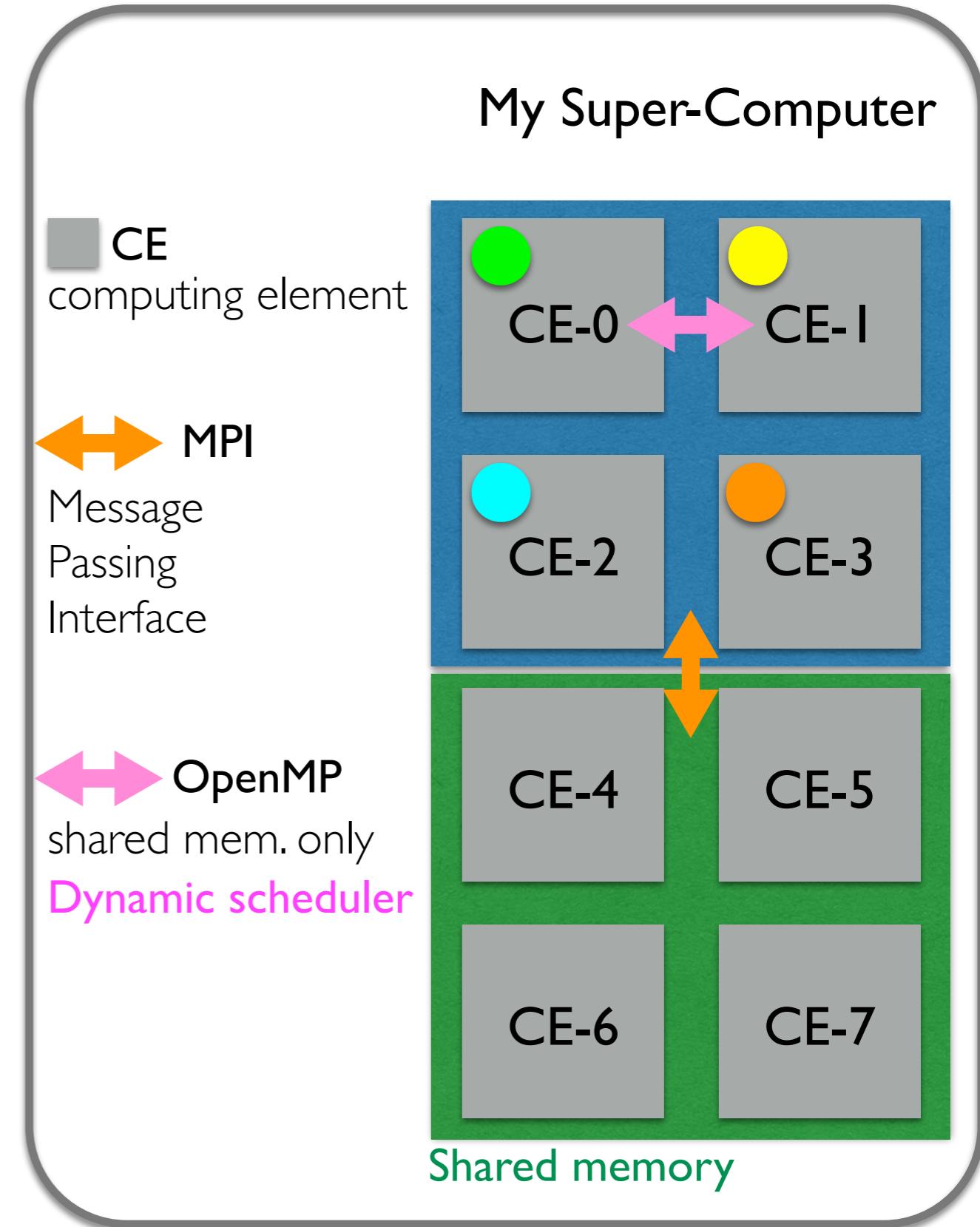
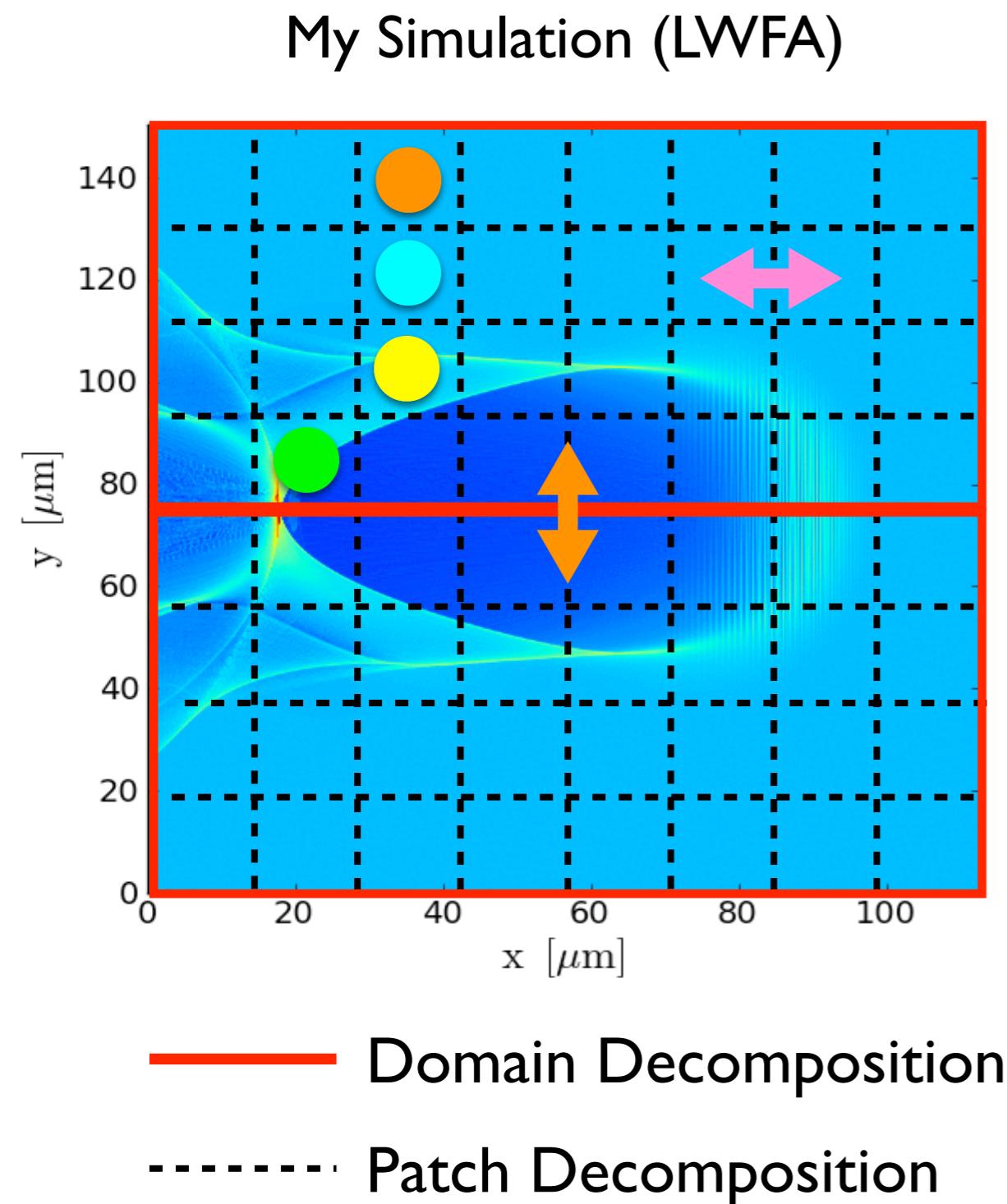
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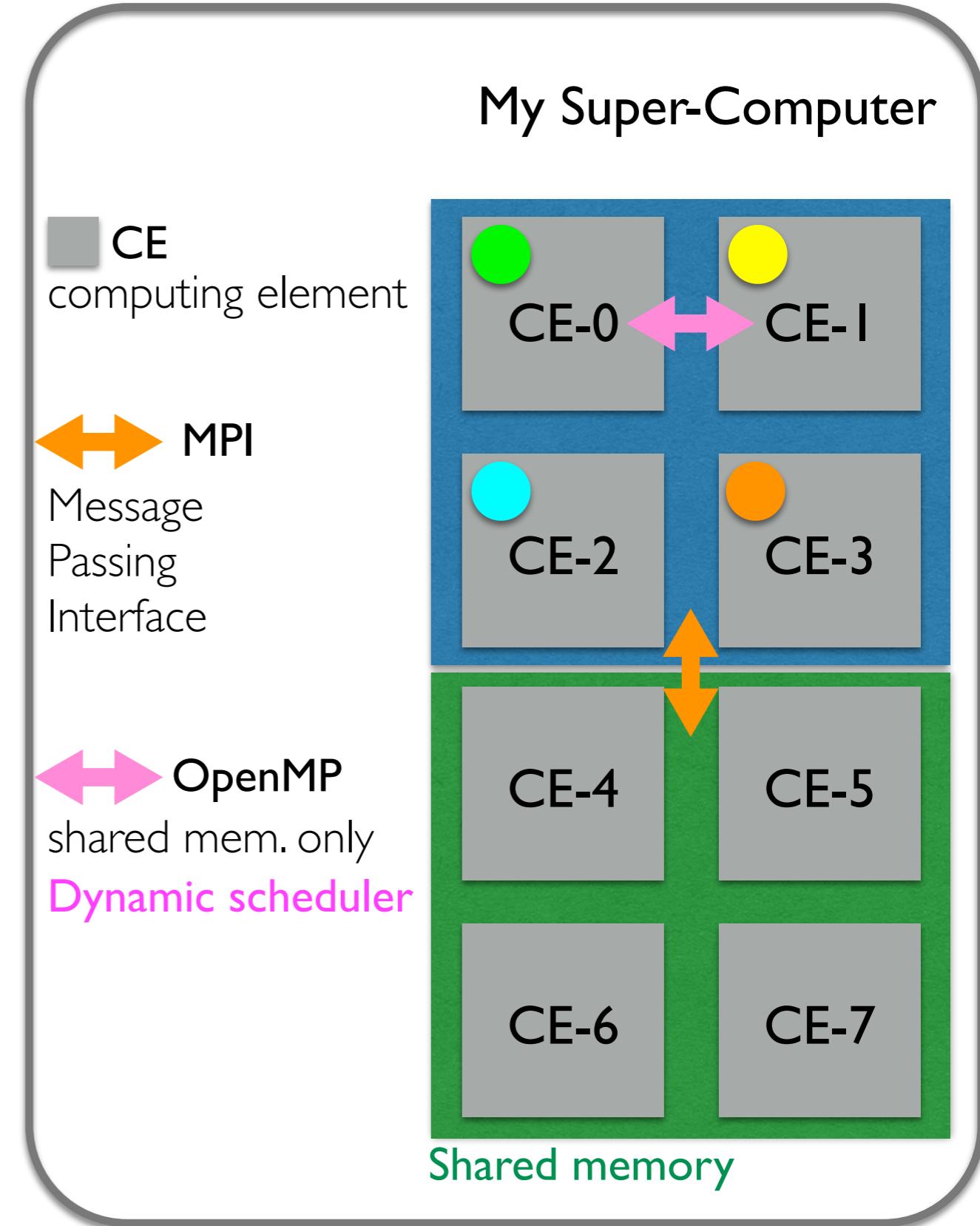
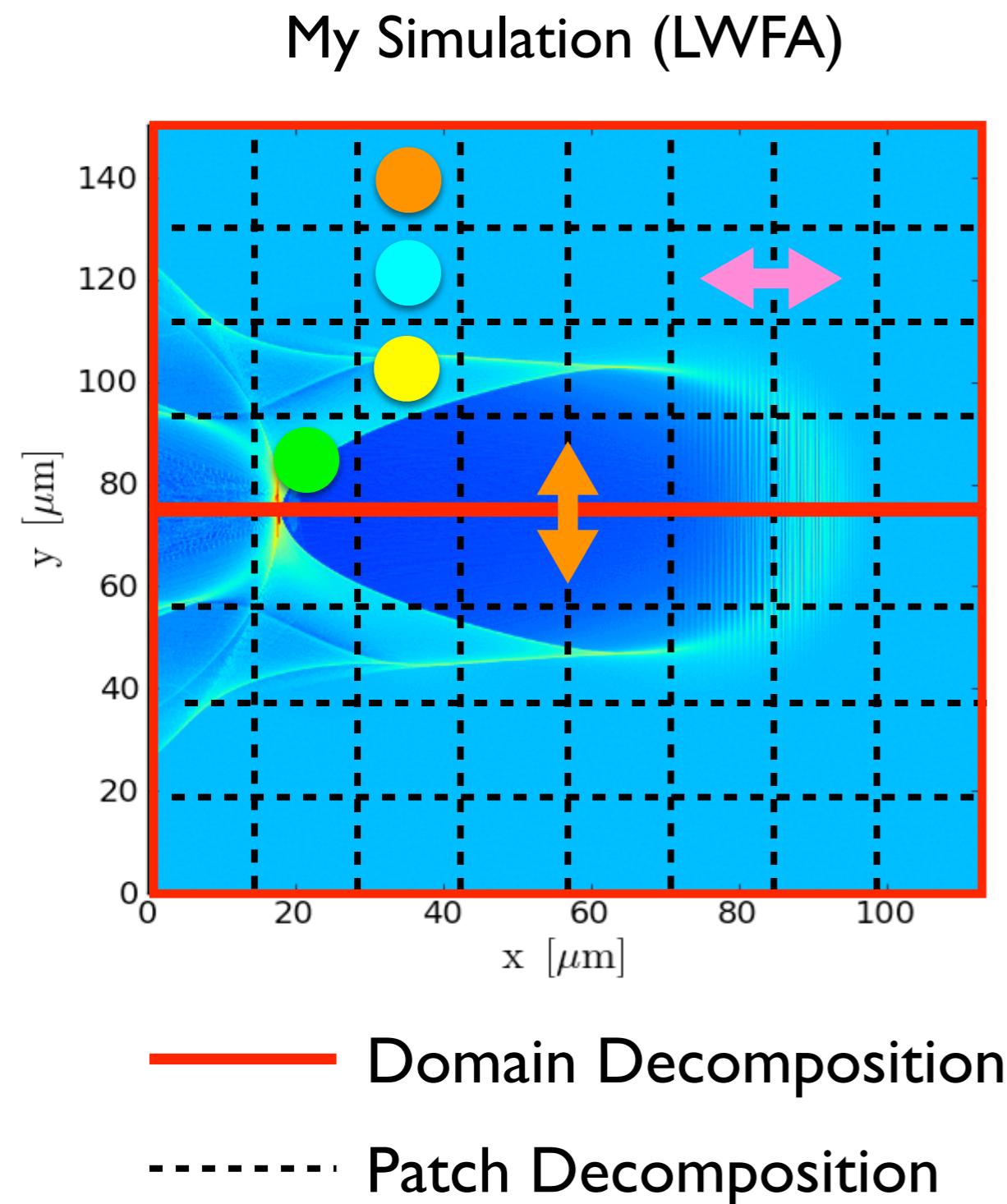
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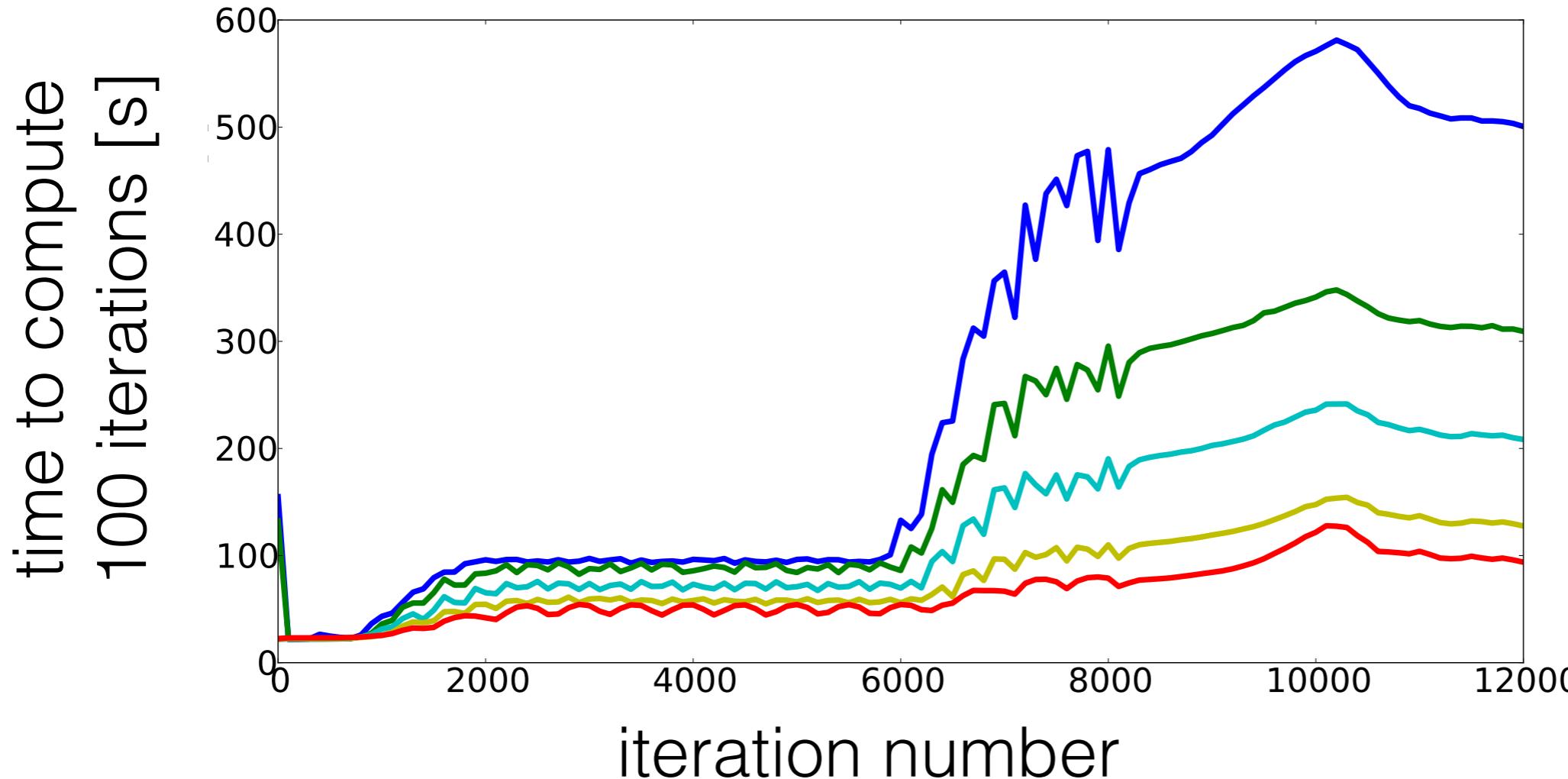
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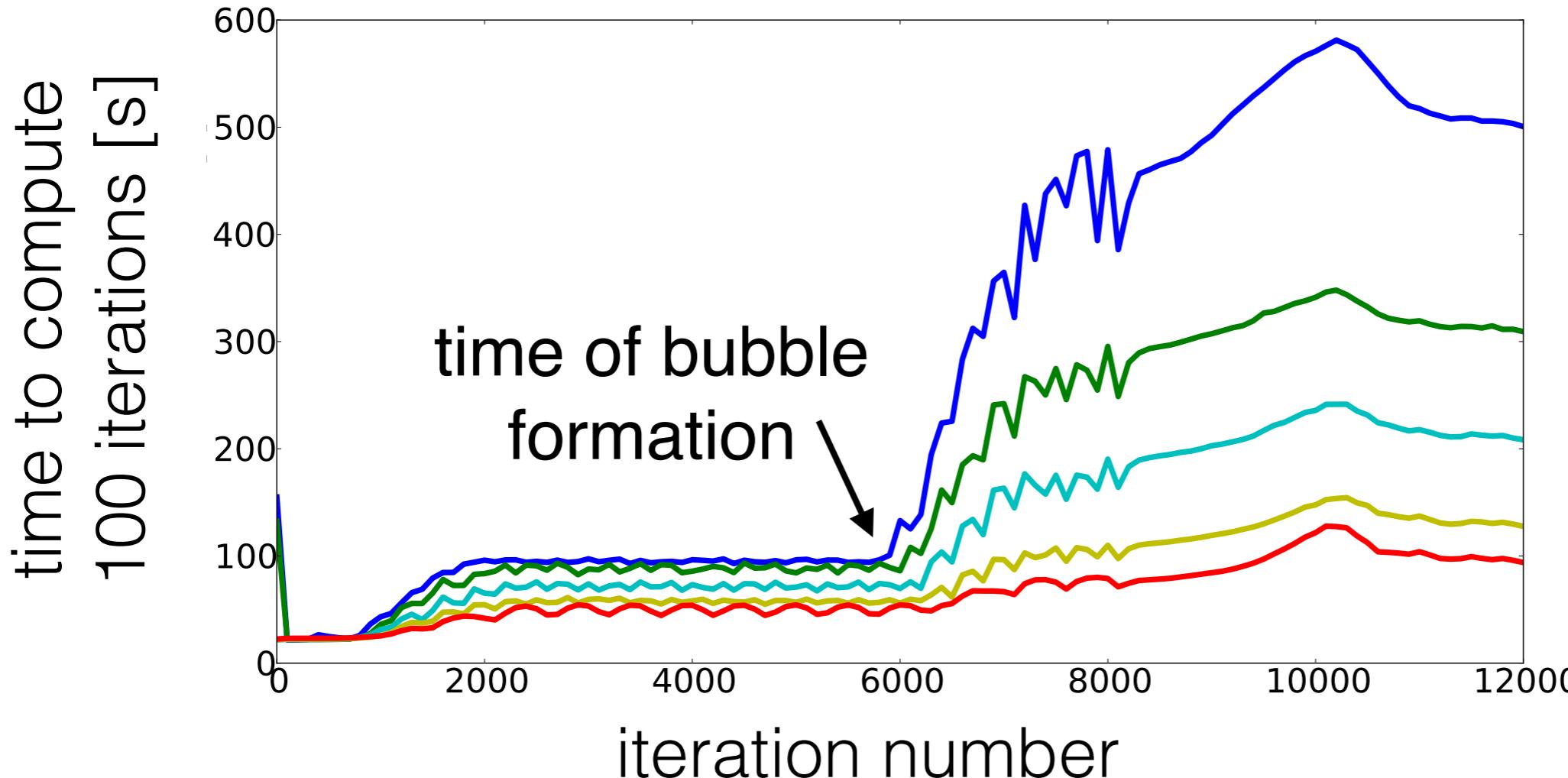
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Hybrid parallelization significantly improves performance



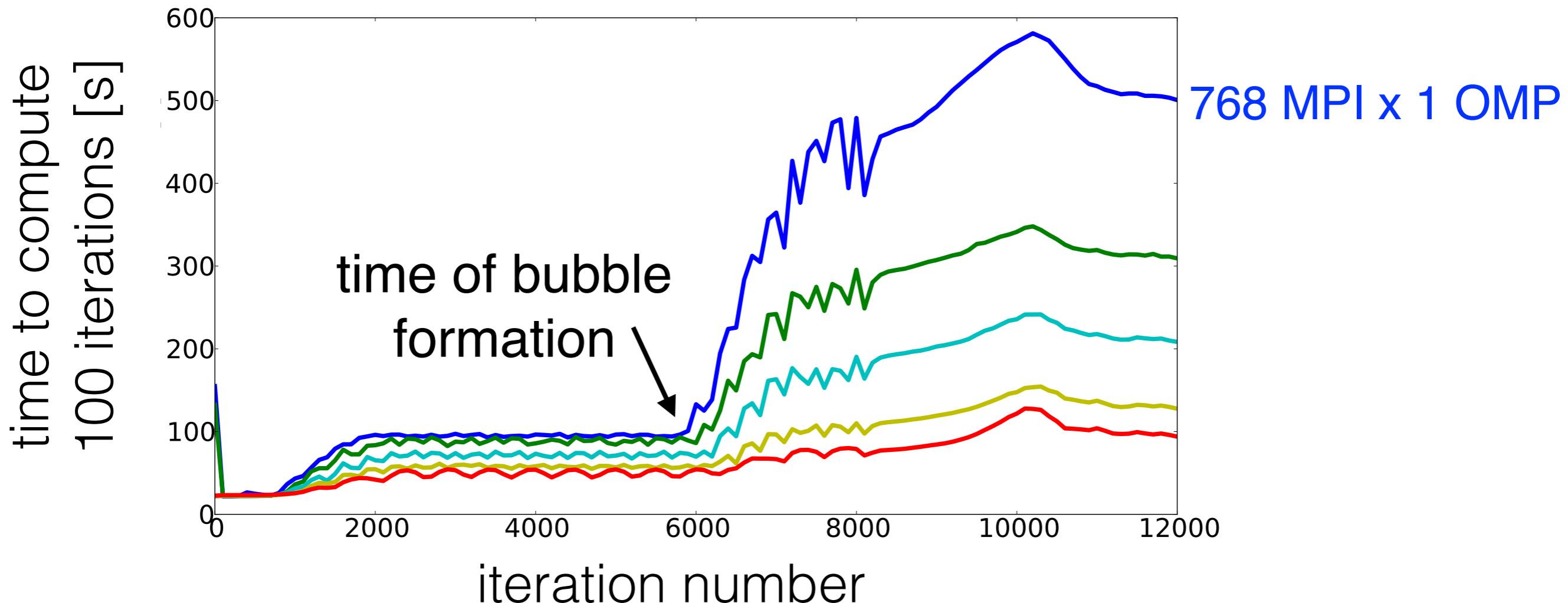
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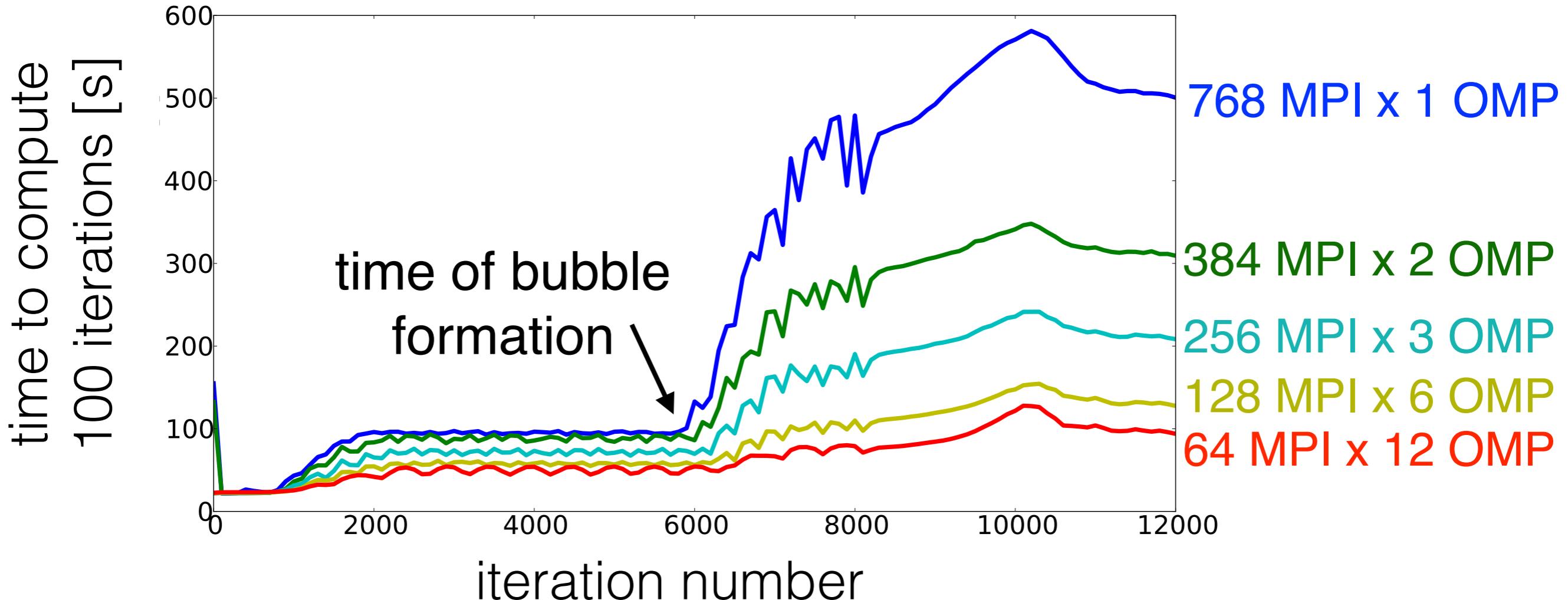
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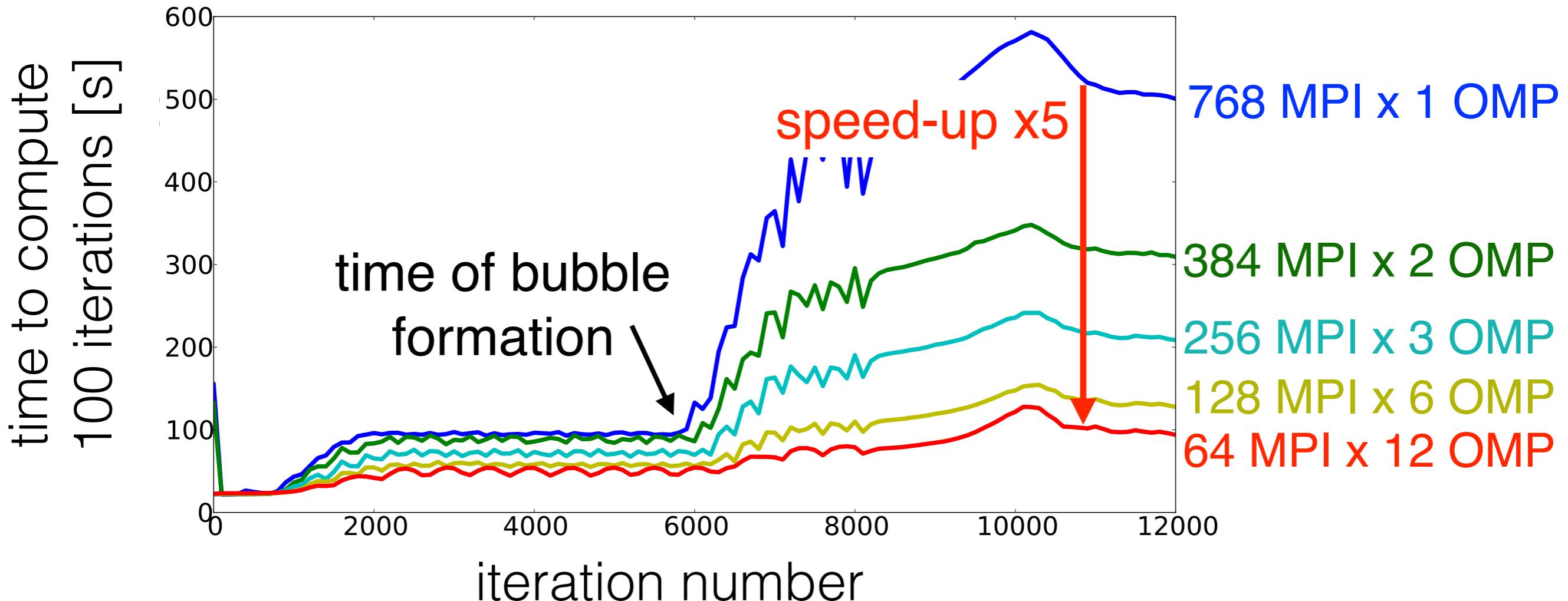
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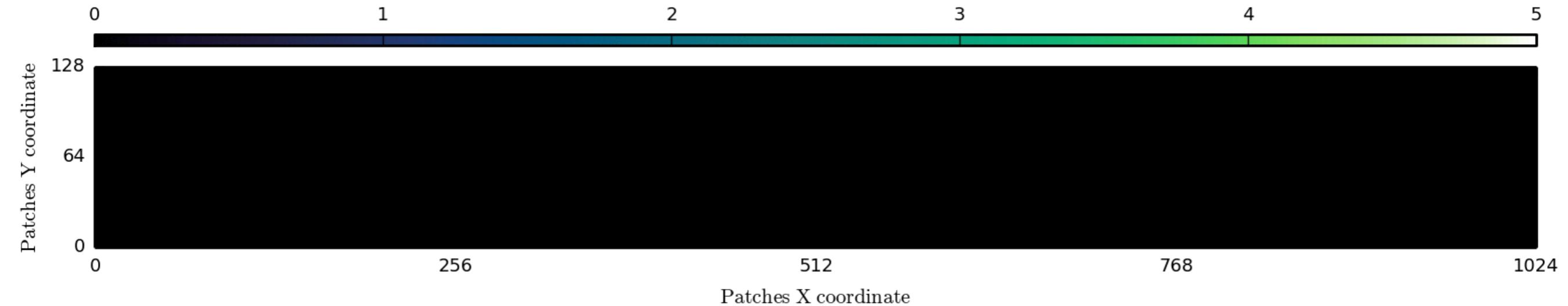


Step 1: Parallelization

Dynamic load balancing further improve performances

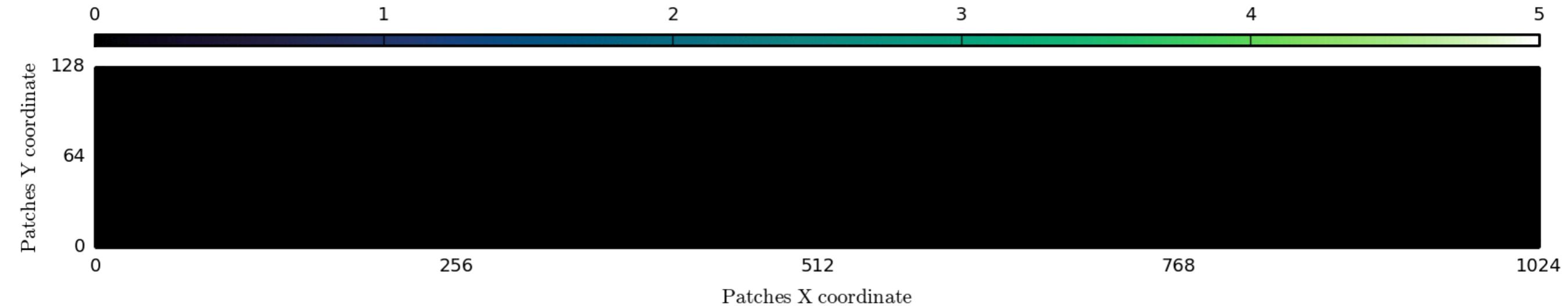
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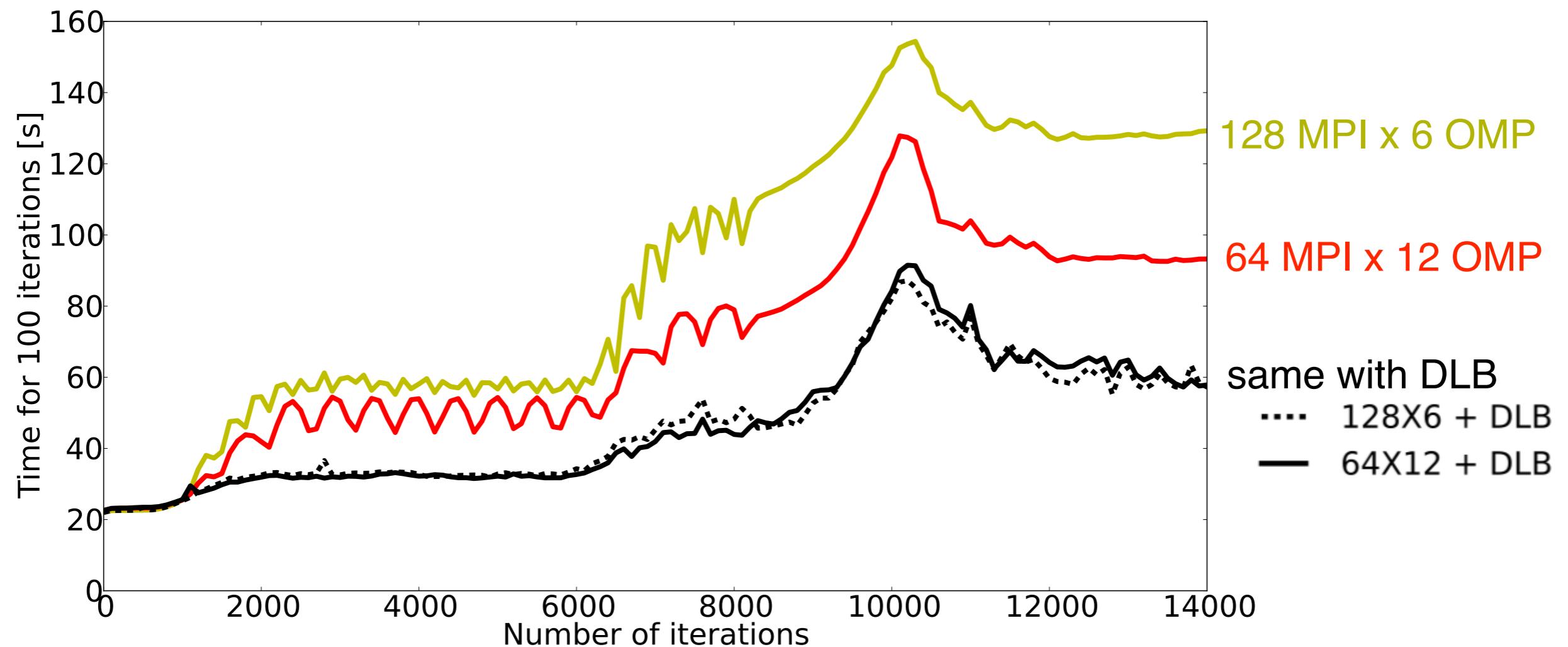
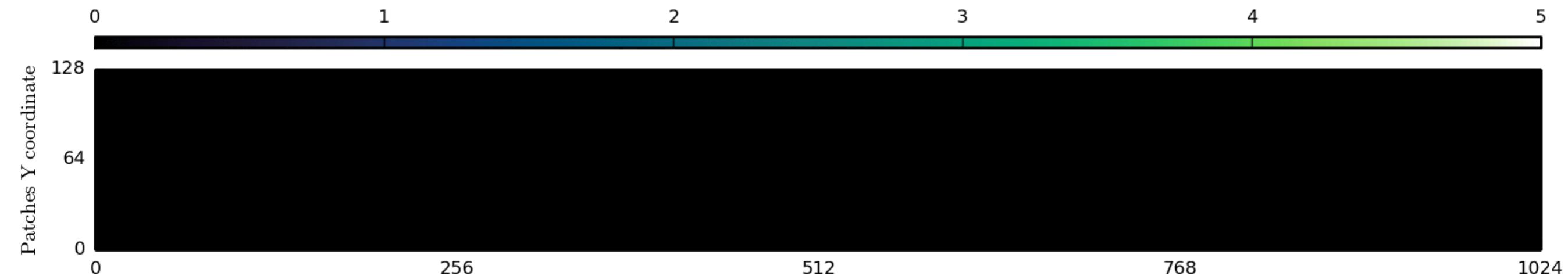
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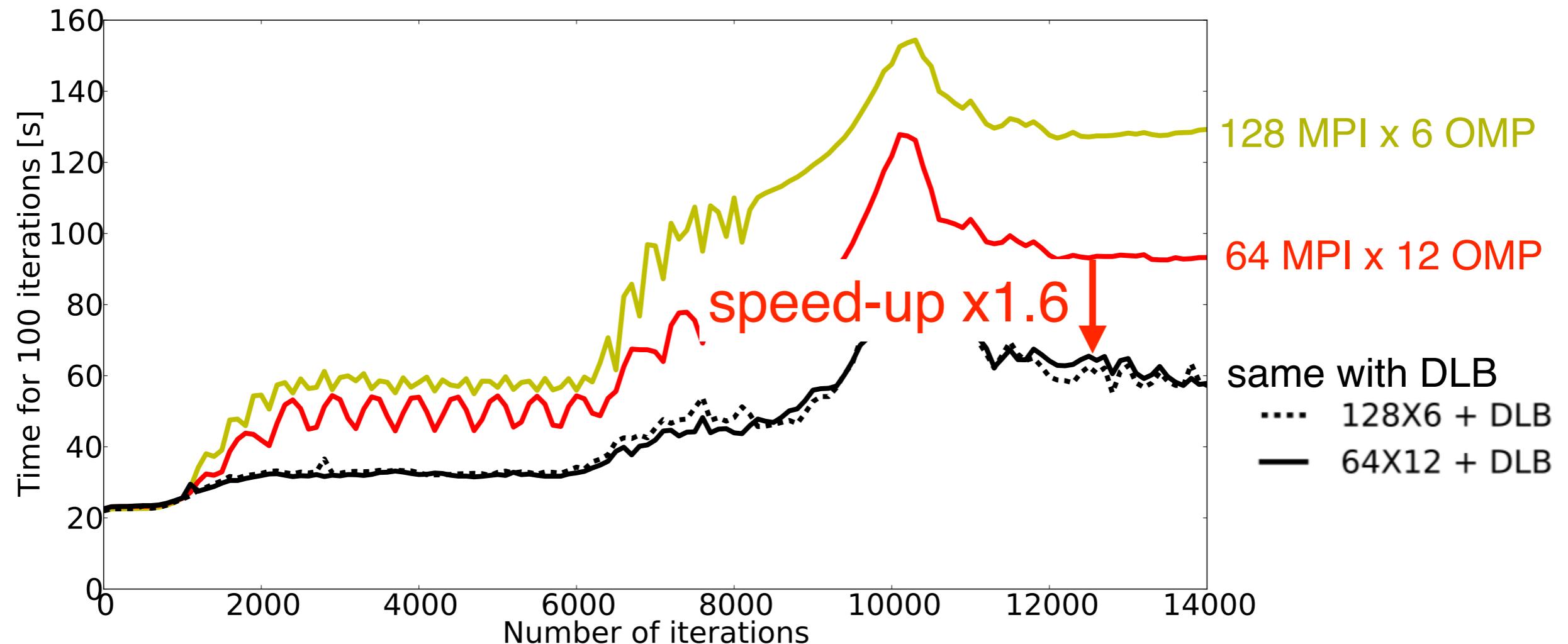
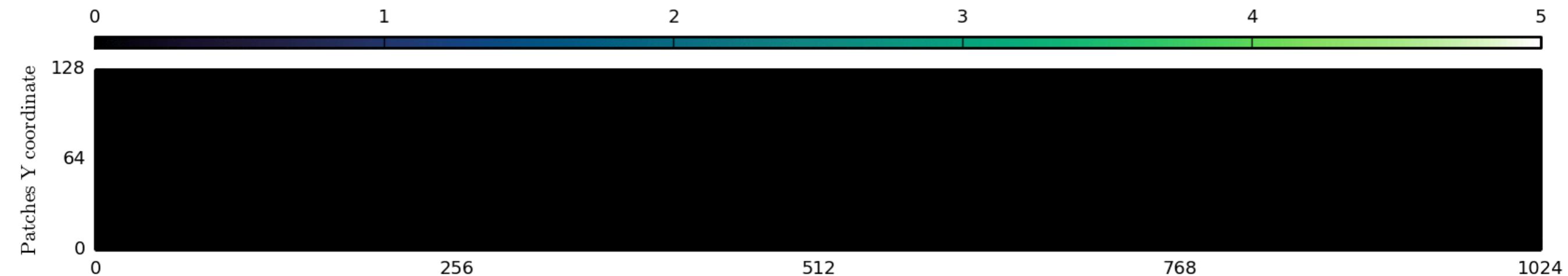
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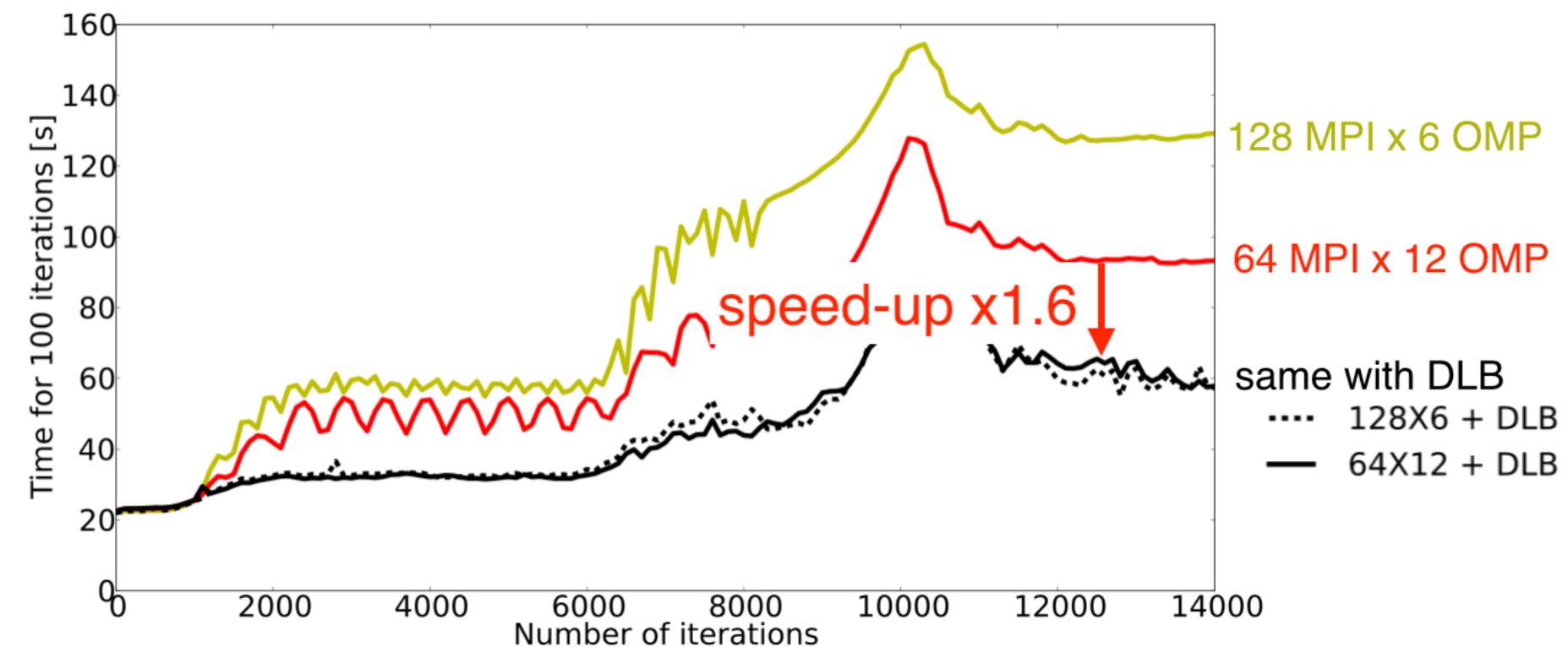
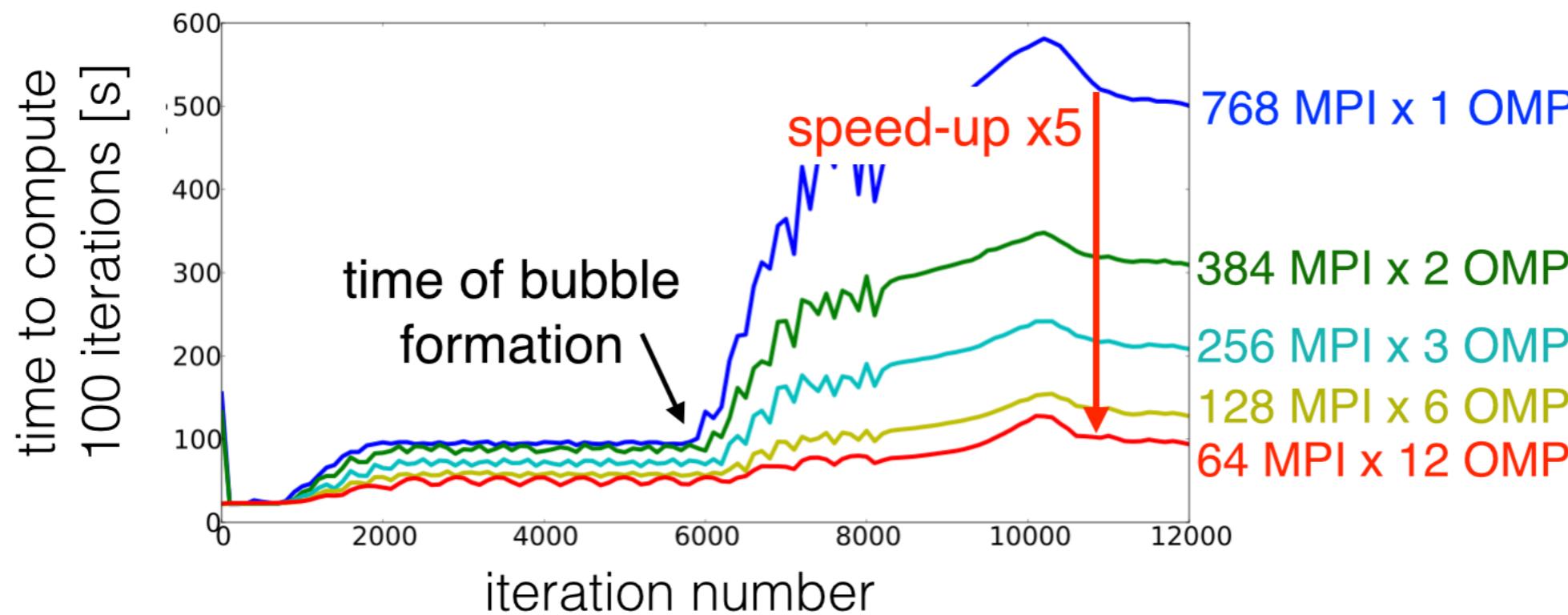
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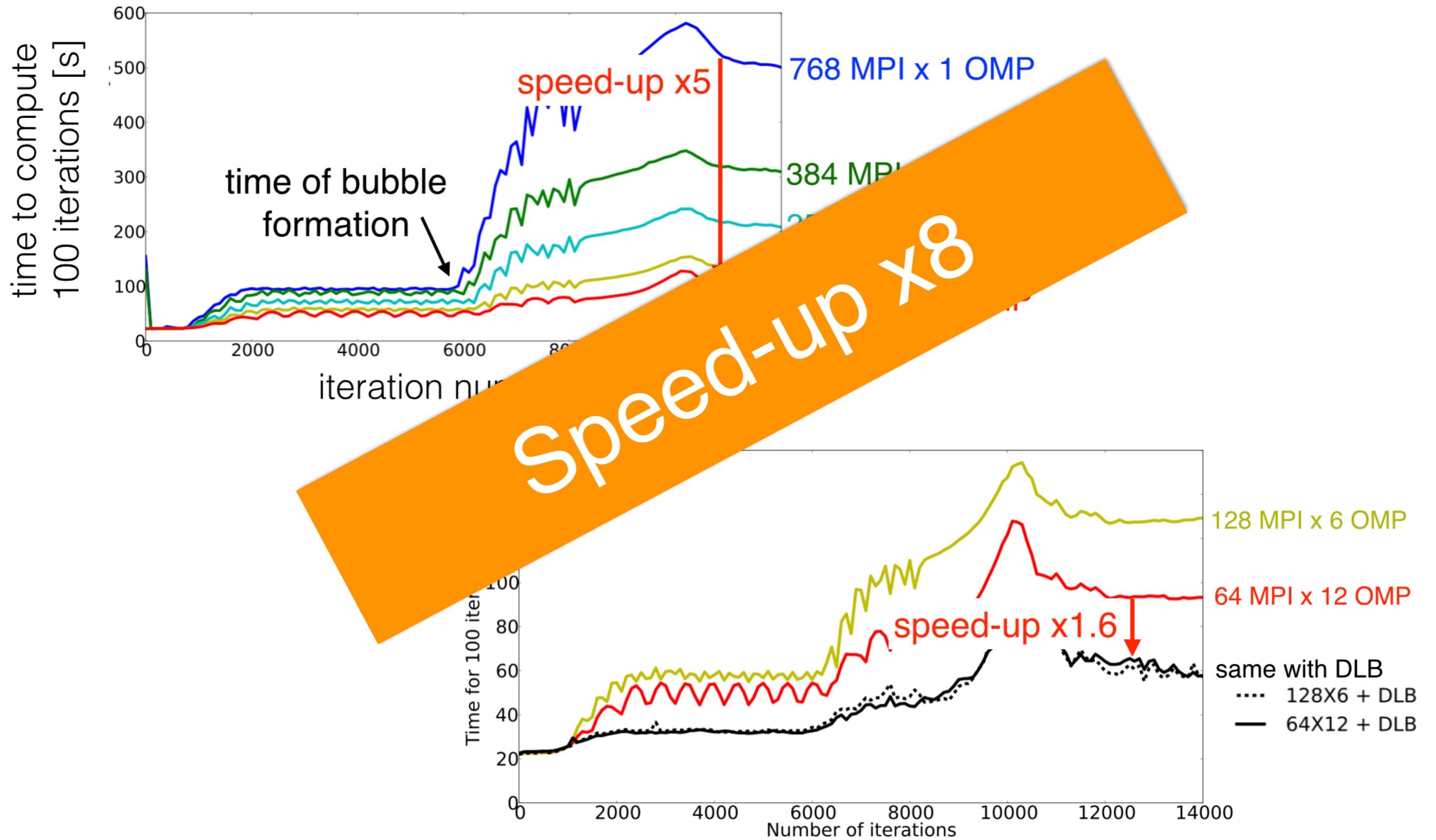


Step 1: Parallelization

Hybrid + Dynamic Load Balancing



Step 1: Parallelization Hybrid + Dynamic Load Balancing



Step 2: Vectorization

Vectorization in a nutshell

Introducing SIMD: Single Instruction, Multiple Data

- Scalar processing

- traditional mode
- one operation produces one result

$$\begin{array}{c} X \\ + \\ Y \\ \hline X + Y \end{array}$$

- SIMD processing

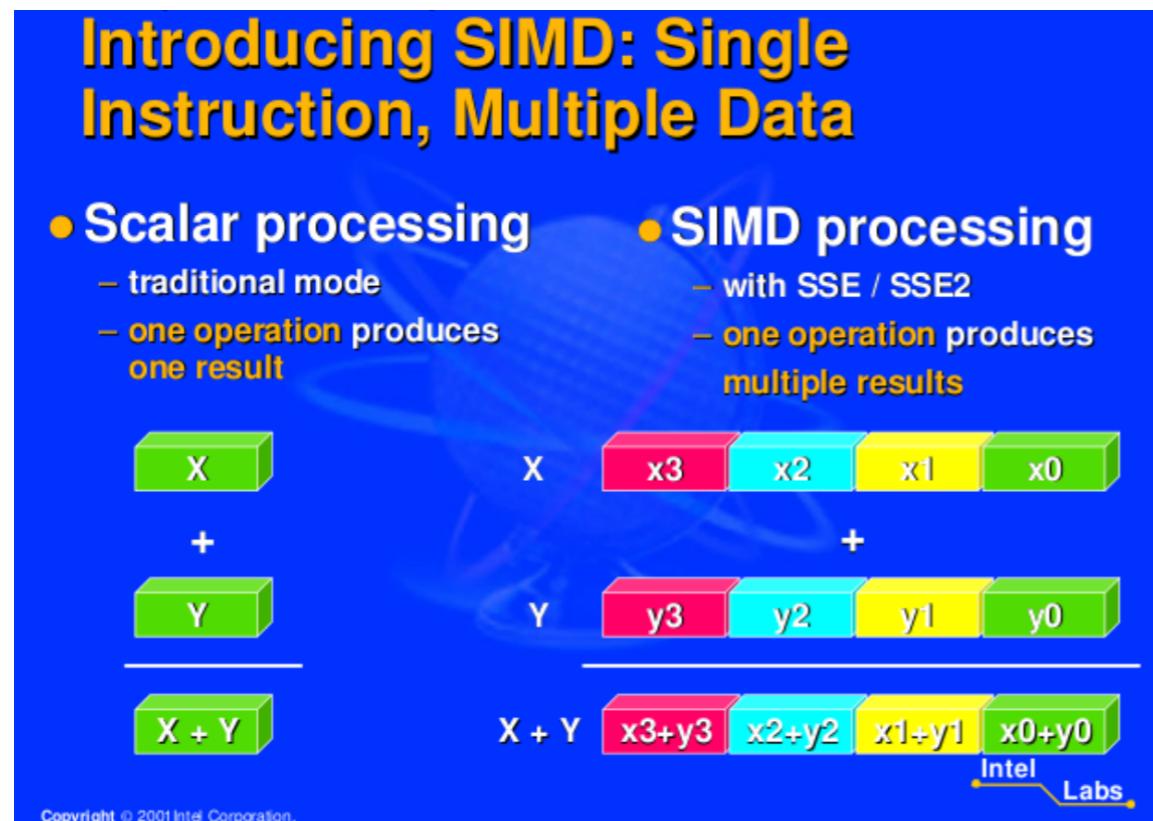
- with SSE / SSE2
- one operation produces multiple results

$$\begin{array}{c} X \quad x3 \quad x2 \quad x1 \quad x0 \\ + \\ Y \quad y3 \quad y2 \quad y1 \quad y0 \\ \hline X + Y \quad x3+y3 \quad x2+y2 \quad x1+y1 \quad x0+y0 \end{array}$$

Intel
Labs

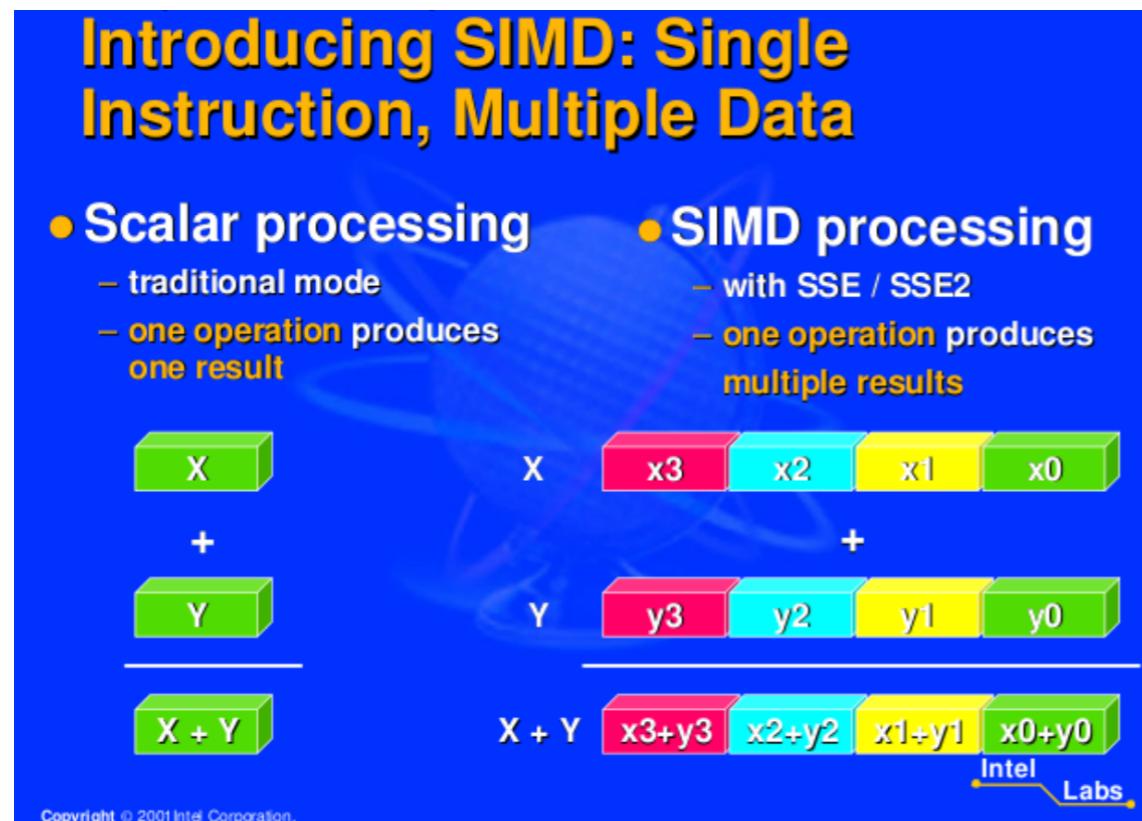
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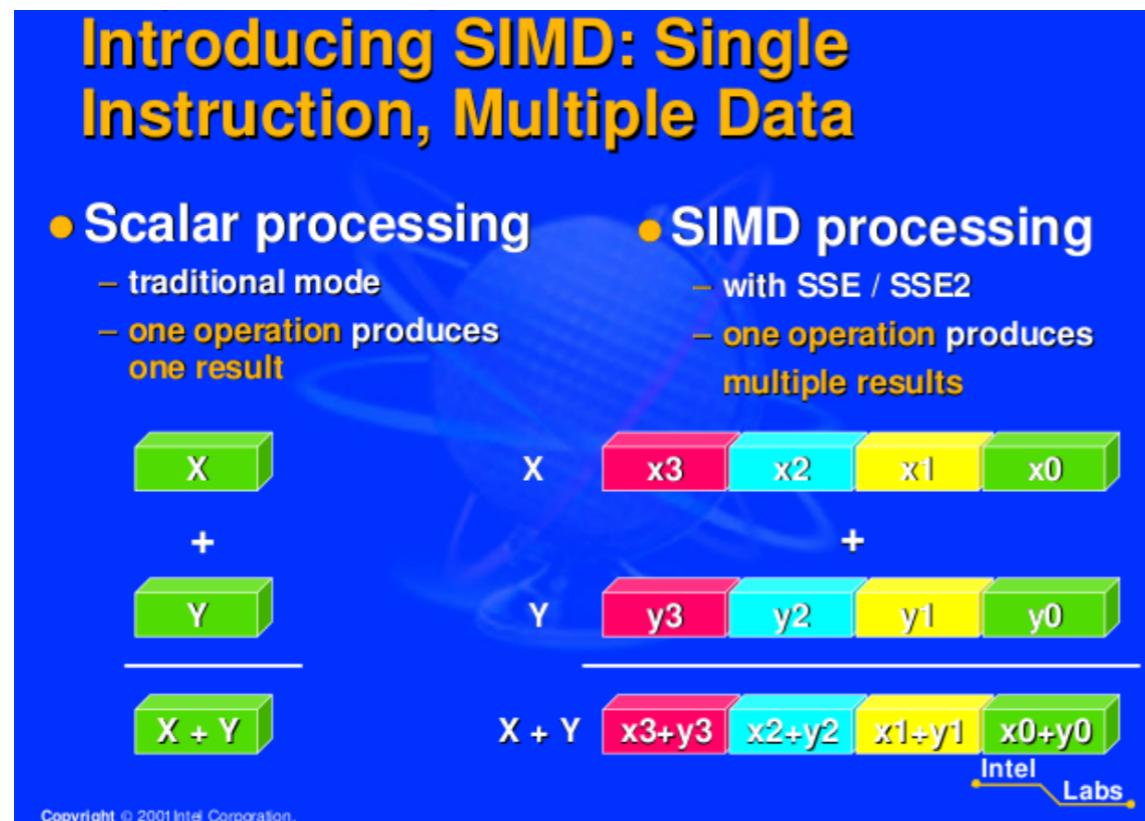
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Smart (particles) operators:
- interpolator, pusher, projector

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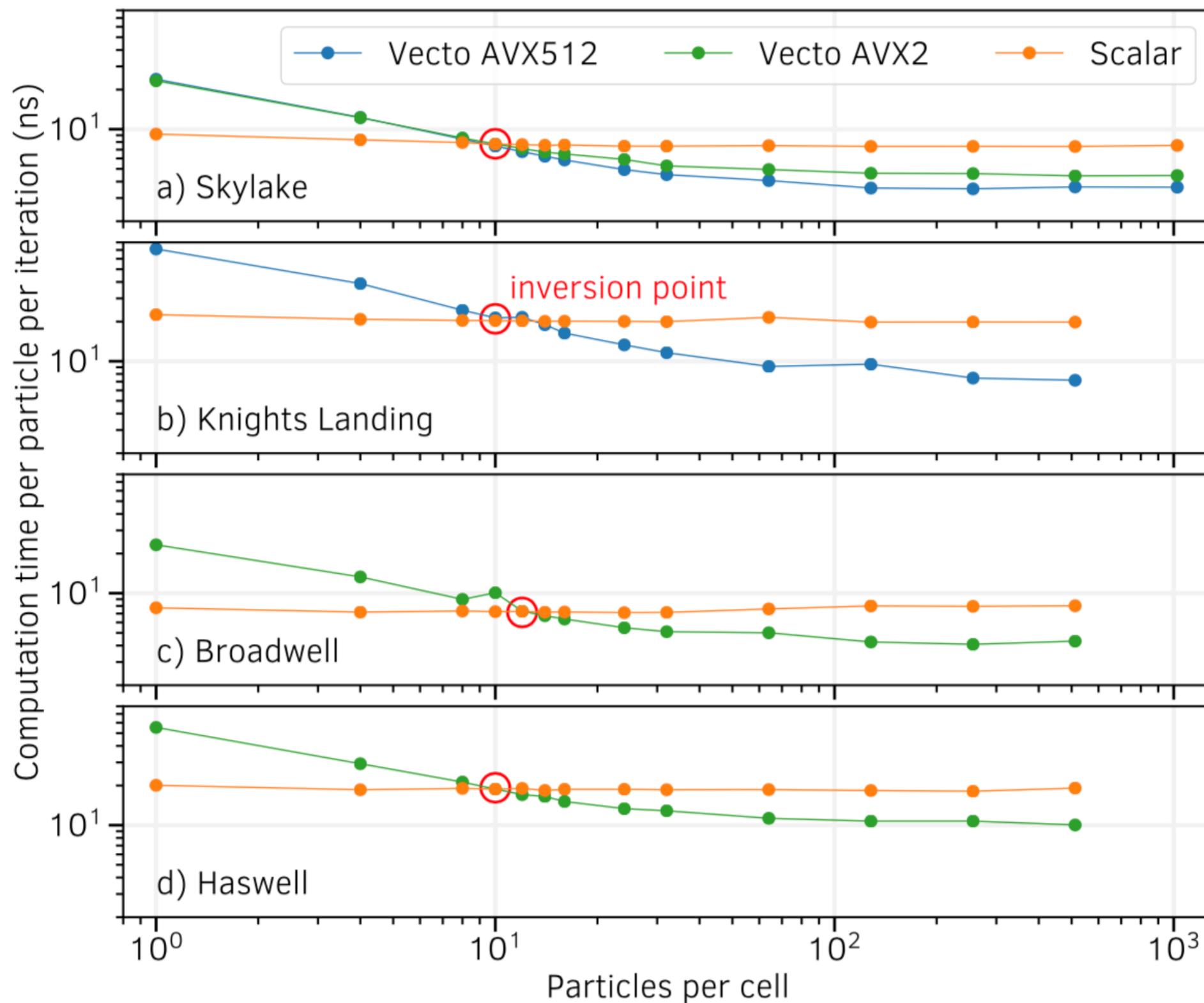
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- Smart (particles) operators:**
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- Smart (particles) data structures:**
- beware random mem. access
 - contiguous memory
 - sort at all times!

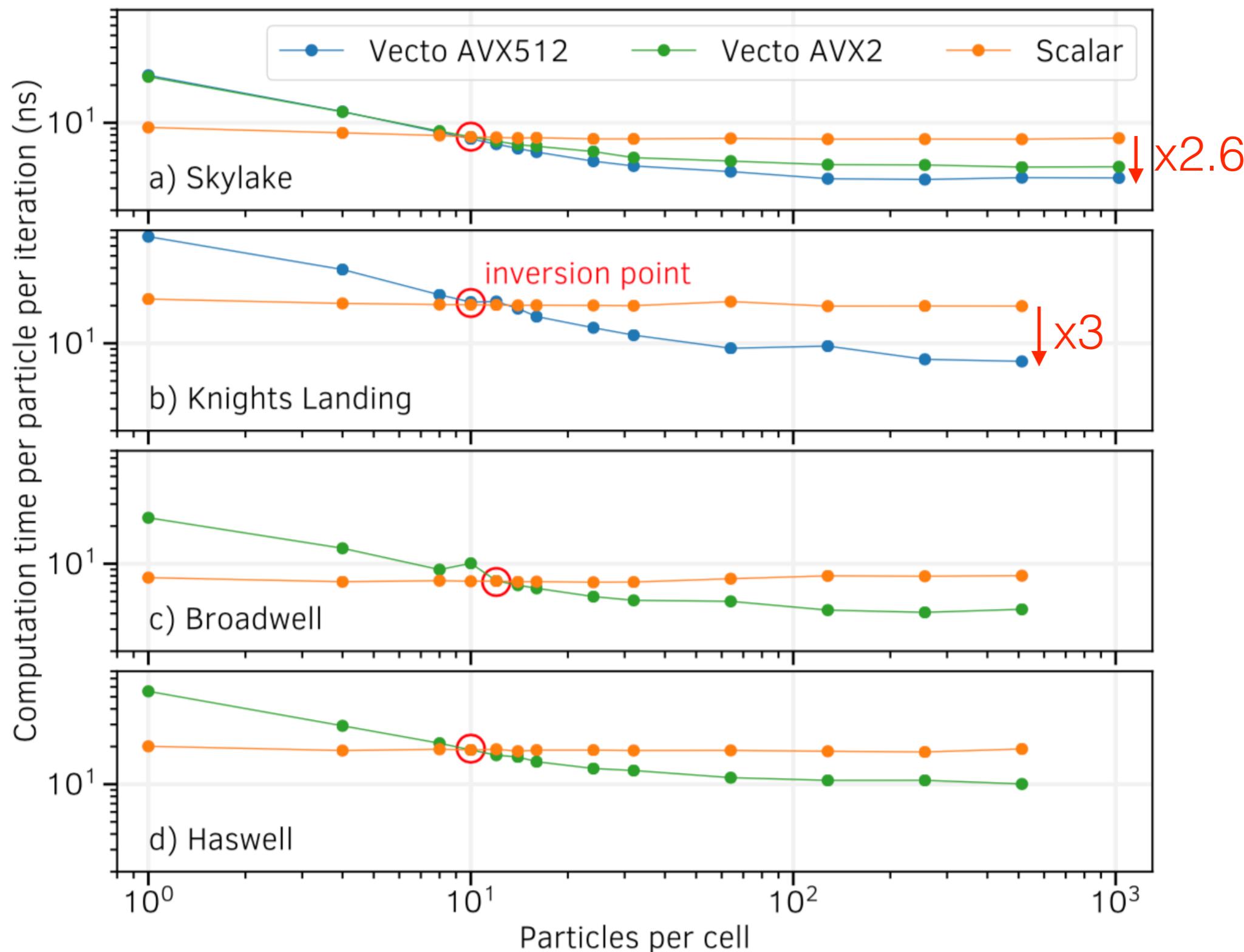
Step 2: Vectorization

SMILEI uses an adaptive vectorization approach



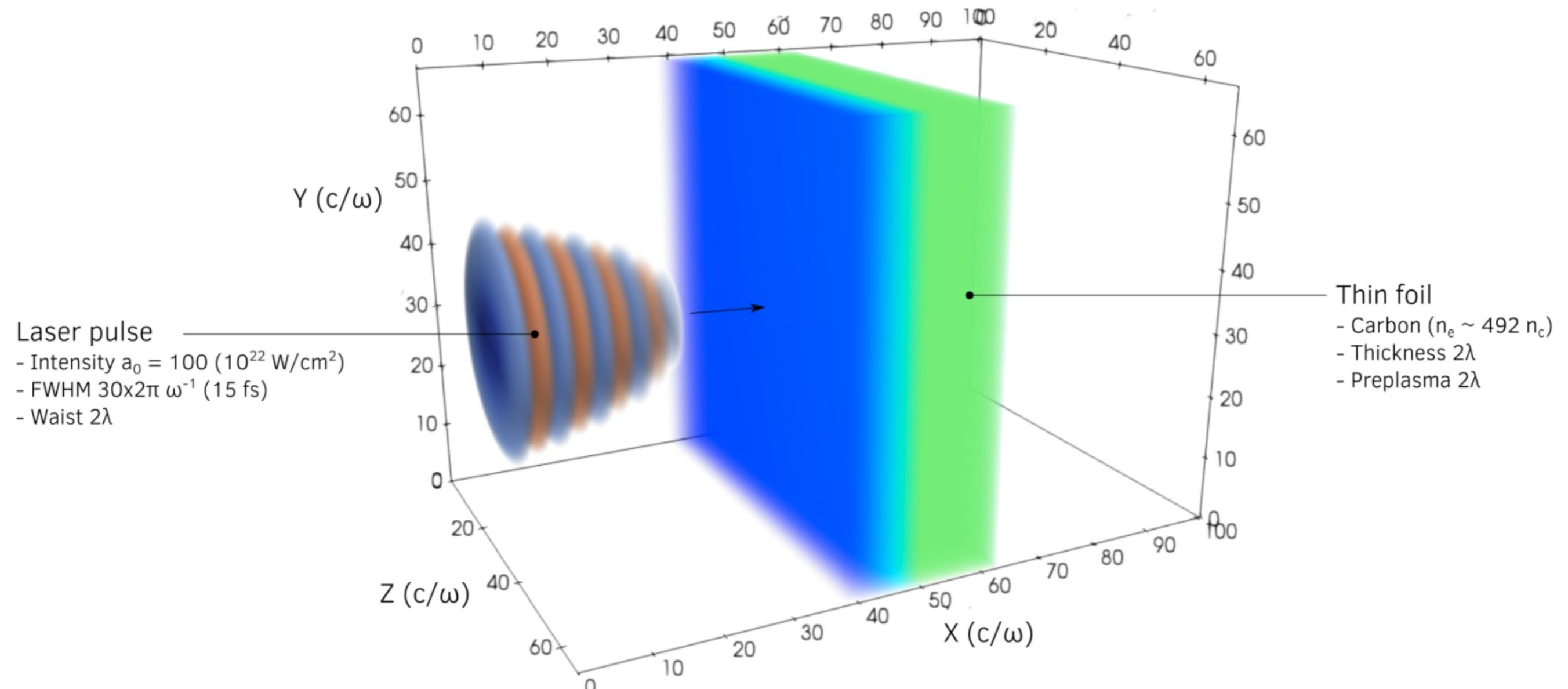
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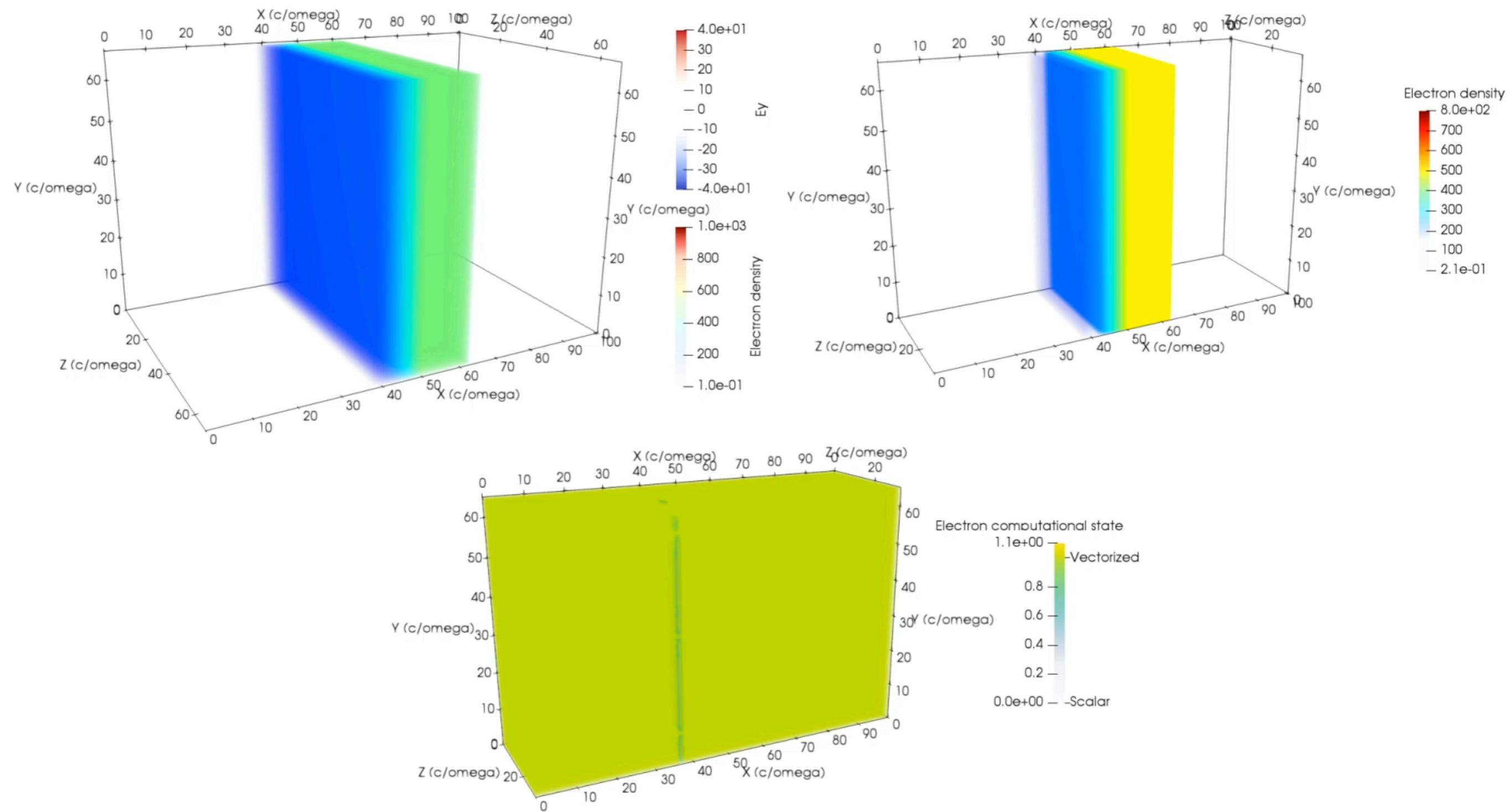
Laser-driven hole-boring



@ 32 PPC : speed-up x 1.5

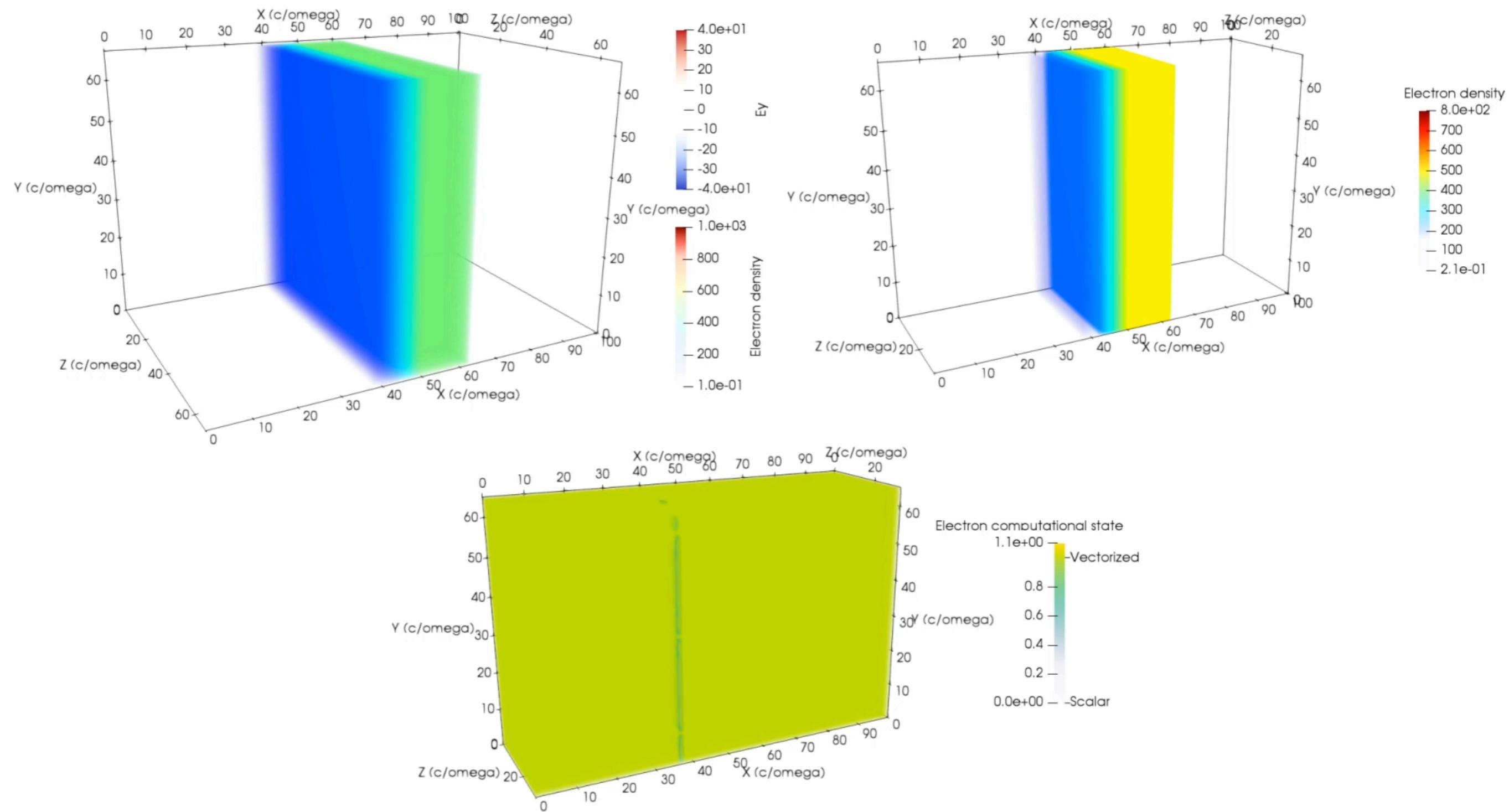
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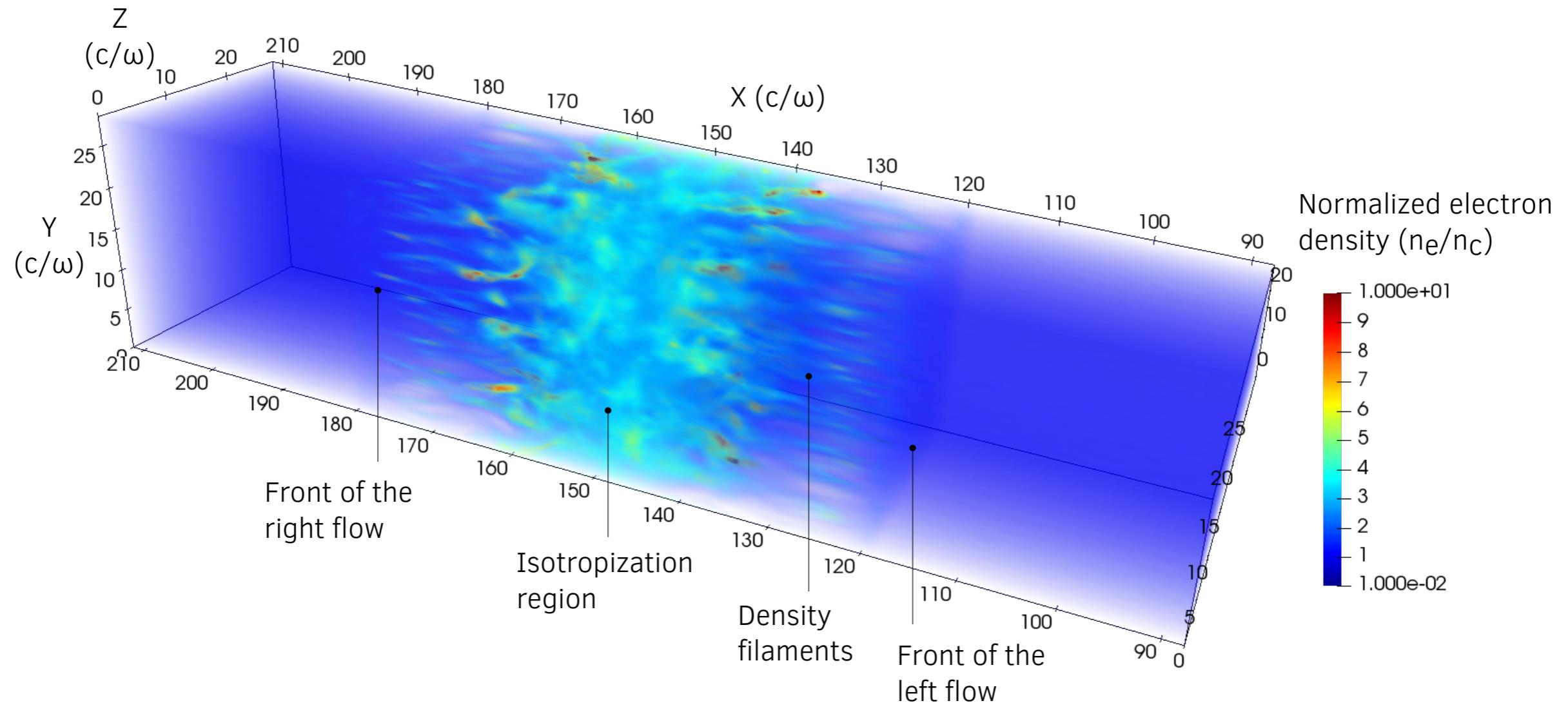
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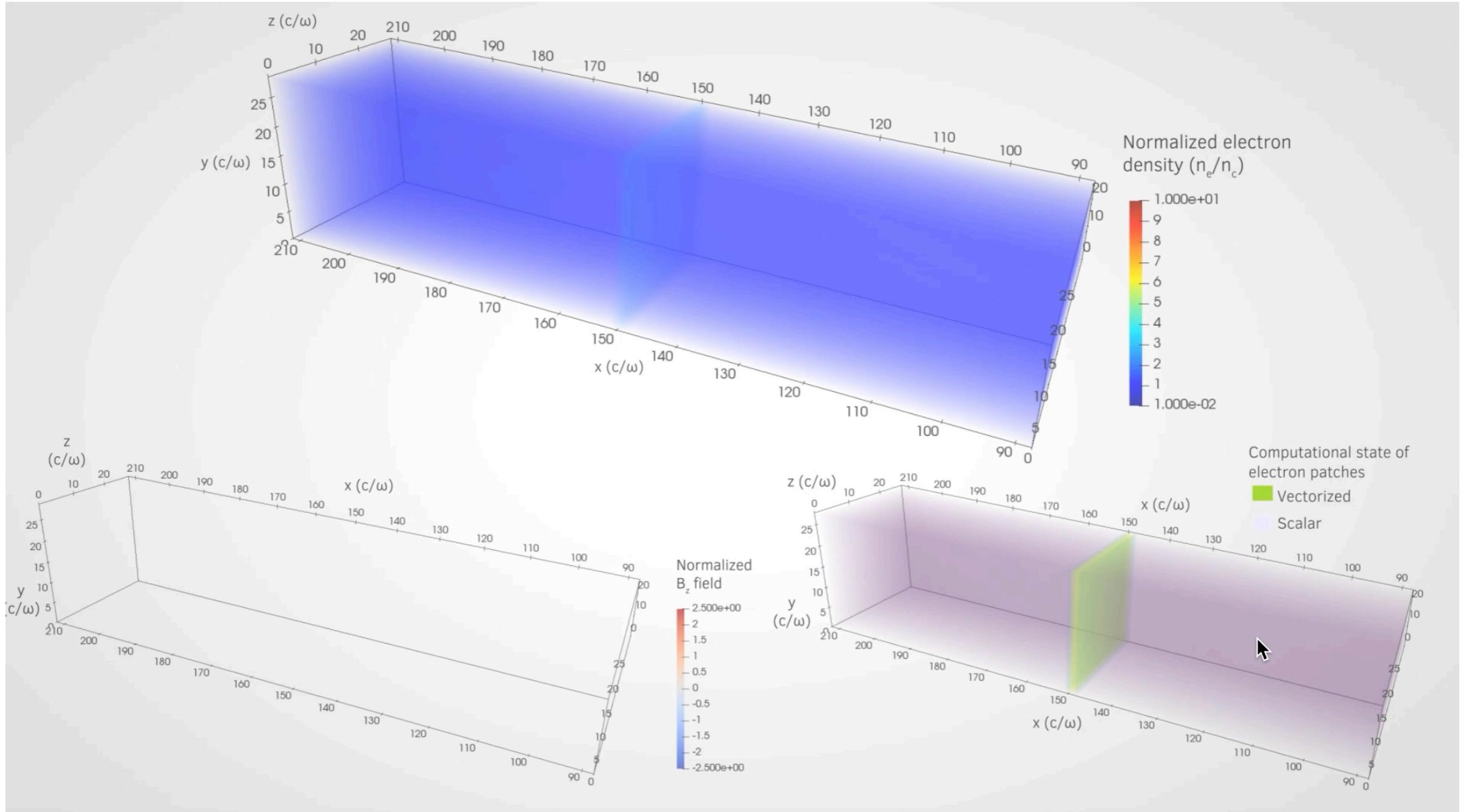
Weibel-mediated collisionless shocks



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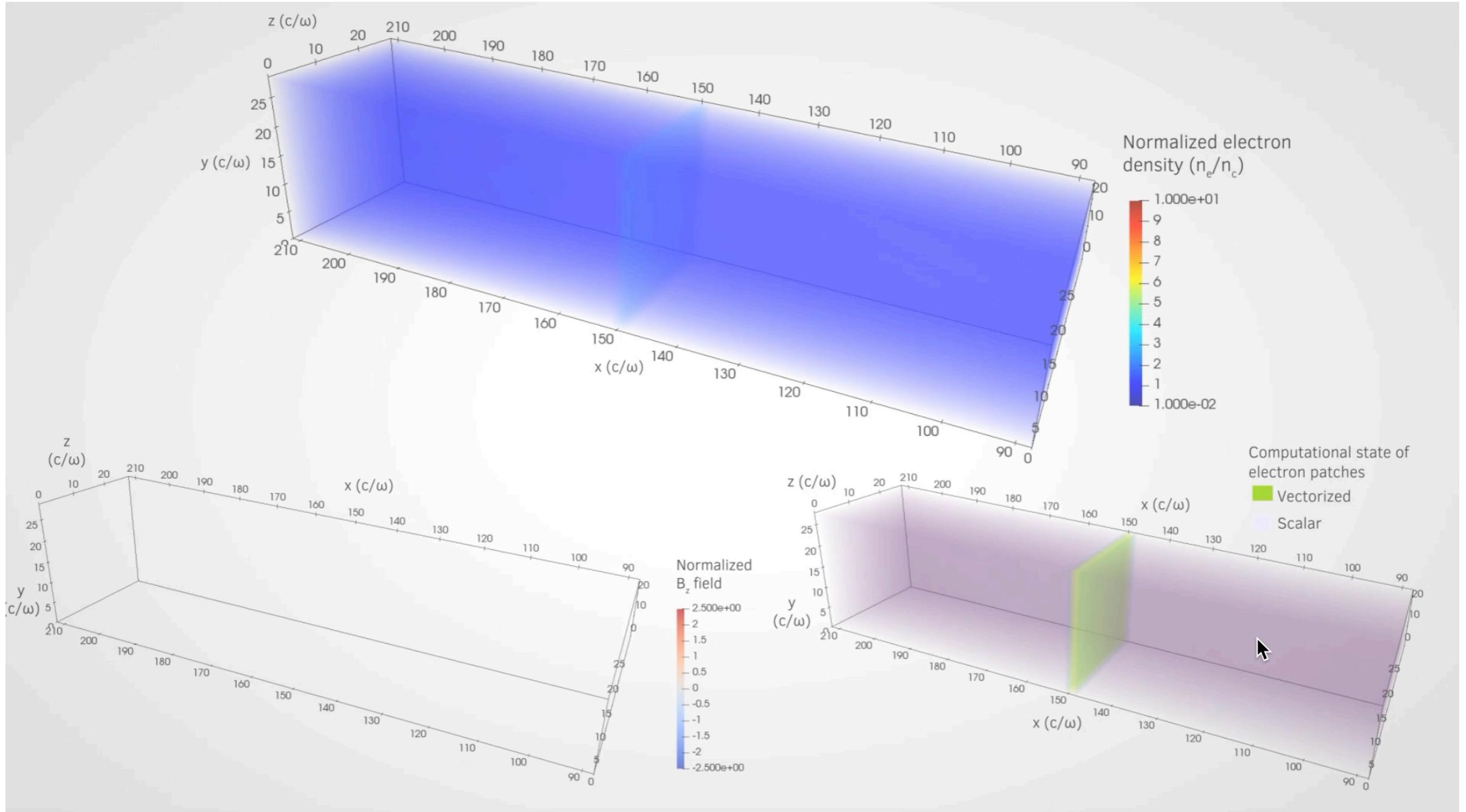
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Collisions can be introduced using an *ad-hoc* Monte-Carlo module

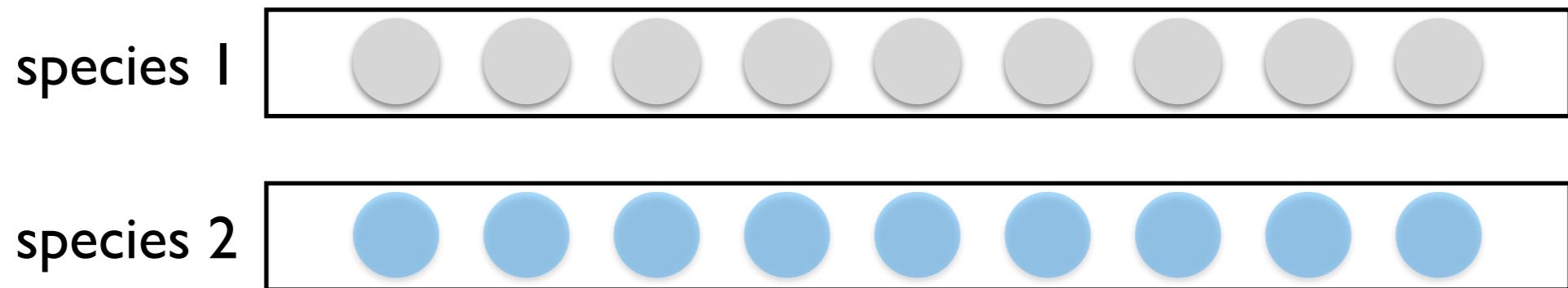
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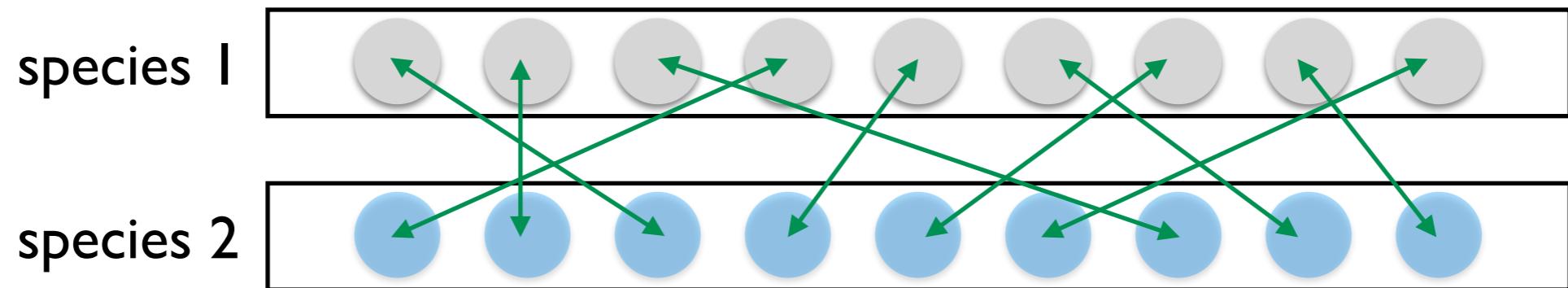
To avoid the N-body problem, *quasi-particles in the cell are randomly “paired”*



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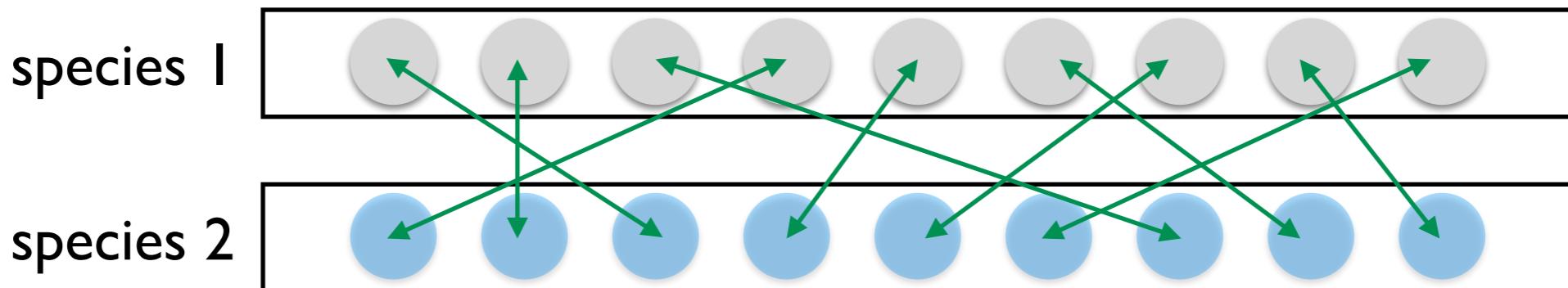
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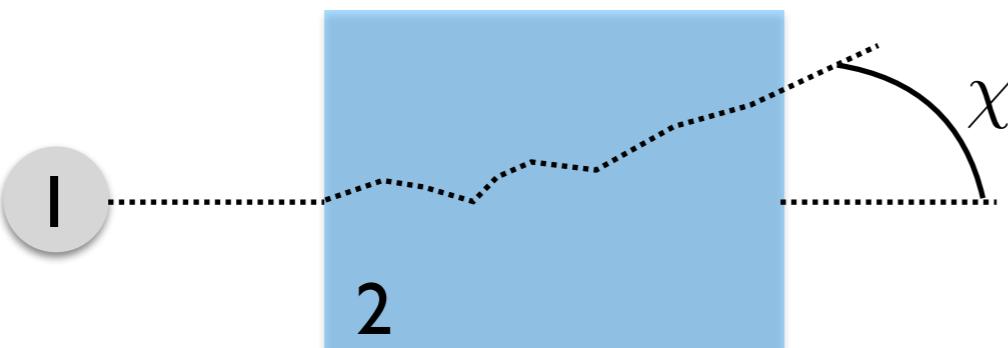
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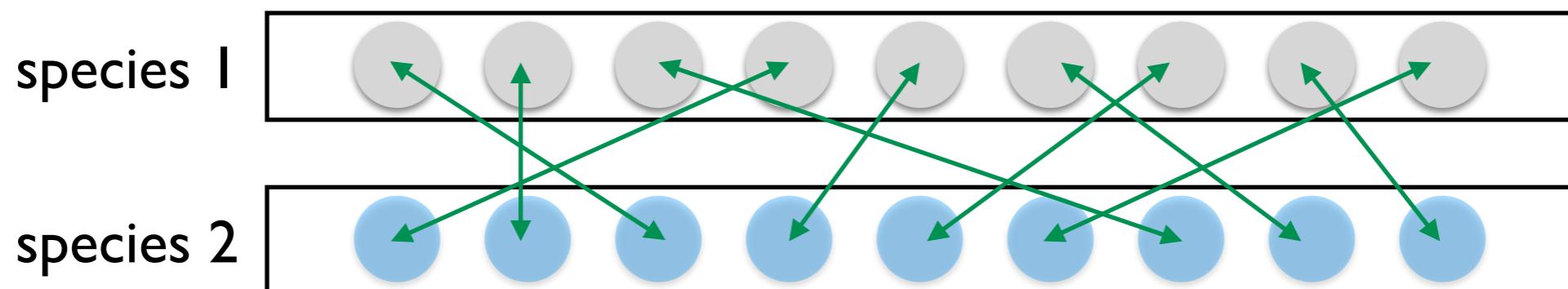
A single particle goes through many ($N \gg 1$) collisions at small angle θ which translates in a total deflection angle χ (not necessarily small)



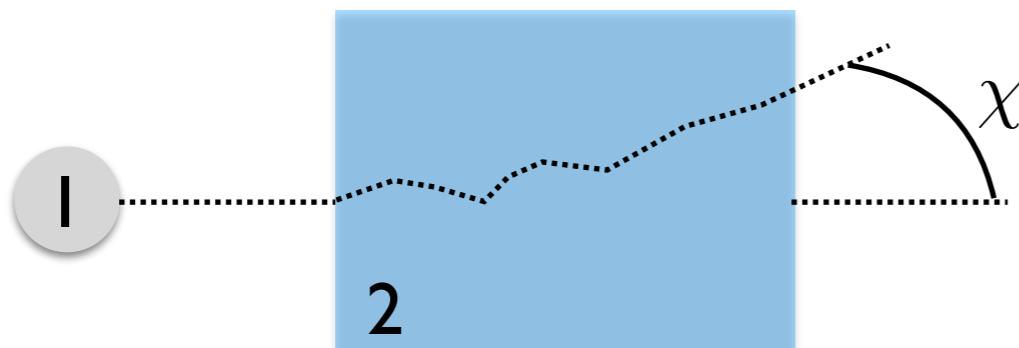
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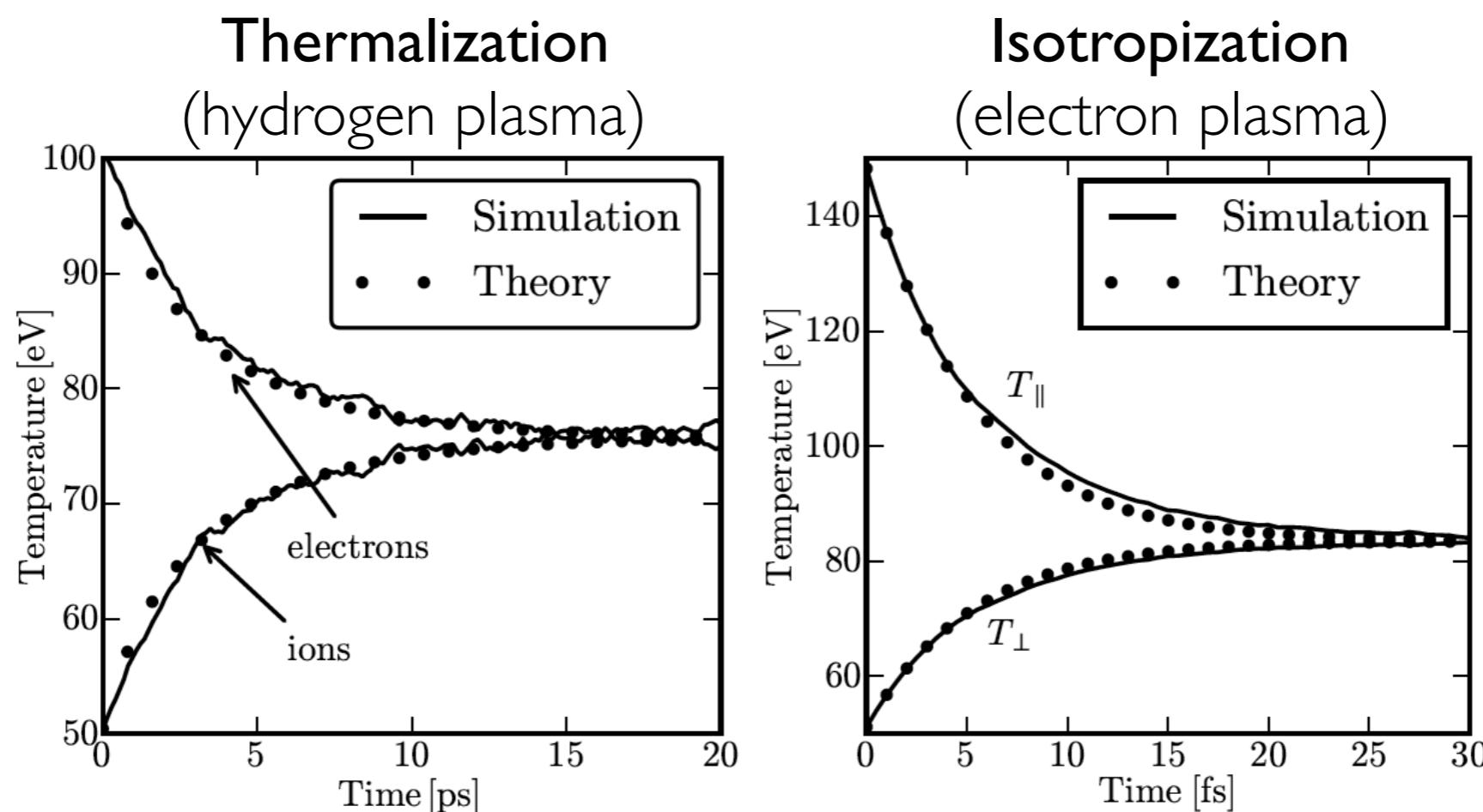


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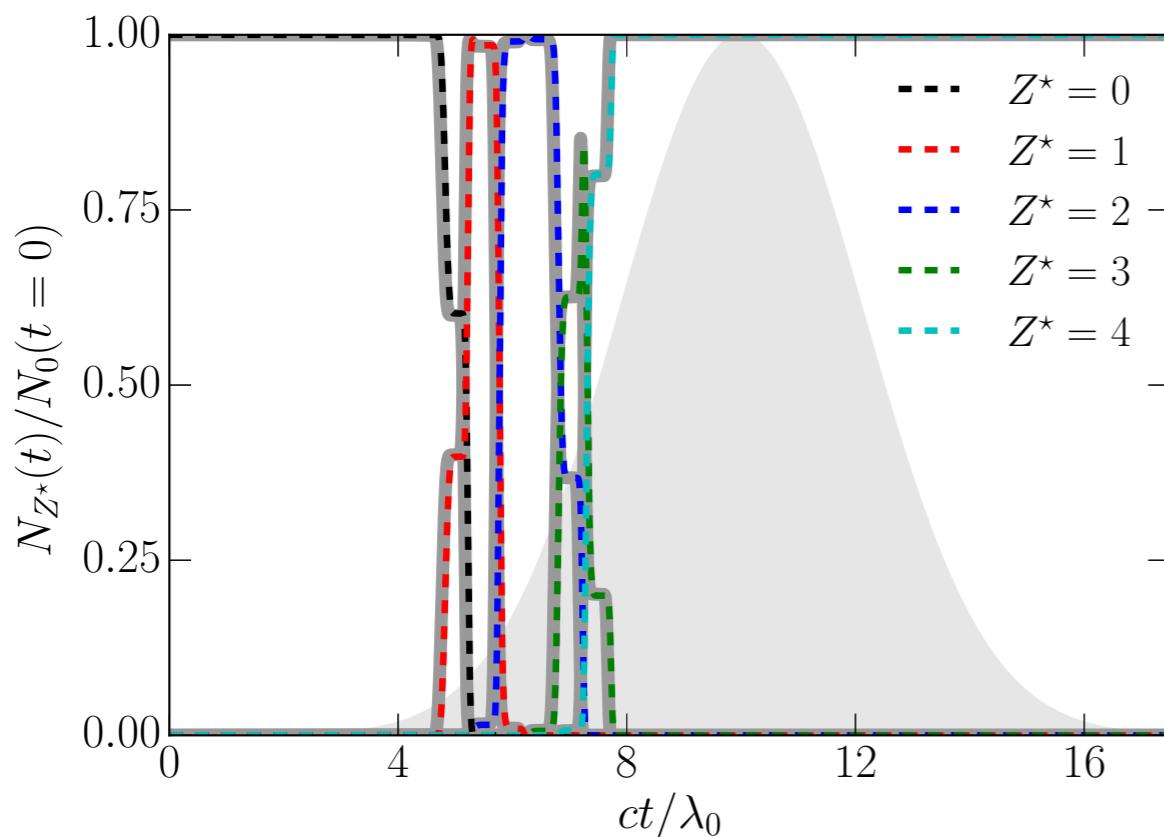
- for each pair (Monte-Carlo)**
- compute the collision rate
 - compute the deflection angle
 - deflect one or both particles

PIC codes are then able to treat purely **collisional processes**

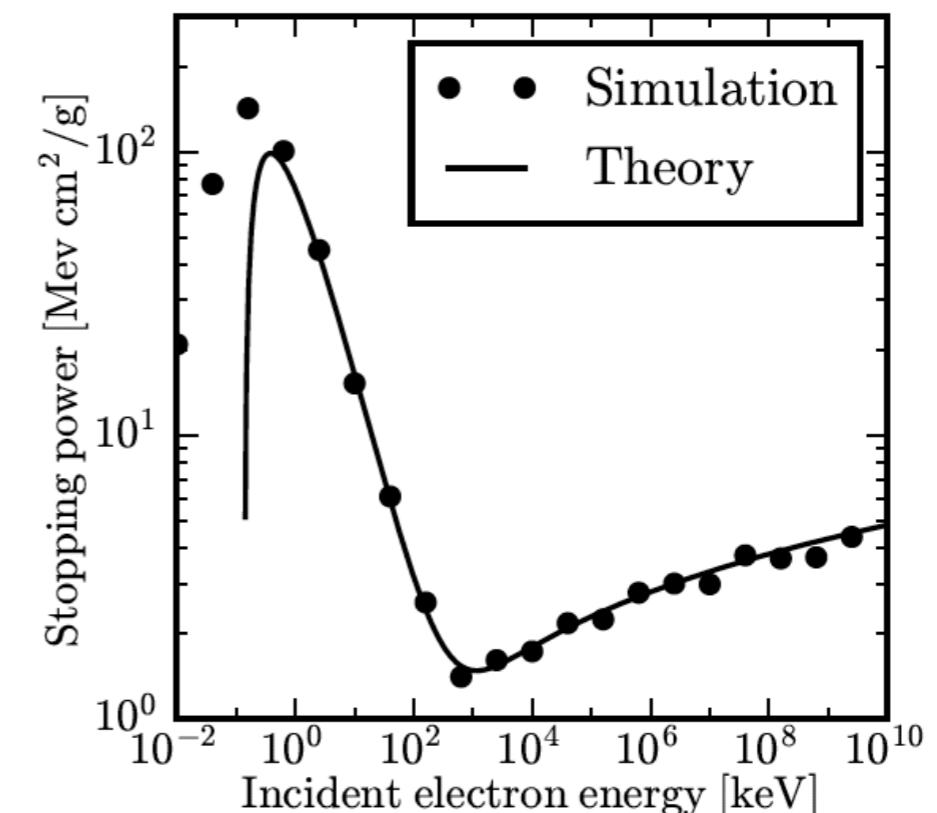


Similarly field and collisional ionization can be treated using a Monte-Carlo approach

Field ionization of Carbon by a $5 \times 10^{16} \text{ W/cm}^2$ 20 fs laser pulse



Stopping power of a cold aluminium plasma of density 10^{21} cm^{-3}



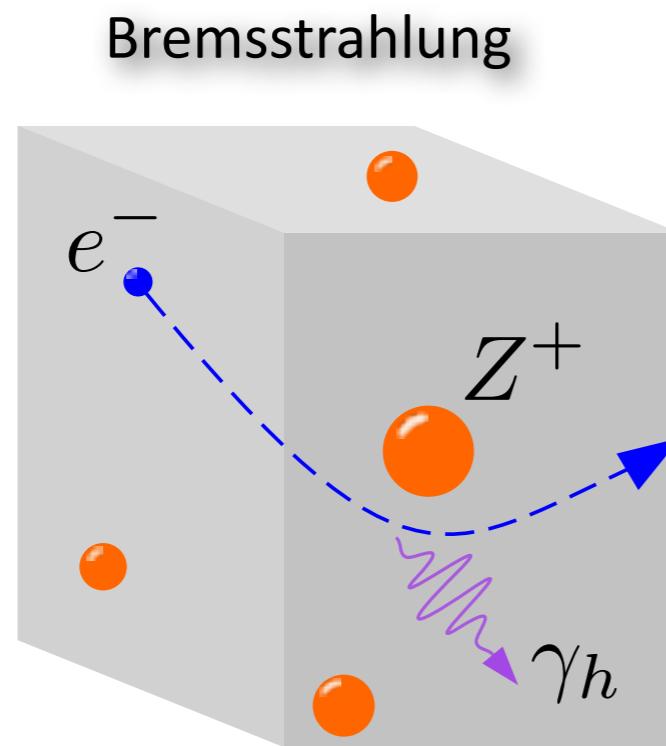
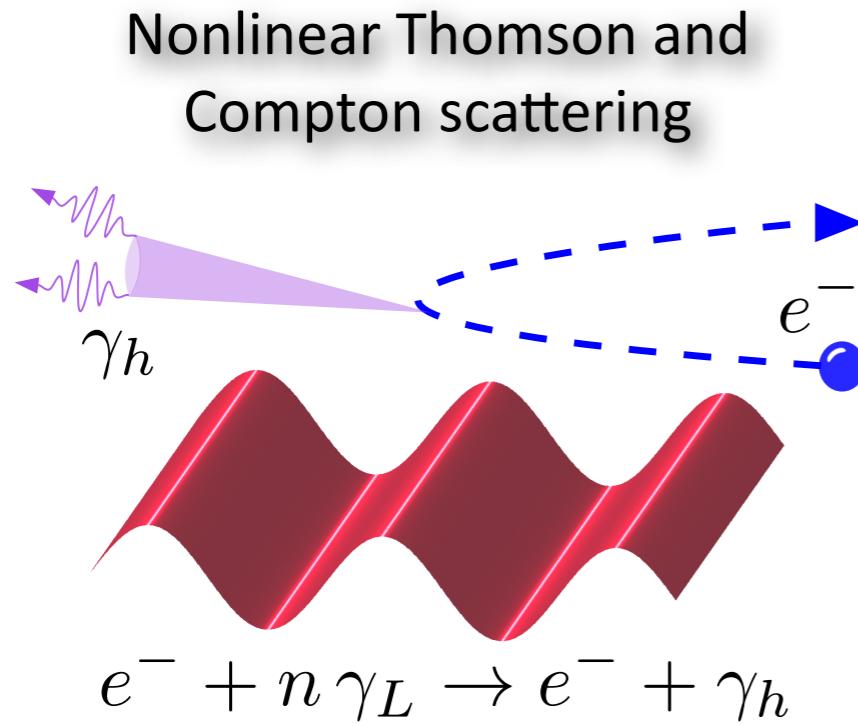
R. Nuter *et al.*, Phys. Plasmas **18**, 033107 (2011); F. Pérez *et al.*, Phys. Plasmas **19**, 083104 (2012)

J. Derouillat *et al.*, *SMILEI: a collaborative, open-source, multi-purpose PIC code for plasma simulation*, to be submitted (available upon request)

Adding Quantum Electrodynamics (QED) effect is also very interesting for forthcoming multi-petawatt facilities

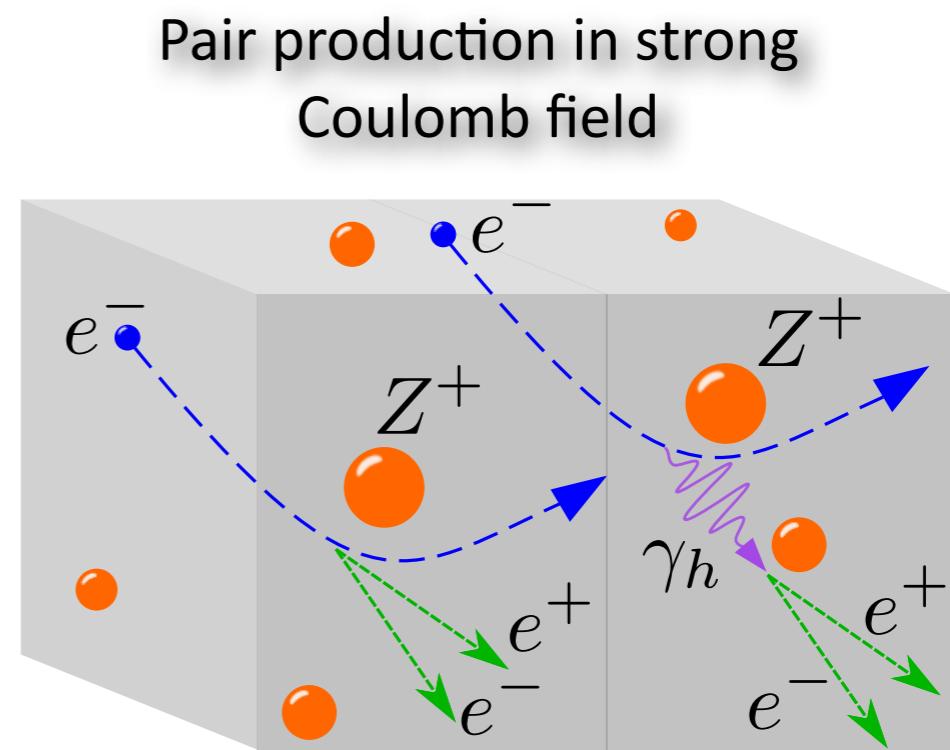
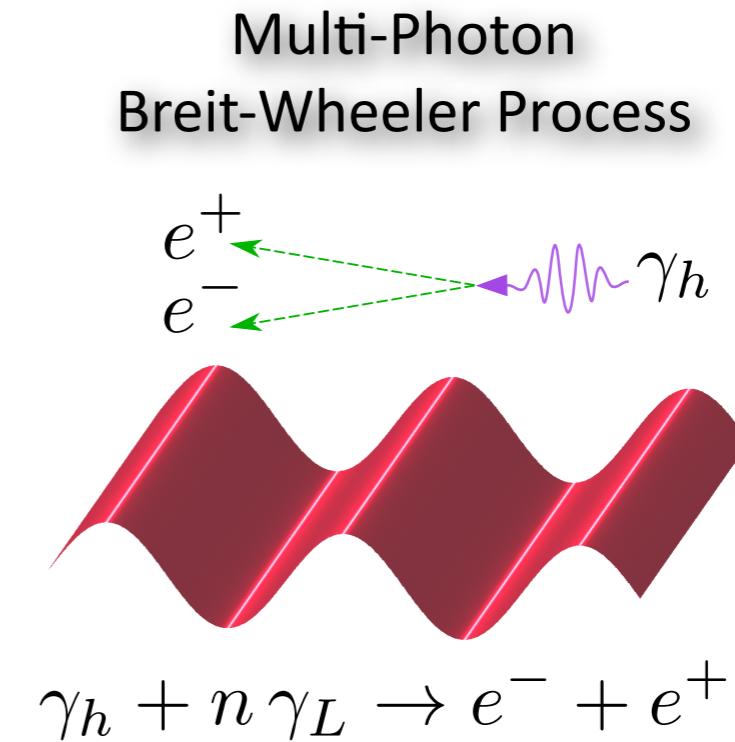
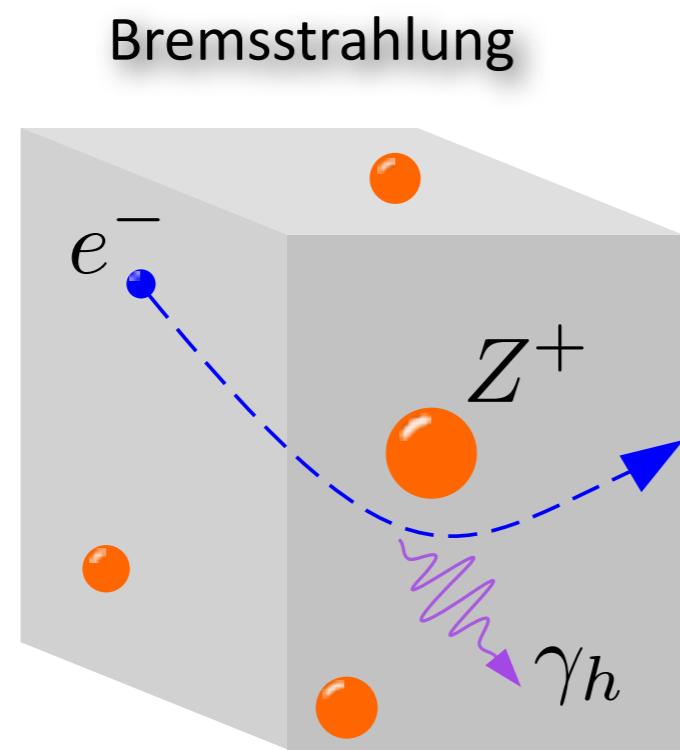
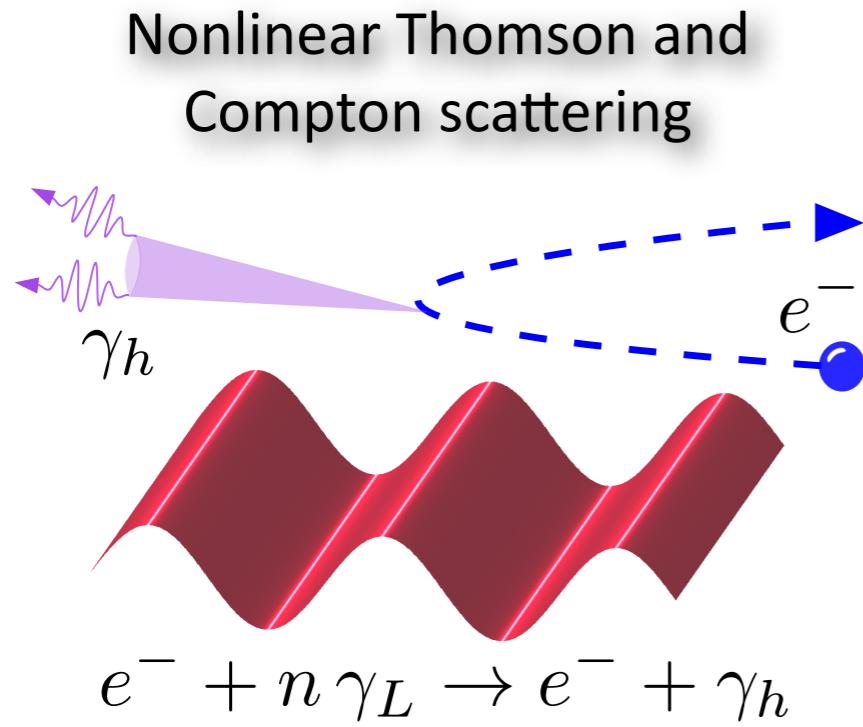
High-Energy Photon Production

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Electron-Positron Pair Production

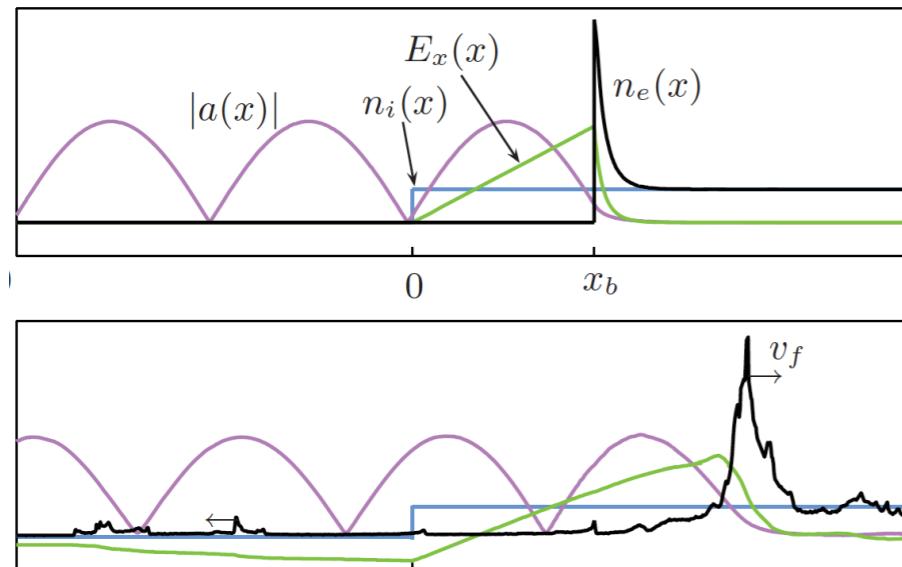
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PIC codes are an excellent tool to support theoretical modelling
Even 1D simulation can bring a deep insight into the physics at play

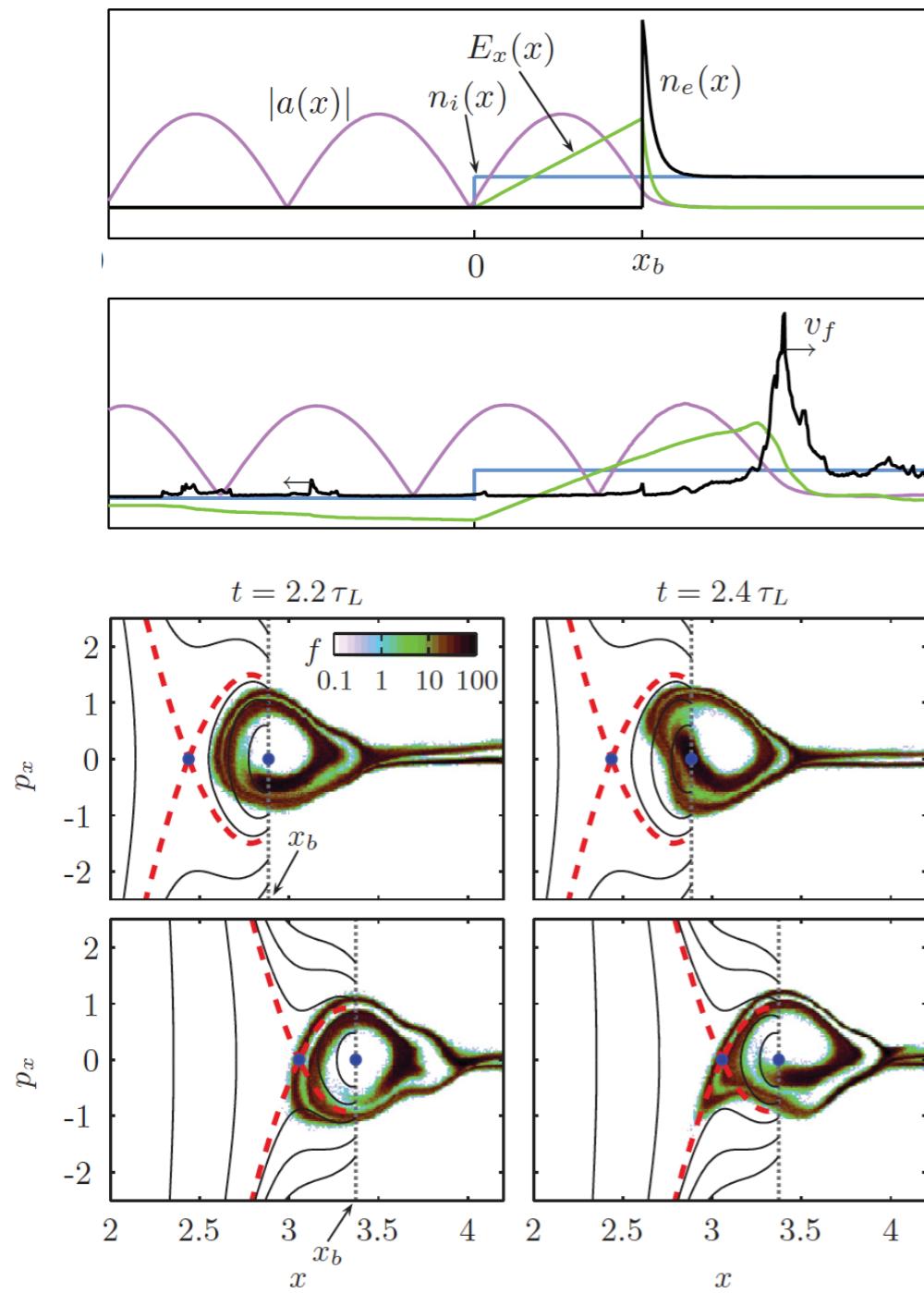
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Relativistically-Induced Transparency



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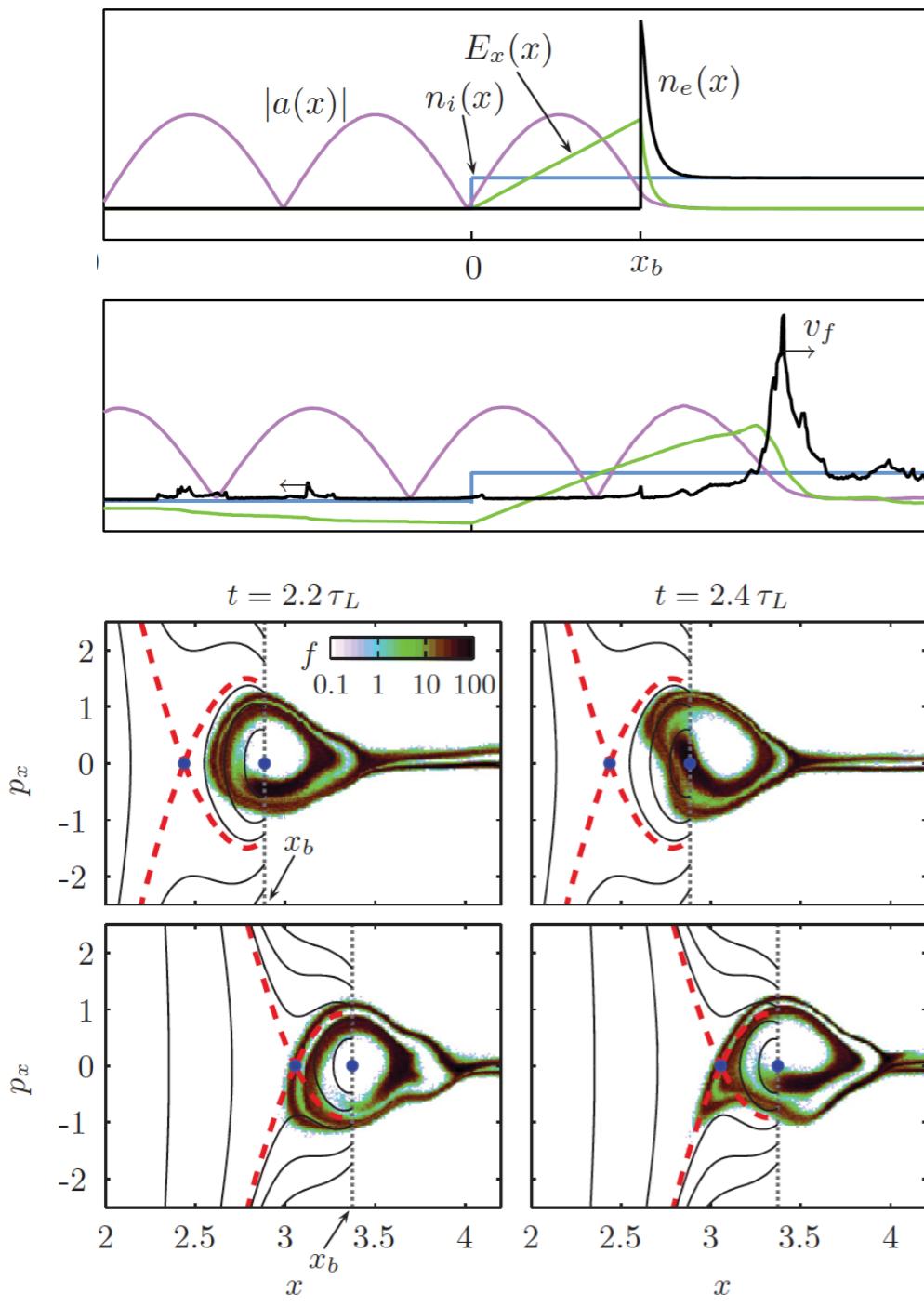
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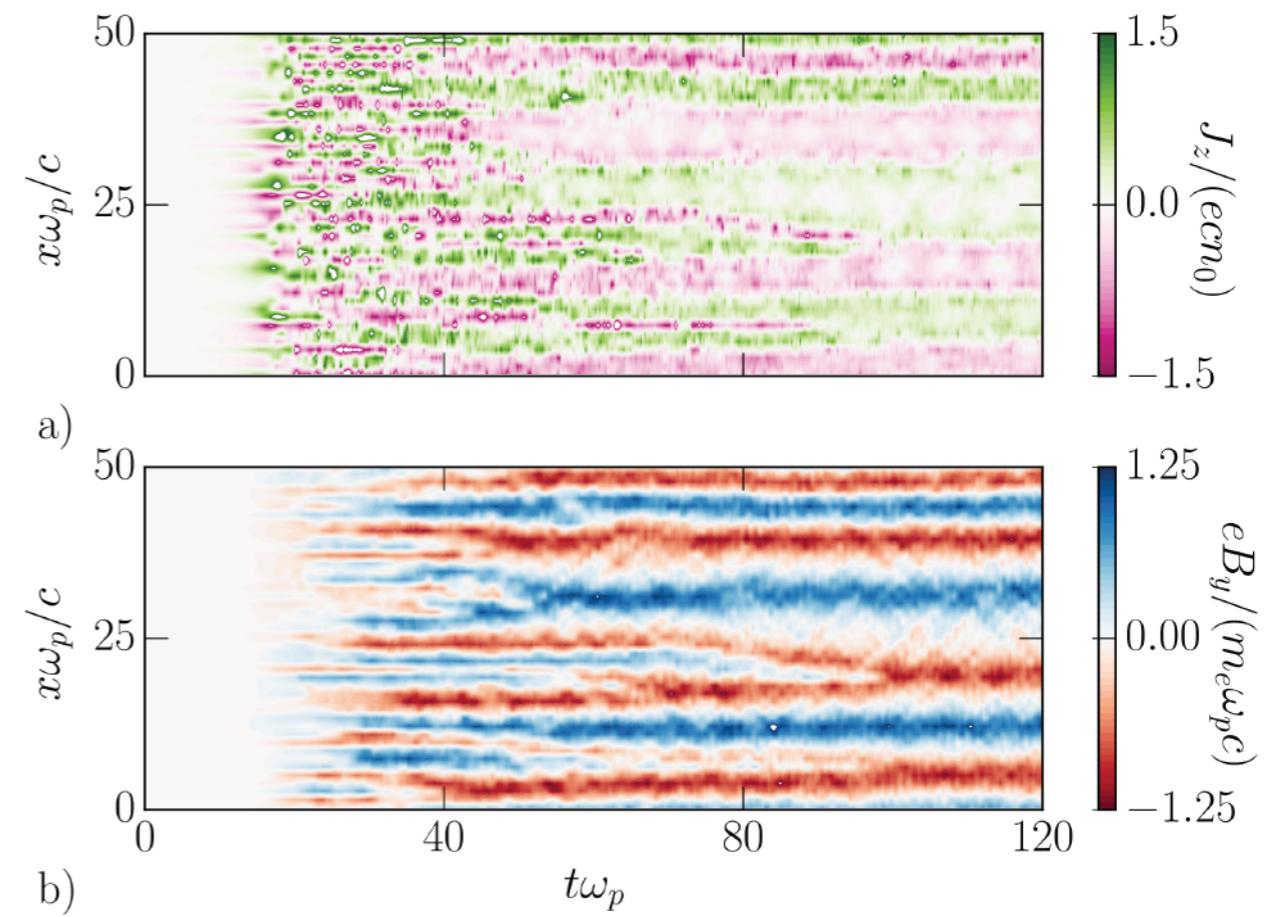
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E. Siminos *et al.*, Phys. Rev. E **86**, 056404 (2012)

Weibel instability in the presence of an external magnetic field



A. Grassi *et al.*, Phys. Rev. E (in press)

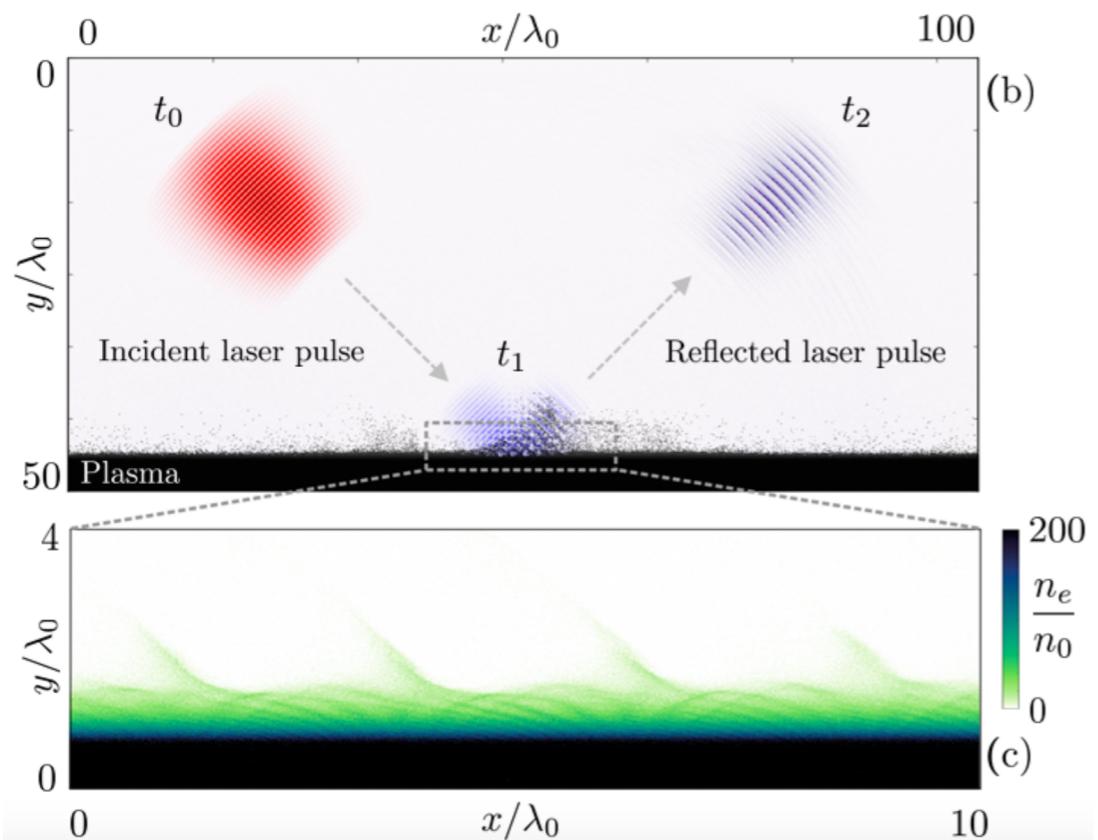
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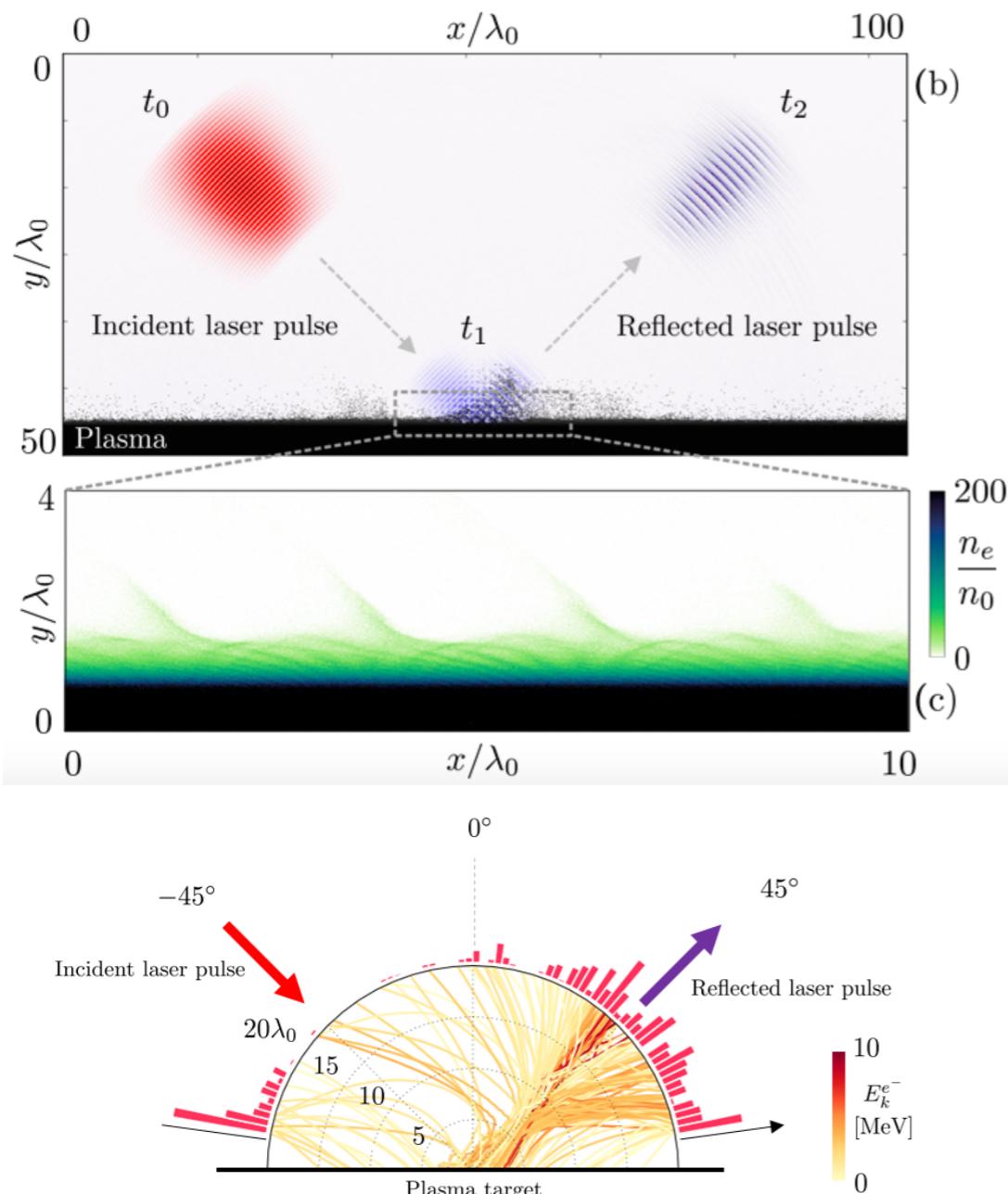
**High-harmonic generation
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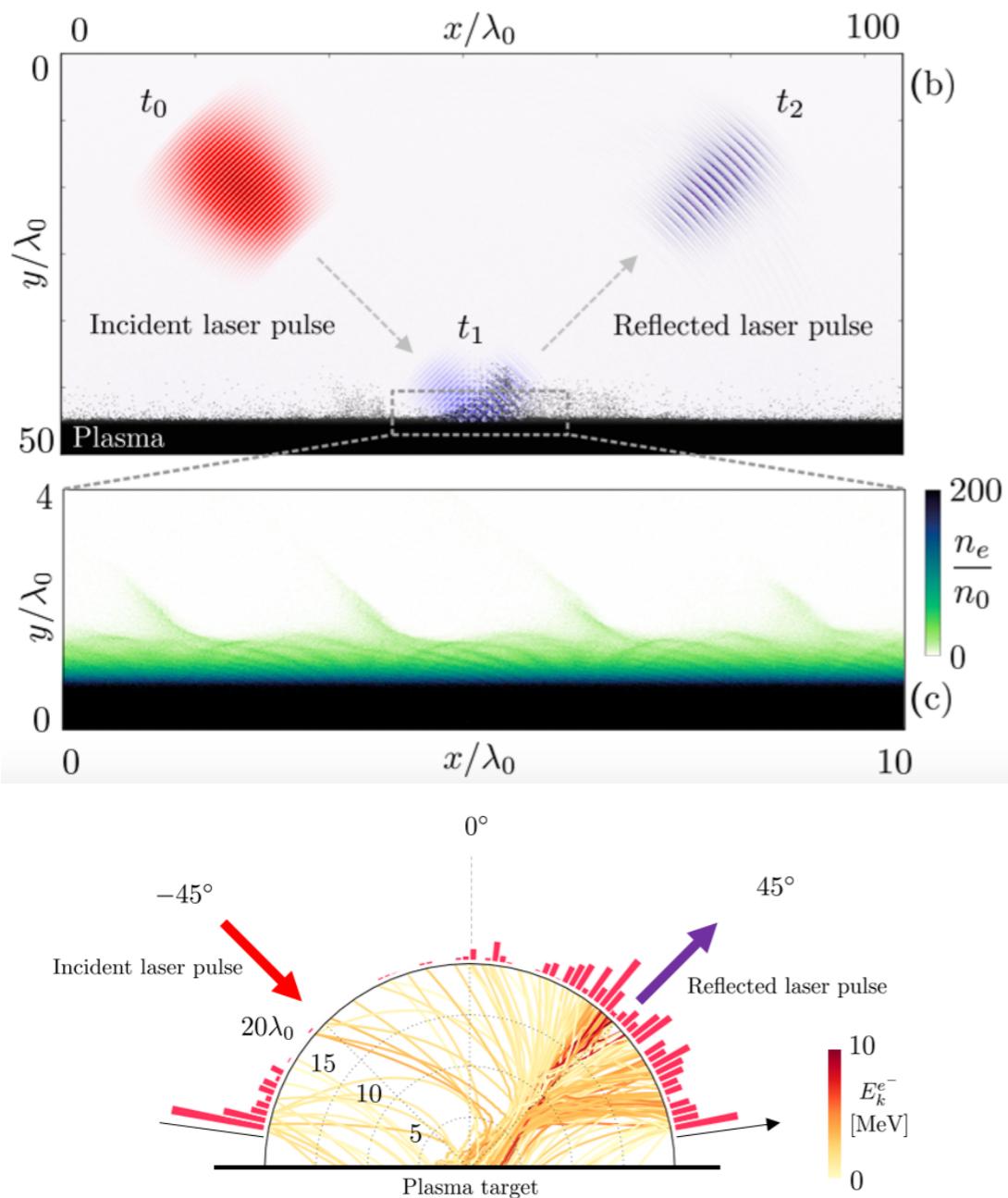
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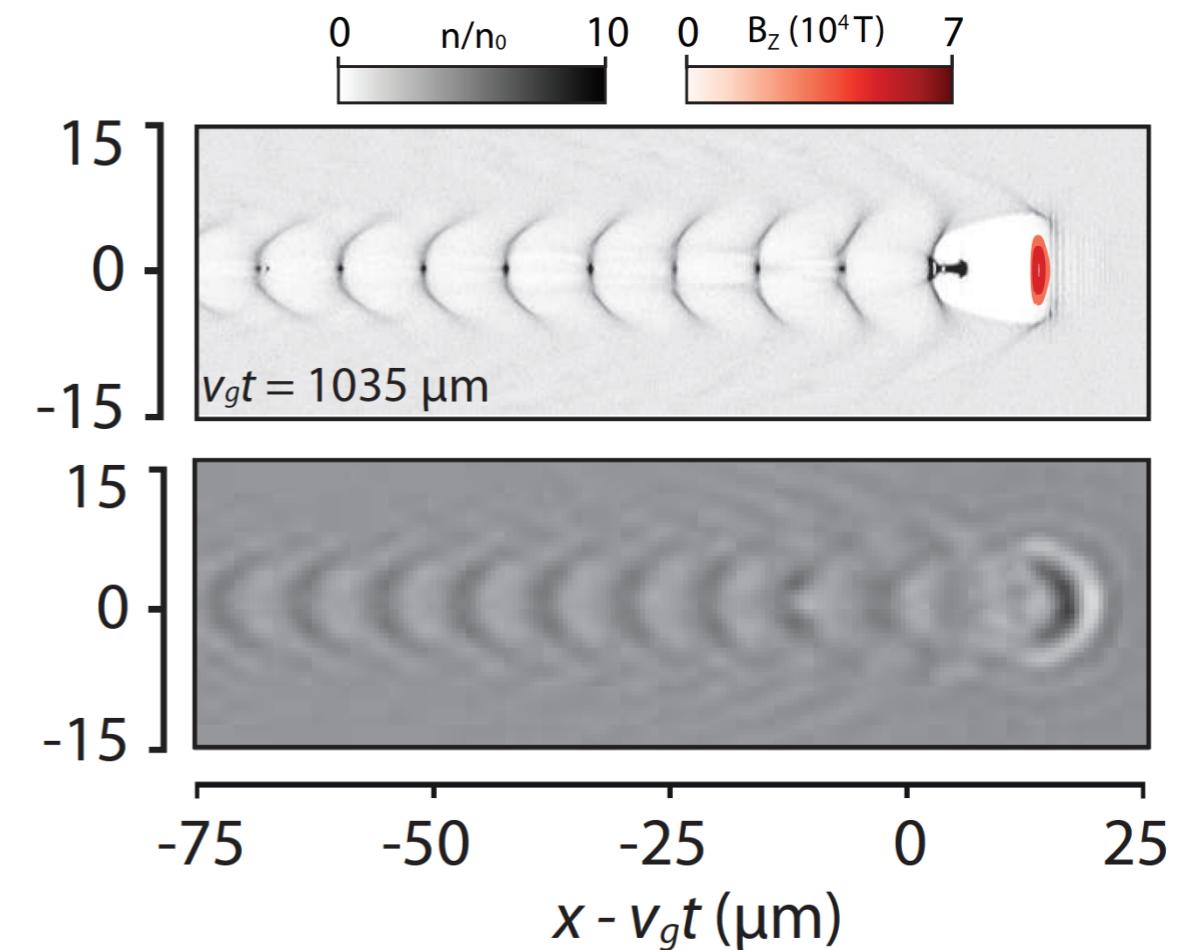
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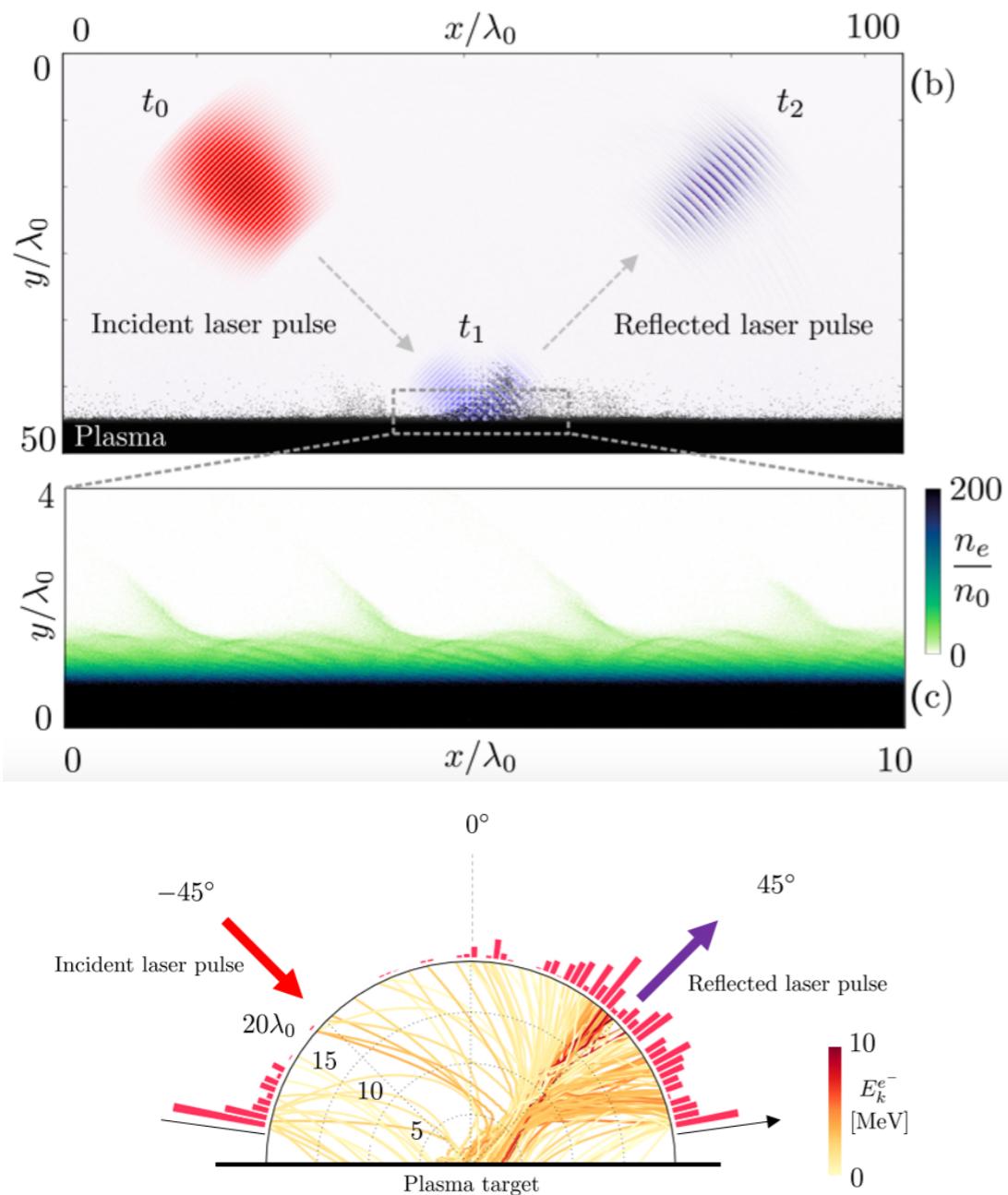
Laser wakefield acceleration of electrons



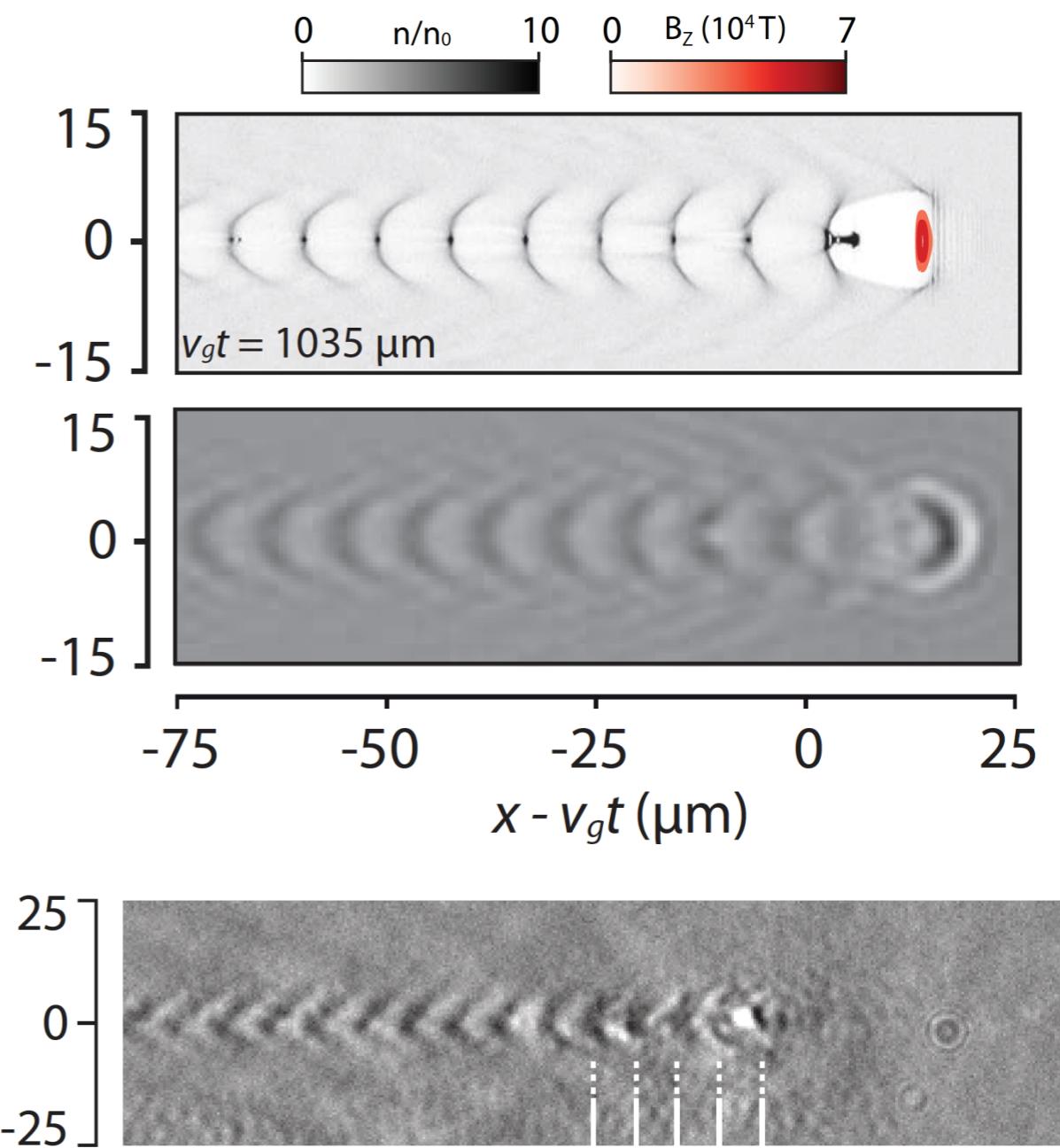
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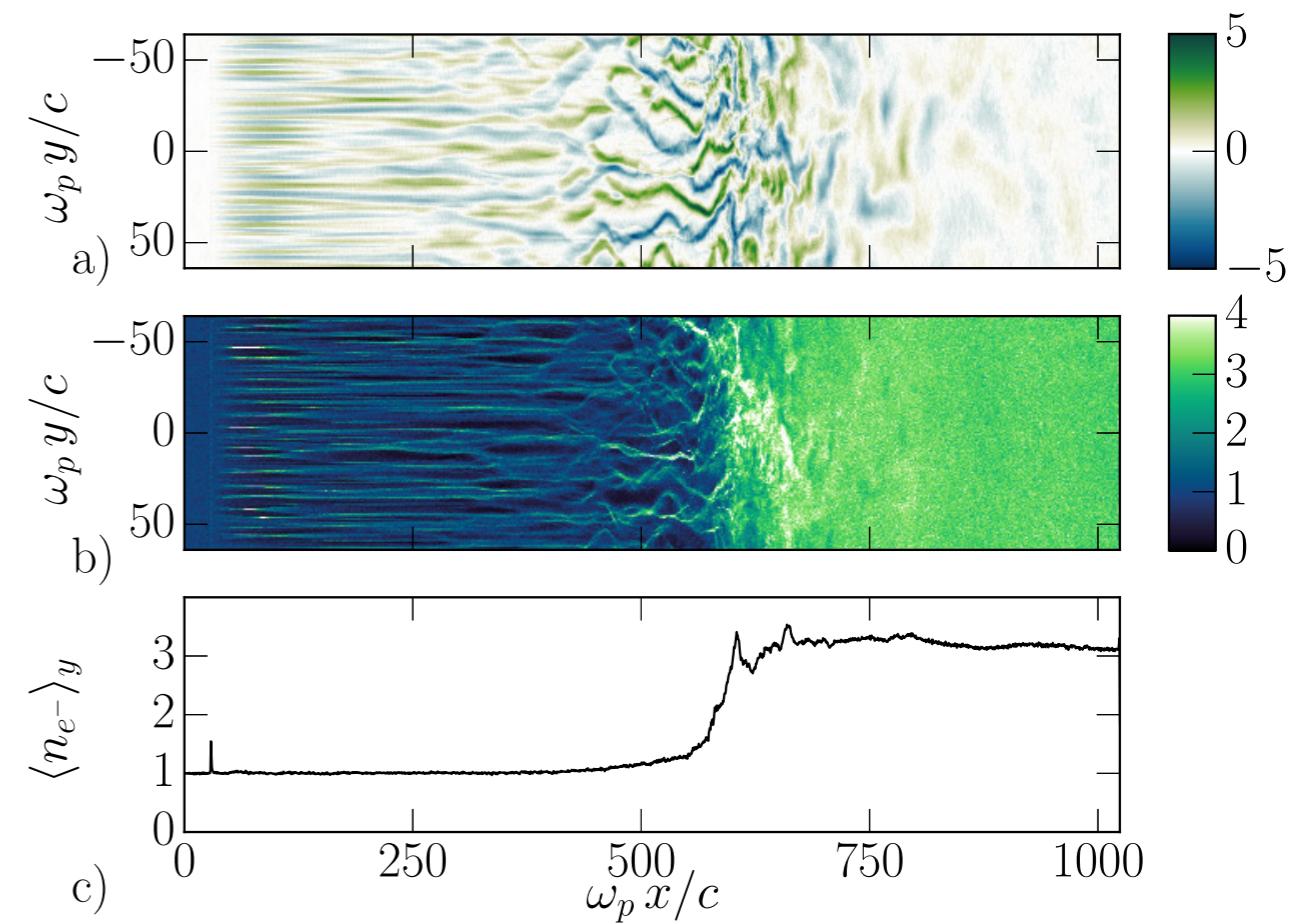
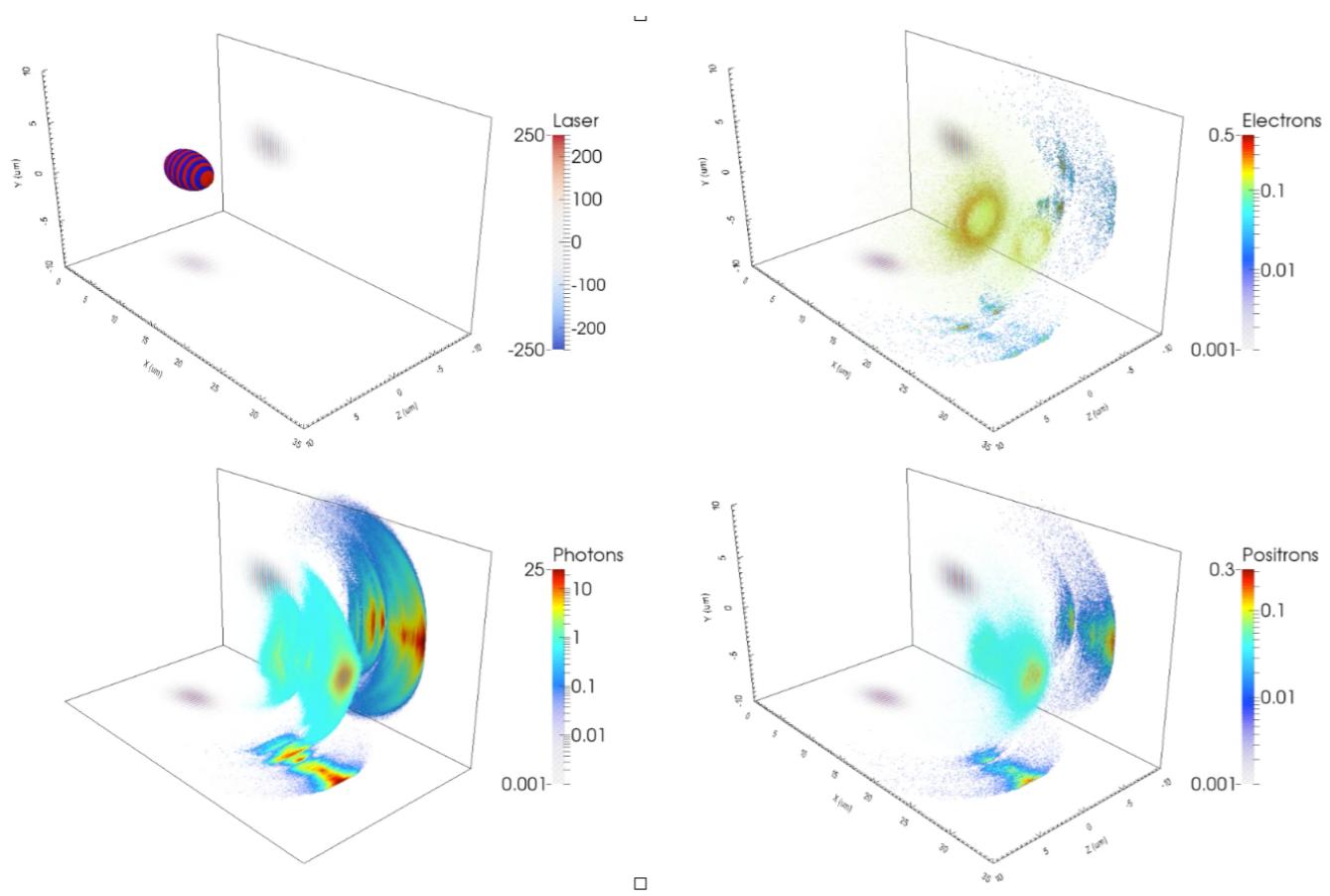
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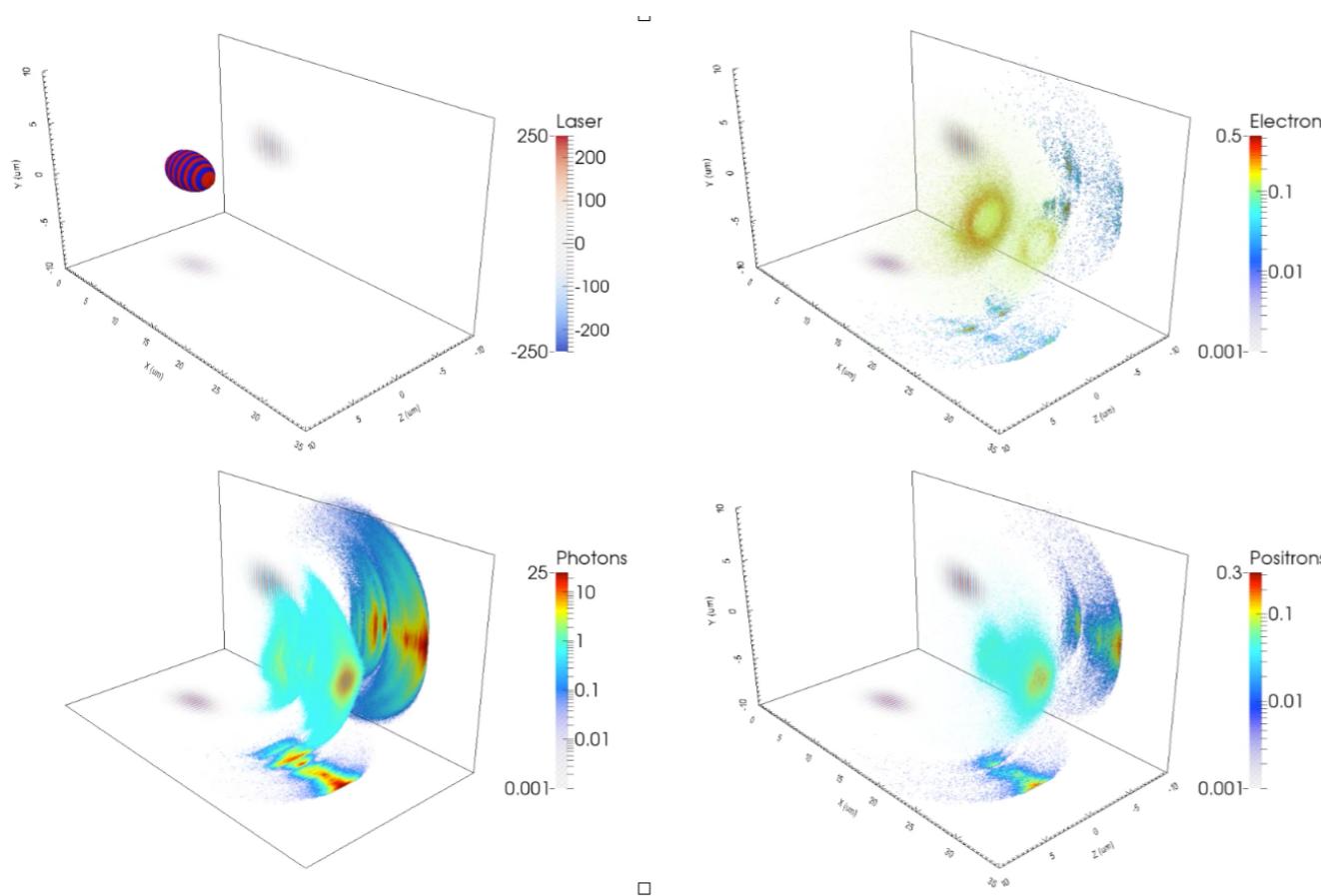


PIC codes are very **versatile**: they can be applied to a wide range of physical scenarii, from laser-plasma interaction to astrophysics

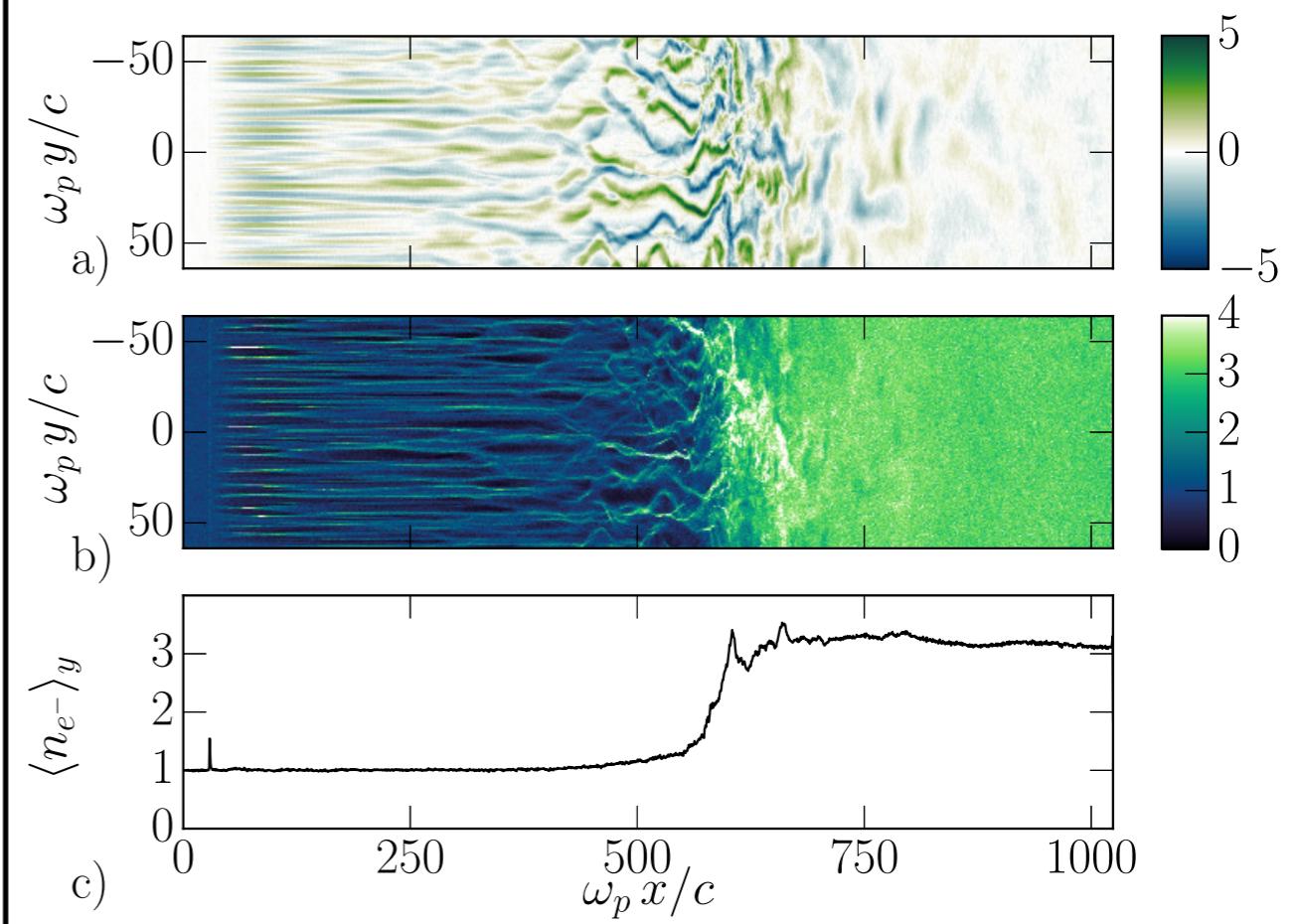


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Pair production on
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Relativistic shocks
in electron-positron plasmas



Conclusions

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- Implementation on new & future architectures requires a strong input (co-development) from HPC specialists

Checkout SMILEI !!!

Code, Diagnostics/visualization tools and tutorials
available online!



Smilei)

Smilei is a Particle-In-Cell code for plasma simulation. Open-source, collaborative, user-friendly and designed for high performances on super-computers, it is applied to a wide range of physics studies: from relativistic laser-plasma interaction to astrophysics.



Download



Github



Partners



Publications



Tutorials

<http://www.maisondelasimulation.fr/smilei>

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