

# Space-time hybrid methods for collisional kinetic equations

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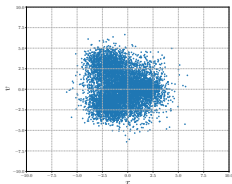
<sup>2</sup>Université Côte d'Azur



# System of particles in interactions

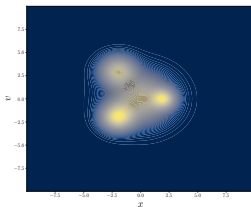
→ Multiscale systems

**Microscopic**



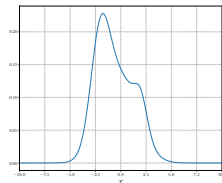
- Exact description
- Too costly simulations

**Mesoscopic**



- Microscopic effects
- Heavy simulations

**Macroscopic**



- Cheaper simulations
- Loss of microscopic effects

# The Vlasov-BGK equation

$f(t, x, v)$  distribution function,  $t > 0$ ,  $x \in \Omega_x \subset \mathbb{R}^{d_x}$  and  $v \in \Omega_v \subset \mathbb{R}^{d_v}$

$$\partial_t f + v \cdot \nabla_x f + E \cdot \nabla_v f = \mathcal{M}_{\rho, u, \theta} - f$$

Maxwellian and moments:

$$\mathcal{M}_{\rho, u, \theta} = \frac{\rho}{(2\pi\theta)^{d_v/2}} \exp\left(-\frac{|v - u|^2}{2\theta}\right)$$

$$\rho = \int_{\mathbb{R}^{d_v}} f \, dv; \quad u = \frac{1}{\rho} \int_{\mathbb{R}^{d_v}} v f \, dv; \quad \theta = \frac{1}{d_v \rho} \int_{\mathbb{R}^{d_v}} |v - u|^2 f \, dv;$$

$$\mathcal{E} = \int_{\mathbb{R}^{d_v}} \frac{|v|^2}{2} f \, dv$$

# From kinetic to fluid: a hydrodynamic limit

$$\text{Knudsen Number } \varepsilon = \frac{\text{Mean free path}}{\text{Characteristic length}}$$

## Hydrodynamic scaling

$$t \rightsquigarrow \frac{t}{\varepsilon} \quad \text{and} \quad x \rightsquigarrow \frac{x}{\varepsilon}$$

$$\partial_t f^\varepsilon + v \nabla_x f^\varepsilon + E \nabla_v f^\varepsilon = \frac{1}{\varepsilon} (\mathcal{M}_{\rho^\varepsilon, u^\varepsilon, \theta^\varepsilon} - f^\varepsilon)$$

$$f^\varepsilon(t, x, v) \xrightarrow{\varepsilon \rightarrow 0} \mathcal{M}_{\rho, u, \theta}$$

Euler with force:

$$\begin{cases} \partial_t \rho + \nabla_x \cdot (\rho u) = 0 \\ \partial_t (\rho u) + \nabla_x \cdot (\rho u \otimes u + pI) - \rho E = 0 \\ \partial_t \mathcal{E} + \nabla_x \cdot ((\mathcal{E} + p)u) - \rho u E = 0 \end{cases}$$

$$p = (\gamma - 1) \left( \mathcal{E} - \frac{\rho |u|^2}{2} \right), \quad \gamma = \frac{d_v + 2}{d_v}$$

# From kinetic to fluid: a diffusive limit

$$\text{Knudsen Number } \varepsilon = \frac{\text{Mean free path}}{\text{Characteristic length}}$$

## Diffusive scaling

$$t \rightsquigarrow \frac{t}{\varepsilon^2} \quad \text{et} \quad x \rightsquigarrow \frac{x}{\varepsilon}$$

$$\varepsilon^2 \partial_t f + \varepsilon (v \cdot \nabla_x f^\varepsilon + E \cdot \nabla_v f^\varepsilon) = \rho M - f$$

$$f(t, x, v) \underset{\varepsilon \rightarrow 0}{\rightarrow} \rho(t, x) M(v), \quad M(v) = \frac{e^{-|v|^2/2}}{(2\pi)^{d_v/2}}$$

Drift diffusion:

$$\partial_t \rho - \nabla_x \cdot (\nabla_x \rho - E \rho) = 0$$

# Kinetic simulations, a heavy numerical cost

- high-dimensional system :  $d_x = d_v = 3 \Rightarrow 7$  dimensions !
- Collision operators, small parameters

## → Hybrid methods

### Space

- Filbet, Rey (2015)
- Diffusive regime
- Domaine indicators
- Interface conditions

### Time

- Grigori, Hirstoaga, Salomon (2023)
- Collisions + hyperbolic
- Multiscale Pred.-corr.
- Parallel efficiency

### Space & Time

- AP scheme → **Stability and consistence**
- Spatial adaptation → **Fast phase-space solver**
- Multiscale parareal → **Parallel efficiency**

Hybridization in space (diffusive scaling)

Hybridization in time (hyperbolic scaling)

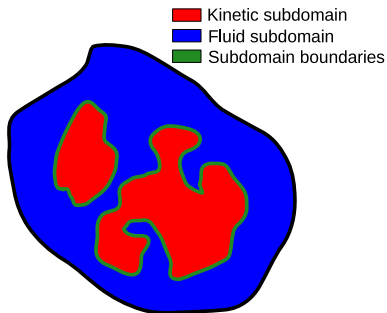
Hybridization in space and time (diffusive scaling)

Conclusion and perspectives

Hybridization in space (diffusive)

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# Domain adaptation<sup>1</sup>



- Maximize the size of the fluid domain
- Limit memory access
- Cell-wise decomposition
- Multiscale communications at interfaces

Should I **Kinetic** or Should I **Fluid**?

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<sup>1</sup>Laidin 2023; Laidin and Rey 2023.

**Diffusive scaling:**  $\varepsilon^2 \partial_t f + \varepsilon (v \cdot \partial_x f^\varepsilon + E \cdot \partial_v f^\varepsilon) = \rho M - f$

**Fluid limit:**  $\partial_t \rho - \partial_x \cdot (\partial_x \rho - E \rho) = 0$

**Micro-macro decomposition:**

$$f^\varepsilon(t, x, v) = \underbrace{\rho^\varepsilon(t, x) M(v)}_{\text{Local equilibria in velocity}} + \underbrace{g^\varepsilon(t, x, v)}_{\text{Perturbation}}$$

where  $M(v)$  is known.

$$\|g^\varepsilon(t, x, \cdot)\|_2 \leq \delta_0 \iff \text{locally at equilibrium}$$

# Macroscopic criterion

Truncated Chapman-Enskog expansion:

$$f^\varepsilon(t, x, v) = \rho^\varepsilon(t, x)M(v) + \sum_{k=1}^K \varepsilon^k h^{(k)}(t, x, v)$$

$$k = 0 : \quad h^{(1)} = -\mathbb{T}(\rho^\varepsilon M)$$

$$k = 1 : \quad h^{(2)} = -\partial_t(\rho^\varepsilon M) - \mathbb{T}(h^{(1)})$$

$$2 \leq k \leq K-1 : \quad h^{(k+1)} = -\partial_t h^{(k-1)} - \mathbb{T}(h^{(k)})$$

A hierarchy of fluid models:

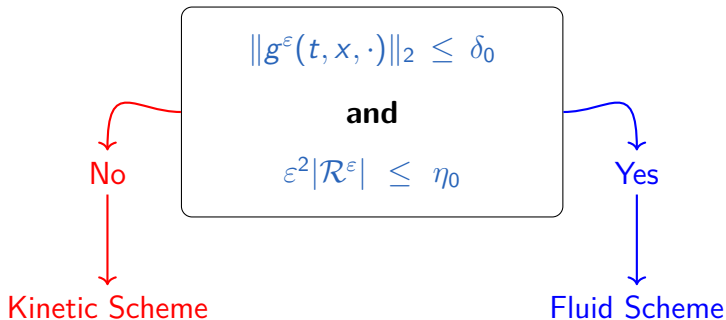
$$K = 1 \quad \partial_t \rho - \partial_x (\partial_x \rho - E\rho) = 0$$

$$K = 3 \quad \partial_t \rho^\varepsilon - \partial_x (\partial_x \rho^\varepsilon - E^\varepsilon \rho^\varepsilon) = \varepsilon^2 \mathcal{R}^\varepsilon$$

where  $\mathcal{R}^\varepsilon$  depends only on  $\rho^\varepsilon, E^\varepsilon$  and moments  $m_k = \int v^k M dv$

$$\varepsilon^2 |\mathcal{R}^\varepsilon| \leq \eta_0 \iff \text{the solution is locally close to the limit model}$$

**At each time step and for each position:**



## Continuous micro-macro model:

$$\begin{cases} \partial_t g^\varepsilon + \frac{1}{\varepsilon^2} g^\varepsilon + \frac{1}{\varepsilon} (\mathbb{T} g^\varepsilon - \partial_x \langle v g^\varepsilon \rangle M + v M J_x^\varepsilon) = 0 \\ \partial_t \rho^\varepsilon + \frac{1}{\varepsilon} \partial_x \langle v g^\varepsilon \rangle = 0 \end{cases}$$

## Continuous limit model ( $\varepsilon \rightarrow 0$ ):

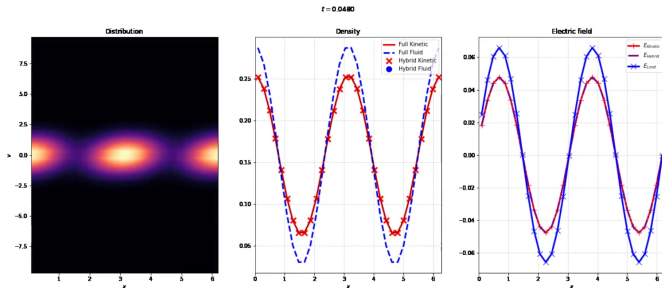
$$\partial_t \rho = \partial_x (\partial_x \rho - E \rho)$$

## Discretization ( $\Delta x, \Delta v, \Delta t > 0$ ):

- Primal cells  $\mathcal{X}_i = (x_{i-\frac{1}{2}}, x_{i+\frac{1}{2}})$ , staggered cells  $\mathcal{X}_{i+\frac{1}{2}} = (x_i, x_{i+1})$
- Velocity cells  $\mathcal{V}_j = (v_{j-\frac{1}{2}}, v_{j+\frac{1}{2}})$ , control volumes  $K_{i+\frac{1}{2},j} = \mathcal{X}_{i+\frac{1}{2}} \times \mathcal{V}_j$
- Unknowns:

$$\rho_i^{\varepsilon,n} \approx \frac{1}{\Delta x} \int_{\mathcal{X}_i} \rho^\varepsilon(t^n, x) dx, \quad g_{i+\frac{1}{2},j}^{\varepsilon,n} \approx \frac{1}{\Delta x \Delta v} \int_{K_{i+\frac{1}{2},j}} g^\varepsilon(t^n, x, v) dx dv$$

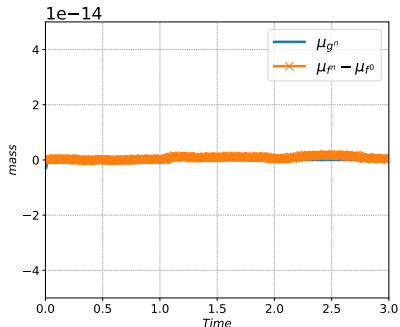
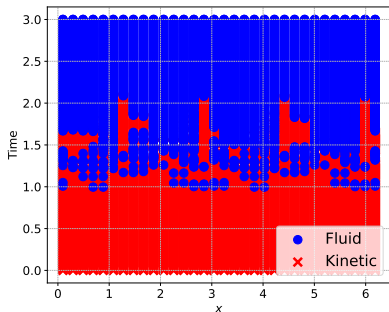
# Hybrid scheme $\varepsilon = 0.1$



$$f_0(x, v) = (2\pi)^{-3/2} (e^{-|v-3|^2/2} + e^{-|v+3|^2}) v_x^4 (1 + 0.05 \cos(2x))$$

# Mass conservation

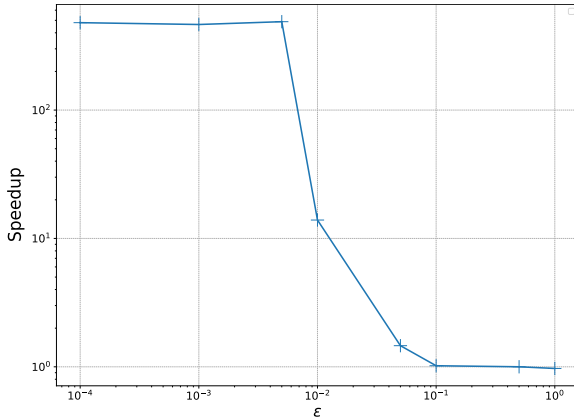
## State of cells and conservation property



$$f_0(x, v) = (2\pi)^{-3/2} (e^{-|v-3|^2/2} + e^{-|v+3|^2}) v_x^4 (1 + 0.05 \cos(2x)), \quad \varepsilon = 0.1$$

# Performances

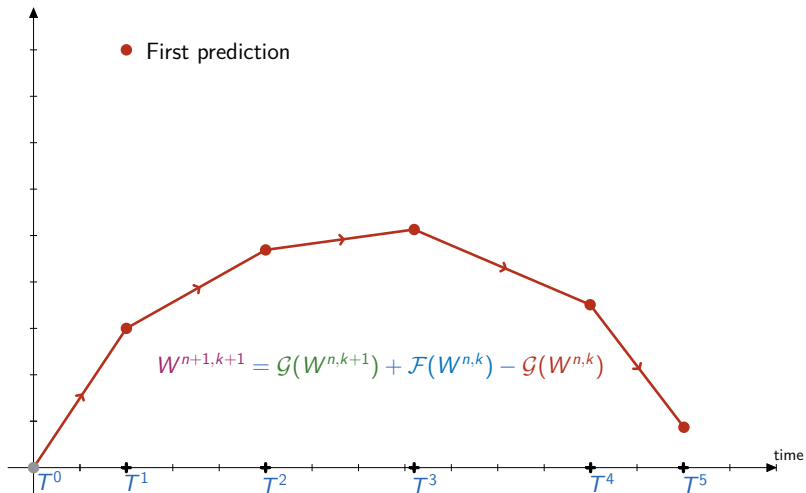
$$\textit{Speedup} = \frac{\textit{Kinetic CPU time}}{\textit{Hybrid CPU time}}$$



Hybridization in time (hyperbolic)

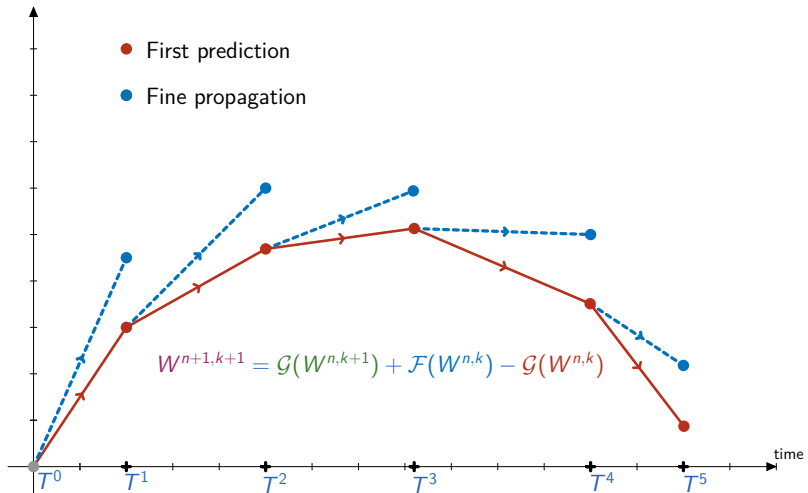
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# The parareal method<sup>2</sup>



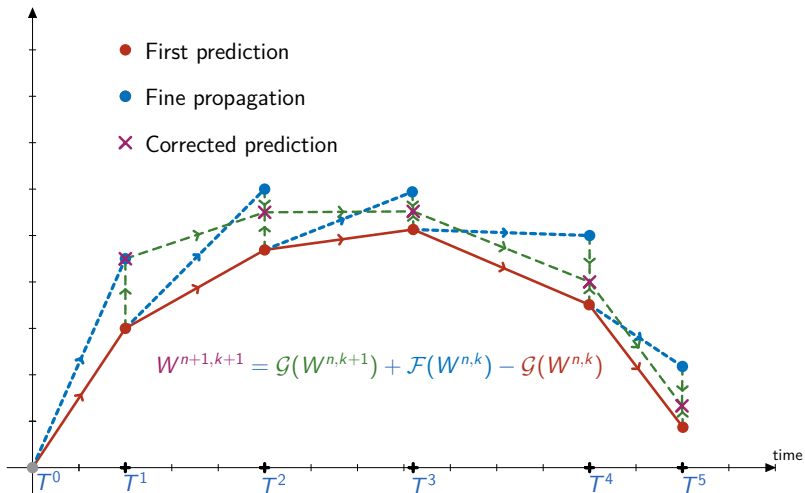
<sup>2</sup>Lions et al. 2001; Maday and Turinici 2002.

# The parareal method<sup>3</sup>



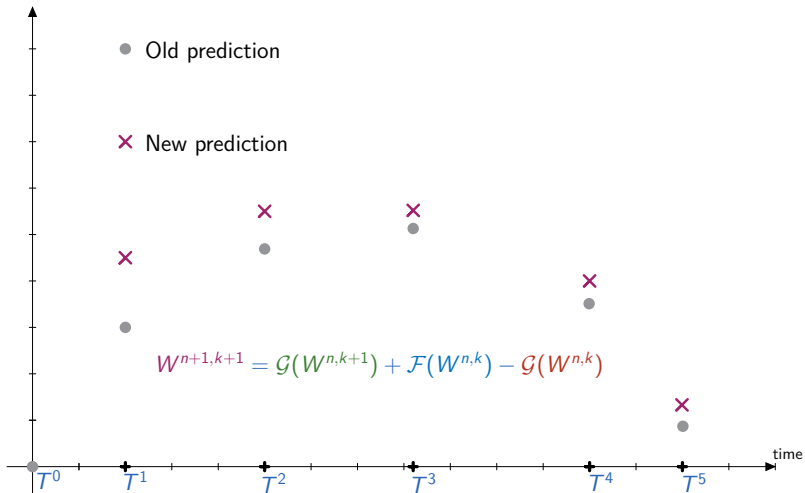
<sup>3</sup>Lions et al. 2001; Maday and Turinici 2002.

# The parareal method<sup>4</sup>



<sup>4</sup>Lions et al. 2001; Maday and Turinici 2002.

# The parareal method<sup>5</sup>



<sup>5</sup>Lions et al. 2001; Maday and Turinici 2002.

# Back to Kinetic: linking the scales

## Definition 1 (Projection to fluid)

For a distribution function  $f(t, x, v) \in L^1((1 + |v|)^2 dv)$ , we define the projection  $\mathcal{P}f(t, x) = U(t, x)$  as

$$\mathcal{P}f = U = \begin{pmatrix} \int_{\mathbb{R}^{d_v}} f dv \\ \frac{1}{\rho} \int_{\mathbb{R}^{d_v}} vf dv \\ \frac{1}{d_v \rho} \int_{\mathbb{R}^{d_v}} |v - u|^2 f dv \end{pmatrix} = \begin{pmatrix} \rho \\ u \\ \theta \end{pmatrix}$$

## Definition 2 (Lifting to kinetic)

For a macroscopic data  $U(t, x) \in \mathbb{R}^5$ , we define the lifting of order  $L$ ,  $\mathcal{L}^{(L)}U(t, x, v) = f(t, x, v)$  as

$$f = \mathcal{L}^{(L)}U = \mathcal{M}_{\rho, u, \theta} \left( 1 + \sum_{l=1}^L \varepsilon^l h^{(l)} \right)$$

The functions  $h^l$  are obtained through a Chapman-Enskog expansion.

# Kinetic Fluid Parareal Algorithm<sup>6</sup>

**Require:**  $U^{0,0}$

- 1: **for**  $n = 1, \dots, N_g$  **do** ▷ First coarse guess
- 2:      $U^{n,0} \leftarrow \mathcal{G}(U^{n-1,0})$
- 3: **end for**
  
- 4: **while**  $k \leq K$  **or** error  $\geq$  tol **do** ▷ Parareal iterations
  
- 5:     **for**  $n = k, \dots, N_g$  **do** ▷ Parallel computation of the jumps
- 6:          $\Delta^n = \mathcal{PFL}(U^{n-1,k-1}) - \mathcal{G}(U^{n-1,k-1})$
- 7:     **end for**
  
- 8:     **for**  $n = k, \dots, N_g$  **do** ▷ Sequential corrections
- 9:          $U^{n,k+1} = \mathcal{G}(U^{n-1,k+1}) + \Delta^n$
- 10:     **end for**
  
- 11:     Compute the successive error and  $k \leftarrow k + 1$
- 12: **end while**

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<sup>6</sup>Laidin and Rey 2025.

# Exterior force

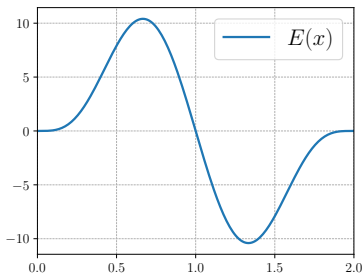
$$f_0(x, v) = \mathcal{M}_{\rho_1(x), u_1(x), \theta_1(x)}(v) + \mathcal{M}_{\rho_2(x), u_2(x), \theta_2(x)}(v), \quad x \in [0, 2], \quad v \in [-8, 8]$$

where

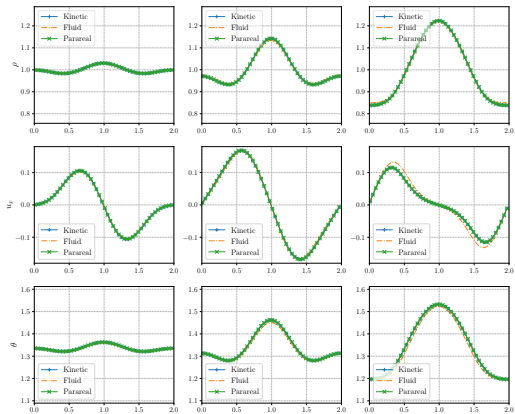
$$(\rho_1(x), u_1(x), \theta_1(x)) = (1, 1, 0, 0, 1)$$

and

$$(\rho_2(x), u_2(x), \theta_2(x)) = (1, -1, 0, 0, 1)$$

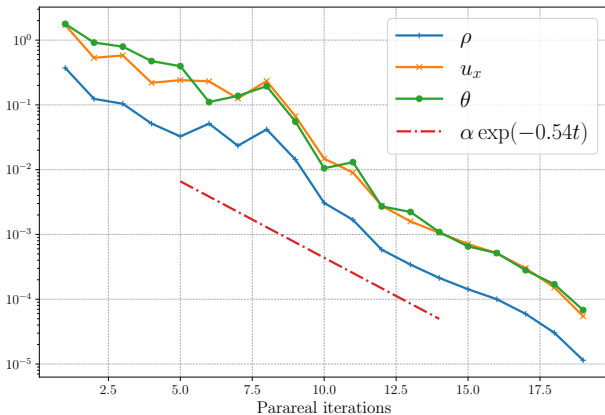


# Exterior force, $\varepsilon = 10^{-5}$ , $K = 20$



**Kinetic (Blue), Parareal (Green), Euler (Orange)**

# Exterior force, convergence



**Convergence of the successive errors**

# Exterior force, Speedup

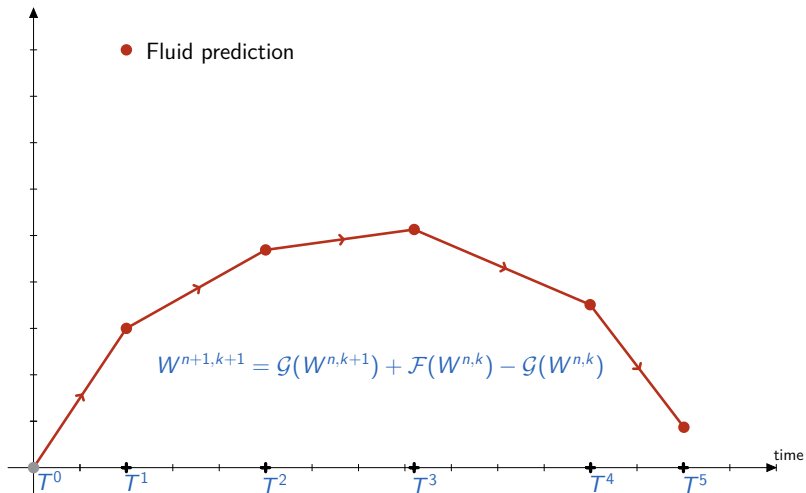
Parareal Iterations	Final error	Runtime (s)	Kinetic Runtime	Speedup
3	1.387E-03	71.7	672.8	9.38
6	7.782E-06	153.0	672.8	4.52
9	2.421E-08	218.4	672.8	3.08

**Performance on 32 threads,  $\varepsilon = 10^{-5}$**

Hybridization in space and time  
(diffusive)

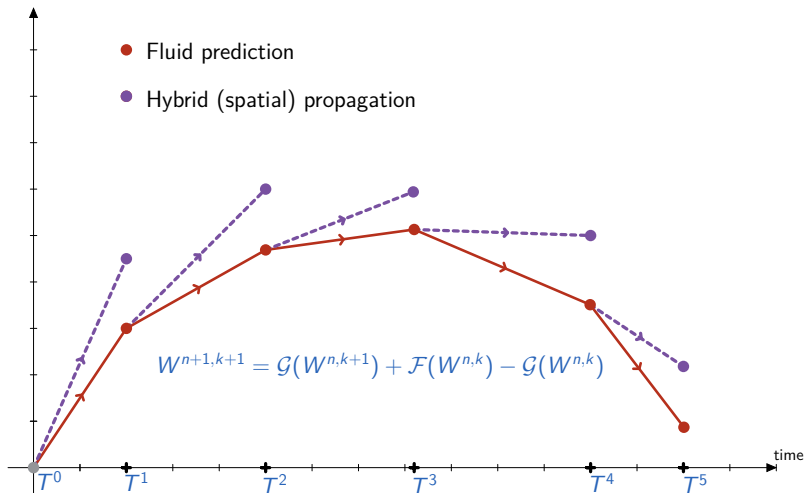
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# Space and time hybridization<sup>7</sup>



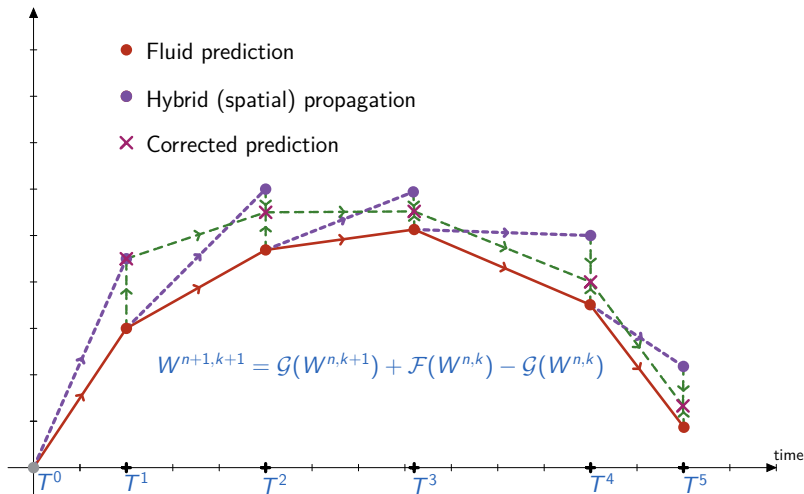
<sup>7</sup>Laidin 2025.

# Space and time hybridization<sup>8</sup>

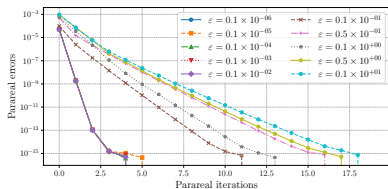
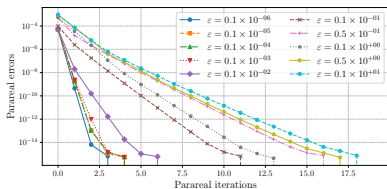


<sup>8</sup>Laidin 2025.

# Space and time hybridization<sup>9</sup>



<sup>9</sup>Laidin 2025.



**Parareal error without (Left) and with (Right) domain adaptation**

<sup>10</sup>Laidin 2025.

$\varepsilon$	Para. Off		Para. On		Fluid
	Adapt. Off	Adapt. On	Adapt. Off	Adapt. On	
1.0	<b>1.00</b>	<b>0.92</b>	<b>2.27</b>	<b>2.16</b>	<b>384</b>
$10^{-1}$	<b>1.00</b>	<b>0.92</b>	<b>2.73</b>	<b>2.56</b>	<b>384</b>
$10^{-2}$	<b>1.00</b>	<b>0.97</b>	<b>3.09</b>	<b>2.96</b>	<b>384</b>
$10^{-3}$	<b>1.00</b>	<b>4.67</b>	<b>5.65</b>	<b>24.02</b>	<b>384</b>
$10^{-4}$	<b>1.00</b>	<b>506</b>	<b>7.39</b>	<b>72.80</b>	<b>384</b>

**Speedups of the space-time hybrid methods**

## Conclusion and perspectives

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## → Dynamic domain adaptation

- Very fast in transient / near fluid regimes
- Robust w.r.t. the Knudsen number

## → Multiscale parareal algorithm

- Parallel in time
- Independent of phase-space solvers
- Robust w.r.t. the Knudsen number

## → Space time hybrid method

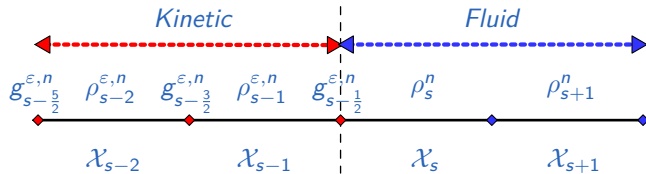
- AP schemes, spatial hybridation, parallel in time
- “Playing on all fields”

## Perspectives

- Higher order lifting procedure / learning
- Learn the domain indicators
- Parallelization: shared  $\rightarrow$  distributed + hybrid paradigm, accelerators
- Different physics: incompressibility, quasineutrality, gyrokinetic
- More complex models: Boltzmann / Landau, energy transport
- Error estimates

Thank you for your attention !

# Interface conditions



# Blast wave, pointwise errors

