Low-rank approximation for nonlinear kinetic problems Part 2: Asymptotic preserving dynamical low-rank schemes

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Transport equation

$$\frac{1}{c}\partial_t I(t,x,\Omega) + \Omega \cdot \nabla_x I(t,x,\Omega) = C(I) - A(I) + G,$$

where $\Omega \in \mathbb{S}^2$ and c is the propagation speed.

Right-hand side depends on the physical problem (Collision, absorption, and source terms).

A **Boltzmann equation** that is of interest in radiation therapy, radiative cooling, neutron transport in nuclear reactors, ...

Main difference with Vlasov equation is lack of long range interaction.

Linear radiative transport equation in diffusive scaling

$$\partial_t f + \frac{1}{\varepsilon} \mathbf{v} \cdot \nabla_{\mathbf{x}} f = \frac{\sigma^S}{\varepsilon^2} \left(\frac{1}{4\pi} \langle f \rangle_{\mathbf{v}} - f \right) - \sigma^A f + G.$$

A up to **5-dimensional** equation for f(t, x, v) ($x \in \mathbb{R}^3$ and $v \in \mathbb{S}^2$) with complex interplay of transport, collision (relaxation to equilibrium), absorption, and source term.

Collision is the stiffest term in the equation (for small ϵ).

For $\epsilon \rightarrow 0$ we obtain the limit

$$\partial_t \rho - \nabla_{\mathsf{x}} \cdot \left(\frac{1}{3\sigma^5} \nabla_{\mathsf{x}} \rho\right) = -\sigma^{\mathsf{A}} \rho + \mathcal{G}, \qquad \rho = \frac{1}{4\pi} \langle f \rangle_{\mathsf{v}} = \frac{1}{4\pi} \int_{\mathbb{S}^2} f \, \mathrm{d} \mathsf{v}.$$

This solution has rank 1 with $f(t, x, v) = \rho(t, x)$.

Diffusive limit

We start with a Hilbert expansion/Chapman-Enskog theory

 $f = f_0 + \epsilon f_1 + \epsilon^2 f_2 + \dots$

Plugging into the radiative transport equation (neglecting absorption and source terms)

$$\partial_t f + \frac{1}{\varepsilon} \mathbf{v} \cdot \nabla_{\mathbf{x}} f = \frac{\sigma}{\varepsilon^2} \left(\frac{1}{4\pi} \langle f \rangle_{\mathbf{v}} - f \right)$$

and collecting terms of the same order in ϵ .

 $\mathcal{O}(1/\epsilon^2)$: $f_0 = \langle f_0 \rangle_v / (4\pi)$ and thus f_0 does not depend on v.

For $\epsilon \to 0$ we have $f = \rho$.

 $\mathcal{O}(1/\epsilon)$:

$$\mathbf{v} \cdot \nabla_{\mathbf{x}} \mathbf{f}_{\mathbf{0}} = \sigma \left(\frac{1}{4\pi} \langle f_{\mathbf{1}} \rangle_{\mathbf{v}} - f_{\mathbf{1}} \right)$$

Diffusive limit

 $\mathcal{O}(1)$:

$$\partial_t f_0 + \mathbf{v} \cdot \nabla_{\mathbf{x}} f_1 = \sigma \left(\frac{1}{4\pi} \langle f_2 \rangle_{\mathbf{v}} - f_2 \right) \implies \partial_t \langle f_0 \rangle_{\mathbf{v}} + \nabla_{\mathbf{x}} \cdot \langle \mathbf{v} f_1 \rangle_{\mathbf{v}} = 0$$

From the $1/\epsilon$ term we get by multiplying with v and integrating

$$\mathbf{v} \cdot \nabla_{\mathsf{x}} \mathbf{f}_{\mathsf{0}} = \sigma \left(\frac{1}{4\pi} \langle f_{\mathsf{1}} \rangle_{\mathsf{v}} - f_{\mathsf{1}} \right) \implies \langle \mathbf{v} f_{\mathsf{1}} \rangle_{\mathsf{v}} = -\frac{1}{\sigma} \langle \mathbf{v} \nabla_{\mathsf{x}} \cdot (\mathbf{v} f_{\mathsf{0}}) \rangle_{\mathsf{v}} = -\frac{1}{3\sigma} \nabla_{\mathsf{x}} \mathbf{f}_{\mathsf{0}}$$

Putting everything together and using $\rho = \langle f_0 \rangle_v / 4\pi + O(\epsilon)$ we get the diffusion equation

$$\partial_t
ho =
abla_x \cdot \left(rac{1}{3\sigma}
abla_x
ho
ight) + \mathcal{O}(\epsilon).$$

- We have to get up to *third order* in ϵ to get the dynamics of f_0 .
- ► The collision operator determines how the equilibrium looks like.
- ► The transport determines the dynamics of the equilibrium.

Interest in the limit

For ϵ small we obtain information on the low-rank structure of the solution

$$f(t,x,v) = f_0(t,x) + \epsilon \left(\frac{1}{4\pi} \langle f_1 \rangle_v - \frac{v}{\sigma} \cdot \nabla_x f_0\right) + \mathcal{O}(\epsilon^2)$$

Up to $\mathcal{O}(\epsilon)$ the solution is **at most rank 4.**

This is one of the few situations where we do know how the low-rank structure of the solution looks like.

Sidenote: If we assume enough smoothness. Then,

$$f(x, v) \approx \sum_{k=1}^{n} \sum_{m=1}^{n} \hat{f}_{km} \mathrm{e}^{ikx} \mathrm{e}^{imv}$$

gives a low-rank approximation with moderate rank. Usually not a valid assumption for hyperbolic problems.

Error analysis for DLR in the diffusion limit

We again use the projector splitting integrator.

L step in the continuous setting

$$\partial_t L_i = -\frac{1}{\epsilon} \sum_{j=1}^r \mathbf{v} \cdot \langle X_i^n, \nabla X_j^n \rangle_{\times} L_j + \frac{1}{\epsilon^2} \sum_{j=1}^r \langle X_i^n \sigma X_j^n \rangle_{\times} (\langle L_j \rangle_{\mathbf{v}} - L_j).$$

For $\epsilon \to 0$ we must have $L_j = \langle L_j \rangle_{\mathbf{v}}$, **L** looses dependence on **v**.

Discretization

Discretization with $L^n = [L_1(v_{\cdot}), L_2(v_{\cdot}), \dots, L_r(v_{\cdot})] \in \mathbb{R}^{n_v \times r}$

$$\frac{\mathsf{L}^{\star} - \mathsf{L}^{n}}{\Delta t} + \frac{1}{\epsilon} \sum_{k=1}^{d} \mathsf{A}_{\partial_{k}}^{n} \mathsf{L}^{\star} \Pi_{v_{k}} = \frac{\mathsf{A}_{\sigma}^{n}}{\epsilon^{2}} \mathsf{L}^{\star} \mathsf{C},$$

QR decomposition: $\mathsf{L}^{\star} = \mathsf{V}^{\star}(\mathsf{S}^{\star})^{\top},$

where

$$\mathsf{A}^n_{\partial_k} = (\mathsf{X}^n)^\top \mathsf{D}_k \mathsf{X}^n, \quad \mathsf{A}^n_{\sigma} = \mathsf{X}^\top \Sigma \mathsf{X}, \quad \Pi_{v_k} = \mathsf{diag}(v_k), \quad \mathsf{C} = \frac{1}{n_v} e e^\top - \mathsf{I}$$

with

$$e = (1, 1, \dots, 1)^{\top}, \qquad \Sigma = \operatorname{diag}(\sigma), \qquad \mathsf{D}_k \approx \partial_{\mathsf{x}_k}.$$

We have used an implicit Euler step here.

Input: X^n , S^n and V^n . **Algorithm:**

- Construct $A^n_{\partial_k}, A^n_{\sigma}$; Compute $L^n = V^n(S^n)^\top$;
- Integrate L step with step size Δt to get L*;
- ▶ Perform QR decomposition to obtain S^{*}, V^{*n*+1}; Construct $\Xi_{v_k}^{n+1}$ and Γ^{n+1} ;
- Integrate S step with step size Δt to get S^{**} with Δt ;
- Construct $K^{\star\star} = X^n S^{\star\star}$;
- Integrate K step with step size Δt to get K^{n+1} ;
- Perform QR decomposition to obtain X^{n+1} , S^{n+1} ;

Output: $X^{n+1}, S^{n+1}, and V^{n+1}$;

We do the L step first because it drives our system into equilibrium.

Theorem (Implicit Euler integrator)

Under technical assumptions there is a constant C independent of ϵ so that:

$$\|f^{n+1}-\rho_0^{n+1}e^{\top}\|_2\leq C\epsilon\,,$$

where ρ_0^{n+1} solves $\frac{\rho_0^{n+1} - \rho_0^{\star\star}}{\Delta t} = \frac{1}{d} \sum_{k=1}^d \mathsf{D}_k \left(\Sigma^{-1} \mathsf{D}_k \rho_0^{n+1} \right)$ and $\|\rho_0^{\star\star} - \rho_0^n\|_2 = O\left(\frac{(\Delta t)^2}{(\Delta x)^4}\right).$ (1)

The good: Captures the correct low-rank structure and correct discretization of limit equation.

The bad: There is a strong constraint on $\Delta t \ll (\Delta x)^2$ that makes the scheme unfeasible in practice.

Proof sketch

All quantities of interested are expanded using the ansatz

$$p=p_0+\epsilon p_1+\epsilon^2 p_2+\cdots$$
.

We have three steps in the projector splitting integrator (equations for L, S, and K).

L step: Asymptotic expansion

$$\begin{cases} \mathcal{O}(1/\epsilon^2): & \mathsf{A}_\sigma^n \mathsf{L}_0^*\mathsf{C} = 0 \,, \\ \mathcal{O}(1/\epsilon): & \sum_{k=1}^d \mathsf{A}_{\partial_k}^n \mathsf{L}_0^* \mathsf{\Pi}_{v_k} = \mathsf{A}_\sigma^n \mathsf{L}_1^*\mathsf{C} \,, \\ \mathcal{O}(1): & \frac{\mathsf{L}_0^* - \mathsf{L}_0^n}{\Delta t} + \sum_{k=1}^d \mathsf{A}_{\partial_k}^n \mathsf{L}_1^* \mathsf{\Pi}_{v_k} = \mathsf{A}_\sigma^n \mathsf{L}_2^*\mathsf{C} \,. \end{cases}$$

This gives the representation of the equilibrium

$$\mathsf{L}_0^{\star} = \mathit{I}_0^{\star} \mathit{e}_{\mathrm{n}}^{\top}.$$

Proof sketch

K step: Asymptotic expansion yields (with $\alpha^{n+1} = (V^{n+1})^{\top} e_n$)

$$\frac{\mathsf{K}_{0}^{n+1}\alpha^{n+1} - \mathsf{K}_{0}^{\star\star}\alpha^{n+1}}{\Delta t} - \frac{1}{d}\sum_{k=1}^{d}\mathsf{D}_{k}\mathsf{\Sigma}^{-1}\mathsf{D}_{k}\mathsf{K}_{0}^{n+1}\alpha^{n+1} = 0$$

and thus

$$\frac{\rho_0^{n+1} - \rho_0^{\star\star}}{\Delta t} - \frac{1}{d} \sum_{k=1}^d \mathsf{D}_k \left(\Sigma^{-1} \mathsf{D}_k \rho_0^{n+1} \right) = 0.$$

S step together with the L step is a complication (ideally $\rho_0^{\star\star} = \rho_0^n$). But

$$\rho_0^{\star\star} = \left(\mathsf{I} + \frac{\Delta t}{d}\mathcal{L}\right)^{-1} \left(\mathsf{I} - \frac{\Delta t}{d}\mathcal{L}\right)^{-1} \rho_0^n = \left(\mathsf{I} - \frac{(\Delta t)^2}{d^2}\mathcal{L}^2\right)^{-1} \rho_0^n,$$

where $\mathcal{L} = \sum_{k=1}^{d} X^{n} A_{\partial_{k}}^{n} (A_{\sigma}^{n})^{-1} A_{\partial_{k}}^{n} (X^{n})^{\top}$.

Input: X^n , S^n and V^n . **Algorithm:**

- Construct $A^n_{\partial_k}, A^n_{\sigma}$; Compute $L^n = V^n(S^n)^\top$;
- Integrate L step with step size Δt to get L^{*} using Crank-Nicolson;
- ▶ Perform QR decomposition to obtain S^{*}, V^{*n*+1}; Construct $\Xi_{\nu_k}^{n+1}$ and Γ^{n+1} ;
- Integrate S step with step size Δt to get S^{**} with Δt using Crank-Nicolson;
- Construct $K^{\star\star} = X^n S^{\star\star}$;
- Integrate K step with step size Δt to get K^{*n*+1};
- Perform QR decomposition to obtain X^{n+1} , S^{n+1} ;

Output: X^{n+1} , S^{n+1} , and V^{n+1} ;

Error analysis

Crank-Nicolson for the L step:

$$\frac{\mathsf{L}^{\star}-\mathsf{L}^{n}}{\Delta t}+\frac{1}{\epsilon}\sum_{k=1}^{d}\mathsf{A}_{\partial_{k}}^{n}\left(\frac{\mathsf{L}^{\star}+\mathsf{L}^{n}}{2}\right)\mathsf{\Pi}_{v_{k}}=\frac{\mathsf{A}_{\sigma}^{n}}{\epsilon^{2}}\frac{\mathsf{L}^{\star}+\mathsf{L}^{n}}{2}\mathsf{C}.$$

Theorem (CNIE integrator)

Assuming that $f^n =
ho_0^n e^\top + \mathcal{O}(\epsilon)$ we get

$$\|\rho^{\star\star}-\rho^n\|_2=O(\epsilon)$$

and

$$\frac{\rho_0^{n+1} - \rho_0^{\star\star}}{\Delta t} - \frac{1}{d} \sum_{k=1}^d \mathsf{D}_k \left(\Sigma^{-1} \mathsf{D}_k \rho_0^{n+1} \right) = 0.$$

Further, we have $f^{n+1} = \rho_0^{n+1} e^\top + \mathcal{O}(\epsilon)$.

The low-rank limit is obtained independent of the time step size.

Much of the proof proceeds analogous to the implicit Euler case.

But, because of the Crank-Nicolson scheme in the L and S step we now have

$$\begin{split} \rho_0^{\star} &= \left(\mathsf{I} - \frac{\Delta t}{2d} \mathcal{L} \right)^{-1} \left(\mathsf{I} + \frac{\Delta t}{2d} \mathcal{L} \right) \rho_0^n, \\ \rho_0^{\star \star} &= \left(\mathsf{I} + \frac{\Delta t}{2d} \mathcal{L} \right)^{-1} \left(\mathsf{I} - \frac{\Delta t}{2d} \mathcal{L} \right) \rho_0^{\star}, \end{split}$$

where as before $\mathcal{L} = \sum_{k=1}^{d} X^n A^n_{\partial_k} (A^n_{\sigma})^{-1} A^n_{\partial_k} (X^n)^{\top}$.

▶ Note that since the S step (second line) is backward in time the signs match up.

Combining these two equations we get $\rho_0^{\star\star} = \rho_0^n$.

Hilbert expansion as discussed here is a formal method.

Proofs can be made rigorous by the techniques in [C. Bardos, R. Santos, and R. Sentis, Trans. Amer. Math. Soc., 284 (1984)] Efficient asymptotic preserving schemes

The previously discussed numerical method are fully implicit.

Requires the solution of a large linear systems in every time step.

► Very expensive!

Macro-micro decomposition

$$f(t, x, v) = \rho(t, x) + \varepsilon g(t, x, v).$$

Evolution equation for ρ and g

$$\partial_t \rho + \frac{1}{4\pi} \nabla_x \cdot \langle \mathbf{v} g \rangle_{\mathbf{v}} = -\sigma^A \rho + G$$
$$\partial_t g + \frac{1}{\varepsilon} \left(I - \frac{1}{4\pi} \langle \rangle_{\mathbf{v}} \right) \left(\mathbf{v} \cdot \nabla_x g \right) + \frac{1}{\varepsilon^2} \mathbf{v} \cdot \nabla_x \rho = -\frac{\sigma^S}{\varepsilon^2} g - \sigma^A g.$$

For $\epsilon \rightarrow 0$ we have

$$g = -\frac{1}{\sigma^S} \mathbf{v} \cdot \nabla_{\mathbf{x}} \rho,$$

which is at most rank d.

We now approximate g using a low-rank representation.

M. Lemou, L. Mieussens, SIAM J. Sci. Comput., 31:334-368, 2008.

Evolution equation for K

$$\partial_{t} \mathcal{K}_{j} = \langle V_{j}, \mathsf{RHS} \rangle_{\nu} \\ = -\frac{1}{\varepsilon} \sum_{l=1}^{r} \left(\langle v V_{j} V_{l} \rangle_{\nu} - \frac{1}{4\pi} \langle V_{j} \rangle_{\nu} \langle v V_{l} \rangle_{\nu} \right) \cdot \nabla_{x} \mathcal{K}_{l} - \frac{1}{\varepsilon^{2}} \langle v V_{j} \rangle_{\nu} \cdot \nabla_{x} \rho - \left(\frac{\sigma^{S}}{\varepsilon^{2}} + \sigma^{A} \right) \mathcal{K}_{j}.$$

First order IMEX discretization

$$\frac{\mathcal{K}_{j}^{n+1} - \mathcal{K}_{j}^{n}}{\Delta t} = -\frac{1}{\varepsilon} \sum_{l=1}^{r} \left(\langle v V_{j}^{n} V_{l}^{n} \rangle_{v} - \frac{1}{4\pi} \langle V_{j}^{n} \rangle_{v} \langle v V_{l}^{n} \rangle_{v} \right) \cdot \nabla_{x} \mathcal{K}_{l}^{n} - \frac{1}{\varepsilon^{2}} \left(\langle v V_{j}^{n} \rangle_{v} \cdot \nabla_{x} \rho^{n} + \sigma^{S} \mathcal{K}_{j}^{n+1} \right) - \sigma^{A} \mathcal{K}_{j}^{n}.$$

is AP and explicit.

Second order

We have the macro-micro equations

 $\partial_t \rho = F(\rho, g), \qquad \partial_t g = G(g, \rho).$

First order can be easily obtained by substituting the following set of equations

$$\partial_t \rho = F(\rho, g^n), \qquad \partial_t g = G(g, \rho^n).$$

For second order we can exploit symmetry. If $g^{n+1/2}$ and $\rho^{n+1/2}$ are first order approximations to $g(t^n + \Delta t/2)$ and $\rho(t^n + \Delta t/2)$ then

$$\partial_t \rho = F(\rho, g^{n+1/2}), \qquad \partial_t g = G(g, \rho^{n+1/2}).$$

is second order accurate (assuming the equations are solved up to second order).

The following method is second order accurate and AP

- Compute $\rho^{n+1/2}$ from ρ^n using g^n (first order is sufficient).
- Solve $\partial_t g = G(g, \rho^{n+1/2})$ using projector splitting and IMEX2.
- Compute ρ^{n+1} from ρ^n using $g(t^{n+1/2})$ (using a second order scheme).

Can be generalized to higher order (almost-symmetric splitting).

L.E., A. Ostermann, Comput. Appl. Math., 271, 2014. L.E., A. Ostermann, Comput. Math. Appl., 67(12), 2014.

Two-material test problem





Two-material test problem





Kinetic equations with a fluid limit

Fluid limit

Collisional kinetic equation with a fluid limit $\partial_t f(t, x, v) + v \cdot \nabla_x f(t, x, v) = \frac{1}{\epsilon} (M(f) - f) \xrightarrow{\epsilon \to 0}$ Euler equations

with

$$M = \frac{\rho}{(2\pi)^{d_v/2}} \exp\left(-\frac{|v-u|^2}{2}\right), \qquad \rho = \int f \,\mathrm{d}v, \qquad u = \frac{1}{\rho} \int v f \,\mathrm{d}v.$$

Chapman–Enskog theory: $f = M(1 + \epsilon f_1 + \epsilon^2 f_2)$

Density and momentum satisfy the (isothermal) Navier-Stokes equation

$$\partial_t \rho + \nabla_x \cdot (\rho u) = 0,$$

$$\partial_t (\rho u) + \nabla \cdot (\rho u \otimes u) + \nabla \rho = \nabla \cdot \left[\mu \left(\nabla u + (\nabla u)^{\mathrm{T}} \right) + \lambda (\nabla \cdot u) I \right].$$

C. Bardos, F. Golse, and D. Levermore. J. Stat. Phys, 63(1-2), 1991.
C. Bardos, F. Golse, and C.D. Levermore. Commun. Pure Appl. Math, 46(5), 1993.

For $\epsilon \rightarrow 0$ we have

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

Dynamics completely determined by the moments ρ and u (fluid regime).

But f is not low-rank.

Interlude: Lattice Boltzmann methods

At least two strategies to solve fluid problems.

- 1. Directly discretize the Navier–Stokes equations.
- 2. Discretize the Boltzmann equation

$$\partial_t f + \mathbf{v} \cdot \nabla_x f = \frac{1}{\epsilon} (M(f) - f)$$

with a very coarse velocity discretization

$$f_j(t,x) \approx f_j(t,x,e_j).$$

Then reconstruct the quantities of interest

$$\rho(t,x) = \int f(t,x,v) \,\mathrm{d}v \approx \sum_j \omega_j f_j(t,x).$$



Low-rank vs Lattice Boltzmann

Why throw away all the information about velocity space?

• Even in a fluid problem this is still important.

Apply the low-rank algorithm with initial value

$$f(0, x, v) = \frac{\rho(0, x)}{(2\pi)^{d_v/2}} \exp\left(-\frac{1}{2}(v - u(0, x))^2\right).$$

Advantages

► For weakly compressible flow we can use (rank 10 in 3d)

$$\frac{\rho}{(2\pi)^{d/2}}\exp\left(-\frac{v^2}{2}\right)\left(1+v\cdot u+\frac{(v\cdot u)^2}{2}-\frac{u^2}{2}\right)+\mathcal{O}\left(u^3\right).$$

- Spectral methods can be incorporated easily.
- ► Straightforward to capture some kinetic effects.
- Additional cost is small.

Low-rank approximation

Evolution equation for K

$$\partial_t K_j(t,x) = -\sum_l c_{jl}^1 \cdot \nabla_x K_l(t,x) + \frac{1}{\epsilon} \left(K_j(t,x) - c_j^3(K)(t,x)\rho(K)(t,x) \right)$$

In the limit $\epsilon \rightarrow 0$ we have

$$K_j - c_j^3(K)(x)
ho(K)(x) = 0$$

which can be written as

$$K_j = \frac{\rho(K)}{(2\pi)^{d/2}} \int V_j(v) \exp\left(-\frac{1}{2}(v-u(K))^2\right) \,\mathrm{d}v.$$

This is simply the projection of the Maxwell–Boltzmann distribution onto the space spanned by the V_i .

Sound waves

Propagation of sound waves

- Fluid solver with $\tau = 7 \cdot 10^{-3}$ (CFL number of 0.9).
- Lie splitting with $\tau = 0.1$ and Strang splitting with $\tau = 0.2$.







The compressible case

Low-rank structure

Is there a low-rank structure in the compressible case?

► Note that equilibrium is uniquely defined by its moments.

In the macro-micro decomposition we have $f = M + \varepsilon f_1$. Plugging this into

$$\partial_t f + \mathbf{v} \cdot \nabla_x f = \frac{\nu}{\varepsilon} (M - f).$$

and using $(M-f)/\varepsilon = f_1$ we get

$$f_1 = -rac{1}{
u}(\partial_t f + v\cdot
abla_{ imes} f) = -rac{1}{
u}(\partial_t M + v\cdot
abla_{ imes} M) + \mathcal{O}(arepsilon).$$

Almost a low-rank structure

$$\frac{1}{M}(\partial_t M + v \cdot \nabla_x M)$$

= $\frac{1}{\rho}(\partial_t \rho + v \cdot \nabla_x \rho) + (v - u) \cdot (\partial_t u + v \cdot \nabla_x u).$

The micro-macro decomposition fails because $f_1 = M(\text{low-rank})$.

We will instead use the **multiplicate decompositon** f = Mg.

Then

$$g = 1 - \frac{\varepsilon}{\nu} \left[\left((v - u) \otimes (v - u) - \frac{|v - u|^2}{d_v} I \right) : \nabla_x u \right] + \mathcal{O}(\varepsilon^2),$$

where $A : B = \sum_{ij} A_{ij} B_{ij}.$

g is low-rank up to at least $\mathcal{O}(\epsilon^2)$.

We have

$$f = Mg,$$
 $g = \sum_{ij} X_i S_{ij} V_j.$

Evolution of the moments to obtain *M*

$$\partial_t U + \nabla_x \cdot \langle v \phi Mg \rangle_v = 0, \quad U = (\rho, \rho u)^T, \quad \phi(v) = (1, v)^T.$$

Treated as any other conservation law.

Evolution of g

$$\partial_t g = -v \cdot \nabla_x g - \frac{1}{M} (\partial_t M + v \cdot \nabla_x M) g + \frac{\nu}{\varepsilon} (1 - g)$$

Treated by dynamical low-rank.

Challenges

For **DLR** this moves the problem into the **coefficients**. E.g.

$$\langle vV_j^n M^n \rangle_v = \frac{\rho^n(x)}{(2\pi)^{d_v/2}} \left\langle vV_j^n(v) \exp\left(-\frac{|v-u^n(x)|^2}{2}\right) \right\rangle_v.$$
$$\langle (v \otimes v)V_j^n M^n \rangle_v = \frac{\rho^n(x)}{(2\pi)^{d_v/2}} \left\langle (v \otimes v)V_j^n(v) \exp\left(-\frac{|v-u^n(x)|^2}{2}\right) \right\rangle_v.$$

Can be treated by fast convolution algorithms

$$\begin{split} g_j^1 &= (v \mapsto vV_j^n) * (v \mapsto \exp(-v^2/2)), & \text{evaluated at} & u^n(x). \\ g_j^2 &= (v \mapsto (v \otimes v)V_j^n) * (v \mapsto \exp(-v^2/2)), & \text{evaluated at} & u^n(x). \end{split}$$

To compute the convolution $h_1 * h_2$ of $h_1(v)$ and $h_2(v)$ we proceed as follows.

Step 1: Compute \hat{h}_1 and \hat{h}_2 by using a FFT.

- Cost: $\mathcal{O}(n^{d_v} \log n^{d_v})$.
- **Step 2:** Compute $g = \mathcal{F}^{-1}(\hat{h}_1 \hat{h}_2)$. \blacktriangleright Cost: $\mathcal{O}(n^{d_v} \log n^{d_v})$.

Step 3: Interpolate g (e.g. using cubic spline interpolation) and evaluate $g(u(x_i))$ for each grid point x_i

• Cost:
$$\mathcal{O}(n^{d_x})$$
.

In the Euler limit ($\epsilon \rightarrow$ 0) shock waves are known to develop.

• Discontinuous solutions or for $\epsilon > 0$ sharp gradients in the solution.

It is well known that standard methods do not work in this case.

This is a significant complication compared to the diffusion limit.

K equation

Discretizing the K equation

$$\partial_t \mathcal{K}_j = -\sum_l (\nabla_x \mathcal{K}_l) \cdot \langle v \mathcal{V}_j \mathcal{V}_l \rangle_v - \sum_l \mathcal{K}_l \langle \mathcal{V}_j \mathcal{V}_l \mathcal{M} \rangle_v + \frac{\rho}{\varepsilon} (\langle \mathcal{V}_j \rangle_v - \mathcal{K}_j)$$

using a first order IMEX scheme

$$\mathcal{K}_{j}^{n+1} = \frac{1}{1 + \Delta t \rho^{n} / \varepsilon} \mathcal{K}_{j}^{n} - \frac{\Delta t}{1 + \Delta t \rho^{n} / \varepsilon} \left[\sum_{l} c_{jl}^{1} \cdot (\nabla_{\mathsf{x}} \mathcal{K}_{l}^{n}) + \sum_{l} c_{jl}^{2} \mathcal{K}_{l}^{n} \right] + \frac{\Delta t \rho^{n}}{\varepsilon + \Delta t \rho^{n}} \langle V_{j} \rangle_{\mathsf{v}}.$$

In 2D $c_{jl}^1 = [c_{jl}^{1;1} c_{jl}^{1;2}]^T$ and the matrices are symmetric. Thus, there exist orthogonal matrices T^m such that $\sum_{jl} T_{ij}^m c_{jl}^{1;m} T_{kl}^m = \lambda_i^m \delta_{ik}$.

K equation

Using
$$\hat{K}_{i}^{n} = \sum_{j} T_{ij}^{1} K_{j}^{n}$$
 we get

$$\hat{K}_{i}^{n+1} = \frac{1}{1 + \Delta t \rho^{n}/\varepsilon} \hat{K}_{i}^{n} - \frac{\Delta t}{1 + \Delta t \rho^{n}/\varepsilon} \left[\lambda_{i}^{1} \partial_{\mathbf{x}} \hat{K}_{i}^{n} + \sum_{jl} T_{ij}^{1} c_{jl}^{1;2} \partial_{\mathbf{y}} K_{l}^{n} + \sum_{lj} T_{ij}^{1} c_{jl}^{2} K_{l}^{n} \right] + \dots$$

Direction of the flow is now obvious.

Replace $\lambda_i^1 \partial_x \hat{K}_i^n$ by an appropriate discrete approximation, which we denote by $\delta_x(K_i^n, \lambda_i^1)$.

$$\mathcal{K}_{j}^{n+1} = \frac{1}{1 + \Delta t \rho^{n}/\varepsilon} \mathcal{K}_{j}^{n} - \frac{\Delta t}{1 + \Delta t \rho^{n}/\varepsilon} \left[\sum_{i} \mathcal{T}_{ij}^{1} \delta_{x}(\hat{\mathcal{K}}_{i}^{n}, \lambda_{i}^{1}) + \sum_{l} c_{jl}^{1;2} \partial_{y} \mathcal{K}_{l}^{n} + \sum_{l} c_{jl}^{2} \mathcal{K}_{l}^{n} \right] + \dots$$

Examples: upwinding, Lax-Wendroff flux with van Leer limiter, ...

For the moments $U = (\rho, \rho u)^T$ we have

$$\partial_t U + \nabla_x \cdot \langle v \phi M g \rangle_v = 0, \qquad \phi(v) = (1, v)^T.$$

This is in the form of a conservation flow, but the flux depends on g.

- Classic methods for conservation flow solve a Riemann problem.
- ► This is difficult to do here.

We consider the one-dimensional conservation law

```
\partial_t U(t,x) + \partial_x F(U(t,x)) = 0.
```

Lax-Friedrichs method

$$U_{j}^{n+1} = \frac{1}{2} \left(U_{j+1}^{n} + U_{j-1}^{n} \right) - \frac{\Delta t}{2\Delta x} \left(F(U_{j+1}^{n}) - F(U_{j-1}^{n}) \right)$$

is stable under the usual CFL condition.

Simplest example of a central scheme.

► Stable scheme that does not require information on the direction of the flow.

Nessyahu-Tadmor

The **Nessyahu–Tadmor** scheme is a second-order (away from discontinuities) central scheme in predictor-corrector form on a staggered grid

$$U_{j}^{\star} = U_{j}^{n} - \frac{\Delta t}{2\Delta x}F_{j}^{\prime},$$

$$U_{j+1/2}^{n+1} = \frac{1}{2}(U_{j}^{n} + U_{j+1}^{n}) + \frac{1}{8}(U_{j}^{\prime} - U_{j+1}^{\prime}) - \frac{\Delta t}{\Delta x}(F(U_{j+1}^{\star}) - F(U_{j}^{\star})).$$

The choice of F'_i and U'_i is free as long as

$$\blacktriangleright F'_j/\Delta x = \partial_x F(U(t^n, x_j)) + \mathcal{O}(\Delta x)$$

$$\blacktriangleright U'_j/\Delta x = \partial_x U(t^n, x_j) + \mathcal{O}(\Delta x).$$

To obtain non-oscillatory solutions that preserve sharp gradient we can choose

$$U'_{j} = \mathsf{MM}(U^{n}_{j+1} - U^{n}_{j}, U^{n}_{j} - U^{n}_{j-1}), \qquad F'_{j} = \mathsf{MM}(F(U^{n}_{j+1}) - F(U^{n}_{j}), F(U^{n}_{j}) - F(U^{n}_{j-1}))$$

with MM the usual minmod limiter.

Generalization to 2D

$$\begin{split} U_{ij}^{\star} &= U_{ij}^{n} - \frac{\Delta t}{2\Delta x} F_{ij}^{\prime \times} - \frac{\Delta t}{2\Delta y} G_{ij}^{\prime y}, \\ U_{i+1/2,j+1/2}^{n+1} &= \frac{1}{4} (U_{ij}^{n} + U_{i+1,j}^{n} + U_{i,j+1}^{n} + U_{i+1,j+1}^{n}) \\ &+ \frac{1}{16} (U_{ij}^{\prime \times} - U_{i+1,j}^{\prime \times} + U_{i,j+1}^{\prime \times} - U_{i+1,j+1}^{\prime \times}) \\ &+ \frac{1}{16} (U_{ij}^{\prime y} - U_{i,j+1}^{\prime y} + U_{i+1,j}^{\prime y} - U_{i+1,j+1}^{\prime y}) \\ &+ \frac{\Delta t}{\Delta x} (F(U_{i+1,j}^{\star}) - F(U_{ij}^{\star})) + \frac{\Delta t}{\Delta y} (F(U_{i,j+1}^{\star}) - F(U_{ij}^{\star})), \end{split}$$

where

$$\begin{aligned} U_{ij}^{\prime x} &= \mathsf{MM}(U_{i+1,j}^n - U_{i,j}^n, U_{ij}^n - U_{i-1,j}^n), \qquad U_{ij}^{\prime y} = \mathsf{MM}(U_{i,j+1}^n - U_{ij}^n, U_{ij}^n - U_{i,j-1}^n), \\ F_{ij}^{\prime x} &= \mathsf{MM}(F(U_{i+1,j}^n) - F(U_{i,j}^n), F(U_{ij}^n) - F(U_{i-1,j}^n)), \\ F_{ij}^{\prime y} &= \mathsf{MM}(F(U_{i,j+1}^n) - F(U_{ij}^n), F(U_{ij}^n) - F(U_{i,j-1}^n)). \end{aligned}$$

In our case we have

$$U = \begin{bmatrix} \rho \\ \rho u \end{bmatrix}, \qquad F(U) = \begin{bmatrix} \sum_{ij} X_i^n S_{ij}^n \langle v V_j^n M^n \rangle_v \\ \sum_{ij} X_i^n S_{ij}^n \langle (v \otimes v) V_j^n M^n \rangle_v \end{bmatrix}$$

To compute the flux we use the fast convolution algorithm.

We found that the staggered variant is more accurate

- 1. Apply the Nessyahu–Tadmor scheme with time step size $\Delta t/2$ and initial value U_{ij}^{n} to obtain $U_{i+1/2,j+1/2}^{n+1/2}$.
- 2. Obtain X^n at the half grid points (i + 1/2, j + 1/2) by computing averages between neighboring grid points.
- 3. Apply the Nessyahu–Tadmor scheme with time step size $\Delta t/2$ and initial value $U_{i+1/2,j+1/2}^{n+1/2}$ and the X^n at the half grid points to obtain $U_{i,j}^{n+1}$.

Shear flow (fluid solver)



Shear flow (dynamical low-rank)

t=12



Error

Error in the moments for different rank r and space discretizations



Shear flow with large Reynolds number





Literature

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► AP dynamical low-rank scheme for the compressible Navier–Stokes limit.