# A semi-Lagrangian AMR scheme for 2D transport problems in conservation form 

Pep Mulet Mestre, Francesco Vecil

Universitat de València
CSASC, Koper/Capodistria (Slovenia), 12 June 2013

## Outline

(1) Introduction
(2) Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## Motivation



No need for fine meshing everywhere in the domain.


Refine only where the important information is.

## Framework

Equations
In dimension $N$, transport equations written in conservtion form:

$$
\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left[a_{1}\left(t, x_{1}, x_{2}\right) u\right]+\frac{\partial}{\partial x_{2}}\left[a_{2}\left(t, x_{1}, x_{2}\right) u\right]=0, \quad u\left(0, x_{1}, x_{2}\right)=u^{0}\left(x_{1}, x_{2}\right)
$$

where $\boldsymbol{a}: \mathbb{R}_{\geq 0} \times \Omega \rightarrow \mathbb{R}^{2}$ is the advection field.

## Example

The three-dimensional Vlasov-Maxwell equation

describes the evolution of $f(t, \boldsymbol{x}, \boldsymbol{p})$, typically representing the concentration of electrons or holes at position $\boldsymbol{x}$ and momentum $\boldsymbol{p}$.

[^0]
## Framework

## Equations

In dimension $N$, transport equations written in conservtion form:

$$
\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left[a_{1}\left(t, x_{1}, x_{2}\right) u\right]+\frac{\partial}{\partial x_{2}}\left[a_{2}\left(t, x_{1}, x_{2}\right) u\right]=0, \quad u\left(0, x_{1}, x_{2}\right)=u^{0}\left(x_{1}, x_{2}\right)
$$

where $\boldsymbol{a}: \mathbb{R}_{\geq 0} \times \Omega \rightarrow \mathbb{R}^{2}$ is the advection field.

## Example

The three-dimensional Vlasov-Maxwell equation

$$
\frac{\partial f}{\partial t}+\boldsymbol{v}(\boldsymbol{p}) \cdot \frac{\partial f}{\partial \boldsymbol{x}}+\boldsymbol{F} \cdot \frac{\partial f}{\partial \boldsymbol{p}}=0, \quad \boldsymbol{v}(\boldsymbol{p}):=\frac{\boldsymbol{p}}{m \sqrt{1+\frac{\mid \boldsymbol{p}^{2}}{m^{2} c^{2}}}}, \quad \boldsymbol{F}:=-e(\boldsymbol{E}+\boldsymbol{v}(\boldsymbol{p}) \wedge \boldsymbol{B})
$$

describes the evolution of $f(t, \boldsymbol{x}, \boldsymbol{p})$, typically representing the concentration of electrons or holes at position $\boldsymbol{x}$ and momentum $\boldsymbol{p}$.

[^1]
## Framework

## Equations

In dimension $N$, transport equations written in conservtion form:

$$
\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left[a_{1}\left(t, x_{1}, x_{2}\right) u\right]+\frac{\partial}{\partial x_{2}}\left[a_{2}\left(t, x_{1}, x_{2}\right) u\right]=0, \quad u\left(0, x_{1}, x_{2}\right)=u^{0}\left(x_{1}, x_{2}\right)
$$

where $\boldsymbol{a}: \mathbb{R}_{\geq 0} \times \Omega \rightarrow \mathbb{R}^{2}$ is the advection field.

## Example

The three-dimensional Vlasov-Maxwell equation

$$
\frac{\partial f}{\partial t}+\boldsymbol{v}(\boldsymbol{p}) \cdot \frac{\partial f}{\partial \boldsymbol{x}}+\boldsymbol{F} \cdot \frac{\partial f}{\partial \boldsymbol{p}}=0, \quad \boldsymbol{v}(\boldsymbol{p}):=\frac{\boldsymbol{p}}{m \sqrt{1+\frac{\mid \boldsymbol{p}^{2}}{m^{2} c^{2}}}}, \quad \boldsymbol{F}:=-e(\boldsymbol{E}+\boldsymbol{v}(\boldsymbol{p}) \wedge \boldsymbol{B})
$$

describes the evolution of $f(t, \boldsymbol{x}, \boldsymbol{p})$, typically representing the concentration of electrons or holes at position $\boldsymbol{x}$ and momentum $\boldsymbol{p}$.

Features
Shocks, large gradients, filamentation, microscopic structures.

## Outline

## (1) Introduction

(2) Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## Grid hierarchy

Resolution levels
We define $L+1$ resolution levels: the coarsest is $\ell=0$, the finest $\ell=L$. In 1D, the meshes are

$$
x_{\ell, j}=x_{\min }+j \Delta x_{\ell}, \quad \Delta x_{\ell}=\frac{x_{\max }-x_{\min }}{2^{\ell} N_{0}}
$$

Grid
The $\ell$-grid at time $t^{n}$ is


We are interested in


## Grid hierarchy

Resolution levels
We define $L+1$ resolution levels: the coarsest is $\ell=0$, the finest $\ell=L$. In 1D, the meshes are

$$
x_{\ell, j}=x_{\min }+j \Delta x_{\ell}, \quad \Delta x_{\ell}=\frac{x_{\max }-x_{\min }}{2^{\ell} N_{0}}
$$

Grid
The $\ell$-grid at time $t^{n}$ is

$$
G_{\ell}^{n}=\left\{x_{\ell, j}\right\}_{j \in \mathcal{G}_{\ell}^{n}}
$$

We are interested in

$$
\mathcal{G}_{\ell}^{n} \subseteq \prod_{i=1}^{N}\left\{0, \ldots, N_{i, \ell}\right\}
$$

## Grid hierarchy

Nesting condition
We are interested in nested meshes:


## Ghost points

Ghost points are added outside the $\ell$-grids to take into account the boundary conditions for the time integration.

## Grid hierarchy

Nesting condition
We are interested in nested meshes:


## Ghost points

Ghost points are added outside the $\ell$-grids to take into account the boundary conditions for the time integration.

## Grid selection

The grid selection is based on two criteria: prediction and gradient.
The prediction criterion
Let $\left\{u_{\ell, j}\right\}_{j \in \mathcal{G}_{\ell}^{n}}$ the point values at reslution level $\ell$. Let $p\left[u_{\ell}\right]$ an interpolator at resolution level $\ell$. We keep point $x_{\ell, j}$ for refinement if $\left|p\left[u_{\ell-1}\right]\left(x_{\ell, j}\right)-u_{\ell, j}\right|>\tau_{p}$.

## The gradient criterion

We estimate the gradient at point $x_{e, j}$ by means of the discrete gradient computed at resolution level $\ell-1$ : if it is over a certain tolerance parameter $\tau_{d, \ell-1}$, then point $x_{\ell, j}$ is selected for refinement.

## Reconstruction

Once the grid has been selected, reconstruct by means of an interpolator the point values that are not assigned yet.

## Grid selection

The grid selection is based on two criteria: prediction and gradient.
The prediction criterion
Let $\left\{u_{\ell, j}\right\}_{j \in \mathcal{G}_{\ell}^{n}}$ the point values at reslution level $\ell$. Let $p\left[u_{\ell}\right]$ an interpolator at resolution level $\ell$. We keep point $x_{\ell, j}$ for refinement if

$$
\left|p\left[u_{\ell-1}\right]\left(x_{\ell, j}\right)-u_{\ell, j}\right|>\tau_{p}
$$

The gradient criterion
We estimate the gradient at point $x_{\ell, j}$ by means of the discrete gradient computed at resolution level $\ell-1$ : if it is over a certain tolerance parameter $\tau_{d, \ell-1}$, then point $x_{\ell, j}$ is selected for refinement.

Reconstruction
Once the grid has been selected, reconstruct by means of an interpolator the point values that are not assigned yet.

## Grid selection

The grid selection is based on two criteria: prediction and gradient.
The prediction criterion
Let $\left\{u_{\ell, j}\right\}_{j \in \mathcal{G}_{\ell}^{n}}$ the point values at reslution level $\ell$. Let $p\left[u_{\ell}\right]$ an interpolator at resolution level $\ell$. We keep point $x_{\ell, j}$ for refinement if

$$
\left|p\left[u_{\ell-1}\right]\left(x_{\ell, j}\right)-u_{\ell, j}\right|>\tau_{p} .
$$

The gradient criterion
We estimate the gradient at point $x_{\ell, j}$ by means of the discrete gradient computed at resolution level $\ell-1$ : if it is over a certain tolerance parameter $\tau_{d, \ell-1}$, then point $x_{\ell, j}$ is selected for refinement.

Reconstruction
Once the grid has been selected, reconstruct by means of an interpolator the point values that are not assigned yet.

## Grid selection

The grid selection is based on two criteria: prediction and gradient.
The prediction criterion
Let $\left\{u_{\ell, j}\right\}_{j \in \mathcal{G}_{\ell}^{n}}$ the point values at reslution level $\ell$. Let $p\left[u_{\ell}\right]$ an interpolator at resolution level $\ell$. We keep point $x_{\ell, j}$ for refinement if

$$
\left|p\left[u_{\ell-1}\right]\left(x_{\ell, j}\right)-u_{\ell, j}\right|>\tau_{p} .
$$

## The gradient criterion

We estimate the gradient at point $x_{\ell, j}$ by means of the discrete gradient computed at resolution level $\ell-1$ : if it is over a certain tolerance parameter $\tau_{d, \ell-1}$, then point $x_{\ell, j}$ is selected for refinement.

## Reconstruction

Once the grid has been selected, reconstruct by means of an interpolator the point values that are not assigned yet.

## Outline

(2) Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## 1D semi-Lagrangian strategy

Characteristic-based solution
The solution to the PDE

$$
\frac{\partial u}{\partial t}+\frac{\partial}{\partial x}[a(t, x) u]=0, \quad u(0, x)=u^{0}(x)
$$

is given by $\quad u(t, x)=u(s, \mathcal{X}(s ; t, x)) J(s ; t, x)$,
with $\mathcal{X}(s ; t, x)$ the characteristic at time $s$, starting from $x$ at time $t$ :

$$
\frac{\mathrm{d} \mathcal{X}(s ; t, x)}{\mathrm{d} s}=a(s, \mathcal{X}(s ; t, x)), \quad \mathcal{X}(t ; t, x)=x, \quad J(s ; t, x):=\frac{\partial \mathcal{X}(s ; t, x)}{\partial x}
$$



## 1D semi-Lagrangian strategy

What do we need?

- A solver for the characteristics $\mathcal{X}(s ; t, x)$ : Runge-Kutta.
- An approximation for the Jacobian through cenetred finite differences:

$$
J(s ; t, x)=\frac{\partial \mathcal{X}(s ; t, x)}{\partial x} \approx \frac{\tilde{\mathcal{X}}(s ; t, x+\delta x)-\tilde{\mathcal{X}}(s ; t, x-\delta x)}{2 \delta x}, \quad \delta x=10^{-m} \Delta x_{\ell}
$$

- An interpolator to reconstruct $u(s, \mathcal{X}(s ; t, x))$ : PVWENO (Point-Value Weighted Essentially Non-Oscillatory).

Error estimate
The local truncation error can be estimated


## 1D semi-Lagrangian strategy

What do we need?

- A solver for the characteristics $\mathcal{X}(s ; t, x)$ : Runge-Kutta.
- An approximation for the Jacobian through cenetred finite differences:

$$
J(s ; t, x)=\frac{\partial \mathcal{X}(s ; t, x)}{\partial x} \approx \frac{\tilde{\mathcal{X}}(s ; t, x+\delta x)-\tilde{\mathcal{X}}(s ; t, x-\delta x)}{2 \delta x}, \quad \delta x=10^{-m} \Delta x_{\ell}
$$

- An interpolator to reconstruct $u(s, \mathcal{X}(s ; t, x))$ : PVWENO (Point-Value Weighted Essentially Non-Oscillatory).


## Error estimate

The local truncation error can be estimated

$$
E=\underbrace{\underbrace{\mathcal{O}\left(\Delta t^{s+1}\right)}_{\text {Runge-Kutta }}+\underbrace{\mathcal{O}\left(\Delta x_{\ell}^{2}\right)}_{\text {Jacobian }}+\underbrace{\mathcal{O}\left(\Delta x_{\ell}^{2 r}\right)}_{\text {PVWENO (advection) }} . . \underbrace{\mathcal{O}} .}_{\text {approximation of the characteristics }}
$$

## The 2D case

Grid hierarchy and selection
We do not give details, but we apply strategies similar to the 1D case.

## The 2D PDE

We solve the 2D PDE $\quad \frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0$
by splitting the $\left(x_{1}, x_{2}\right)$-domain thanks to the second-order Strang scheme:

- Solve for a $\frac{\Delta t}{2}$ time sten $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$;
- Solve for a $\Delta t$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0$;
- Solve for a $\frac{\Delta t}{2}$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$.


## Error estimate

## The Strang splitting constrains the accuracy:



## The 2D case

Grid hierarchy and selection
We do not give details, but we apply strategies similar to the 1D case.

The 2D PDE
We solve the 2D PDE

$$
\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0
$$

by splitting the $\left(x_{1}, x_{2}\right)$-domain thanks to the second-order Strang scheme:

- Solve for a $\frac{\Delta t}{2}$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$;
- Solve for a $\Delta t$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0$;
- Solve for a $\frac{\Delta t}{2}$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$.

Error estimate
The Strang splitting constrains the accuracy:


## The 2D case

Grid hierarchy and selection
We do not give details, but we apply strategies similar to the 1D case.

## The 2D PDE

We solve the 2D PDE $\quad \frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0$
by splitting the $\left(x_{1}, x_{2}\right)$-domain thanks to the second-order Strang scheme:

- Solve for a $\frac{\Delta t}{2}$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$;
- Solve for a $\Delta t$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{2}}\left(a_{2} u\right)=0$;
- Solve for a $\frac{\Delta t}{2}$ time step $\frac{\partial u}{\partial t}+\frac{\partial}{\partial x_{1}}\left(a_{1} u\right)=0$.


## Error estimate

The Strang splitting constrains the accuracy:

$$
E=\mathcal{O}\left(\Delta t^{\min (s+1,3)}\right)+\underbrace{\mathcal{O}\left(\Delta x_{1, \ell}^{2}\right)+\mathcal{O}\left(\Delta x_{2, \ell}^{2}\right)}_{\text {approximated Jacobian }}+\underbrace{\mathcal{O}\left(\Delta x_{1, \ell}^{2 r}\right)+\mathcal{O}\left(\Delta x_{2, \ell}^{2 r}\right)}_{\text {PVWENO }} .
$$

## Outline

(2) Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## Goals

The Adaptive-Mesh-Refinement (AMR) framework is compared to the equivalent Fixed-Mesh (FM) results.

Of course, AMR cannot be more accurate than FM. Rather, it achieves faster computational times in exchange of a loss of precision.

## Outline

## (1) Introduction

(2) Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## Variable-coefficient advection

The test case $\quad \frac{\partial u}{\partial t}+\frac{\partial}{\partial x}(\sin (x) u)=0 \quad$ produces a blow-up.

(a) at time 0
(c) at time 0.89

(b) at time 0.65

(d) at time 1.48

## Variable-coefficient advection

Speedup
For parameters

$$
\begin{aligned}
N_{0} & =128 \\
L & =4 \\
\Delta t_{0} & =0.125 \\
\tau_{p} & =10^{-4} \\
\tau_{d, 0} & =0.5
\end{aligned}
$$

AMR reaches a speedup of 35 times with respect to the equivalently-resolved FM , with a loss of precision from $10^{-9}$ to roughly $10^{-6}$ (the $L^{2}$-error w.r.t. the analytical solution).

## Outline

(2)

Numerical tools

- Multiresolution framework
- Time integration
(3) Experiments
- Introduction
- 1D tests
- 2D tests


## Deformation flows

The system

$$
\frac{\partial f}{\partial t}+\frac{\partial}{\partial x}\left[\sin ^{2}(\pi x) \sin (2 \pi y) g(t) f\right]+\frac{\partial}{\partial y}\left[-\sin ^{2}(\pi y) \sin (2 \pi x) g(t) f\right]=0, \quad(x, y) \in[0,1]^{2}
$$


for $g(t)=\cos \left(\frac{\pi t}{T}\right)$, periodically recovers the initial datum after alternate clockwise and counterclockwise twistings.

Performances
Speedup $\approx 2$.

## Deformation flows

The ODE integrator for the characteristics


As announced by the error estimate

$$
E=\mathcal{O}\left(\Delta x_{1, \ell}^{2}\right)+\mathcal{O}\left(\Delta x_{2, \ell}^{2}\right)+\mathcal{O}\left(\Delta t^{\min (s+1,3)}\right)
$$

the Strang-splitting order constrains the accuracy.

## Kelvin-Hemlholtz instabilities

## The model

The guiding-center model (omitting some details)

$$
\frac{\partial \rho}{\partial t}+\frac{\partial}{\partial x_{1}}\left[\frac{\partial \Phi}{\partial x_{2}} \rho\right]+\frac{\partial}{\partial x_{2}}\left[-\frac{\partial \Phi}{\partial x_{1}} \rho\right]=0, \quad \Delta_{x_{1}, x_{2}} \Phi=\rho
$$

for initial condition $\quad \rho\left(0, x_{1}, x_{2}\right)=1.5 \operatorname{sech}\left(\frac{x_{2}}{0.9}\right) \cdot\left(1+0.08 \sin \left(2 k x_{1}\right)\right)$, periodic $x_{1}$ - and Dirichlet $x_{2}$-boundaries, produces vortices and filamentation.

(f) first instability

## Kelvin-Hemlholtz instabilities



(h) steady state

## SPEEDUP

The AMR strategy achieves a speedup of about 2.8, for parameters $\left(N_{\star, 0}, L\right)=(32,4)$ (the maximum resolution is $512 \times 512), \tau_{p}=10^{-3}$, $\tau_{d, 0}=1.5$.
(g) second instability

# GRÀCIES! HVALA! DANKE! ĎAKUJEM! DĚKUJI! 


[^0]:    Features

[^1]:    Features

