

Structure preserving numerical methods for the Vlasov equation

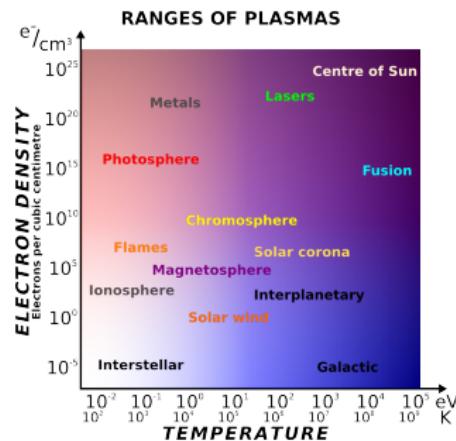
Lukas Einkemmer
University of Innsbruck

Oberwolfach, 2016

Vlasov equation

Main simplifying assumption

- ▶ collisionless plasma
- ▶ described by a (classic) particle density function



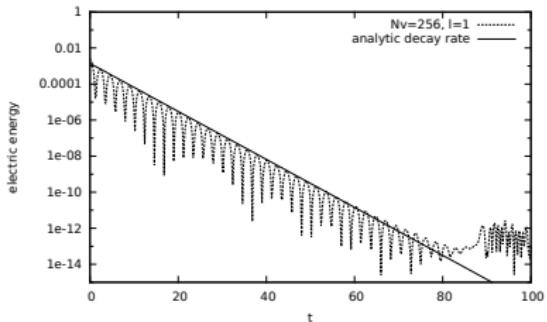
Kinetic equation describes evolution of particle density f

$$\partial_t f(t, x, v) + v \cdot \nabla_x f(t, x, v) + F \cdot \nabla_v f(t, x, v) = 0$$
$$F = q(E + v \times B),$$

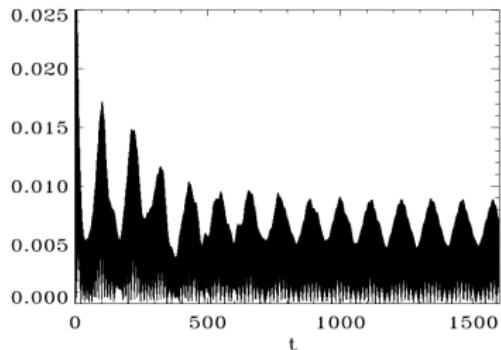
coupled to Gauss's law or Maxwell's equations.

Applications include plasma physics, radiative heat transfer, astrophysics.

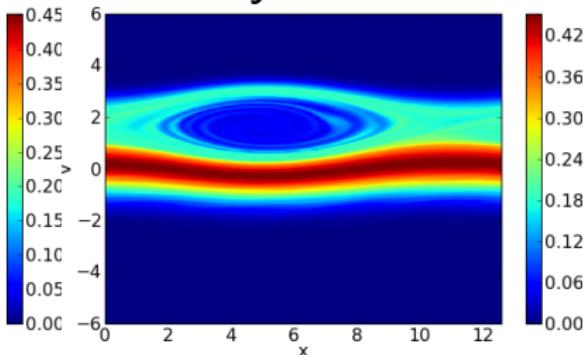
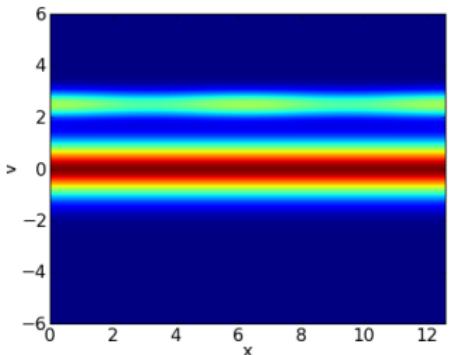
Landau damping



Nonlinear Landau damping and BGK equilibria



Two-stream instability



Time splitting for the Vlasov–Poisson equations

Free streaming part

$$\partial_t f(t, \mathbf{x}, \mathbf{v}) + \mathbf{v} \cdot \nabla_{\mathbf{x}} f(t, \mathbf{x}, \mathbf{v}) = 0$$

Solution:

$$f(\tau, \mathbf{x}, \mathbf{v}) = f(0, \mathbf{x} - \tau \mathbf{v}, \mathbf{v})$$

Acceleration part

$$\partial_t f(t, \mathbf{x}, \mathbf{v}) + \mathbf{F} \cdot \nabla_{\mathbf{v}} f(t, \mathbf{x}, \mathbf{v}) = 0$$

Solution for constant \mathbf{F} :

$$f(\tau, \mathbf{x}, \mathbf{v}) = f(0, \mathbf{x}, \mathbf{v} - \tau \mathbf{F})$$

But: \mathbf{F} is not constant – nonlinear equation

Idea: Freeze force term at half the time step, i.e. at $\tau/2$.

- ▶ only translations have to be computed

Generalized to the Vlasov–Maxwell equations.

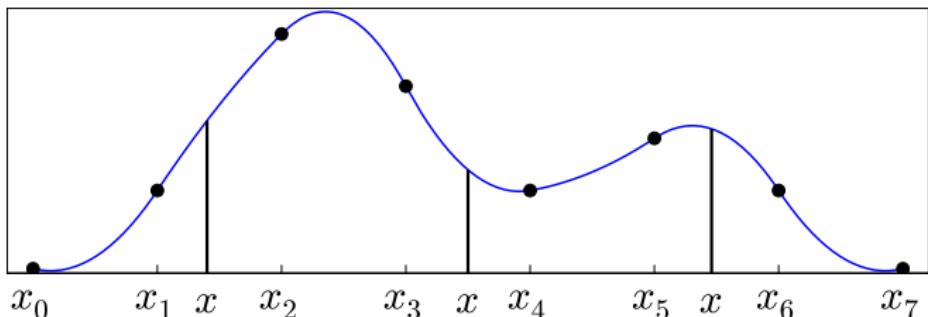
Cubic spline interpolation

Follow the **characteristics** backward in time

- ▶ semi-Lagrangian approach

The feet of the characteristics do not necessarily coincide with the numerical grid

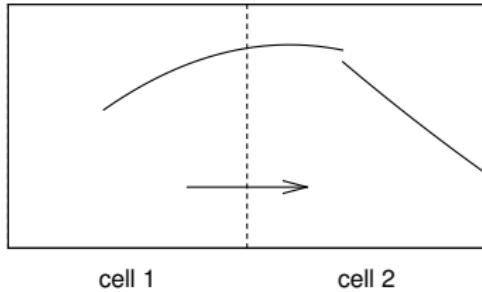
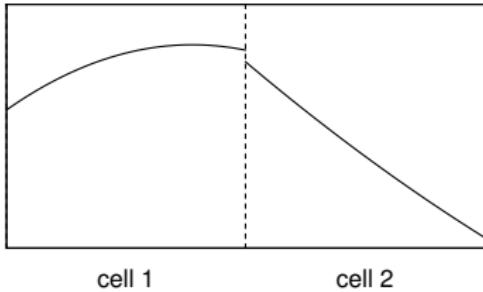
- ▶ Interpolation using **cubic splines**



Results in a **mass conservative** scheme.

Discontinuous Galerkin semi-Lagrangian method

Idea: translation and projection to a subspace of piecewise polynomials (without any continuity constraint).



Yields a **conservative**, **local**, and only **slightly-diffusive** scheme.

Well suited for **parallelization**.

J.A. Rossmanith, D.C. Seal, J. Comput. Phys. (2011)

J.M. Qiu, C.W. Shu, J. Comput. Phys. (2011)

N. Crouseilles, M. Mehrenberger, F. Vecil, ESAIM: Proceedings (2011)

L. Einkemmer, A. Ostermann, SIAM J. Numer. Anal. (2014)

Weak scaling for Vlasov–Poisson

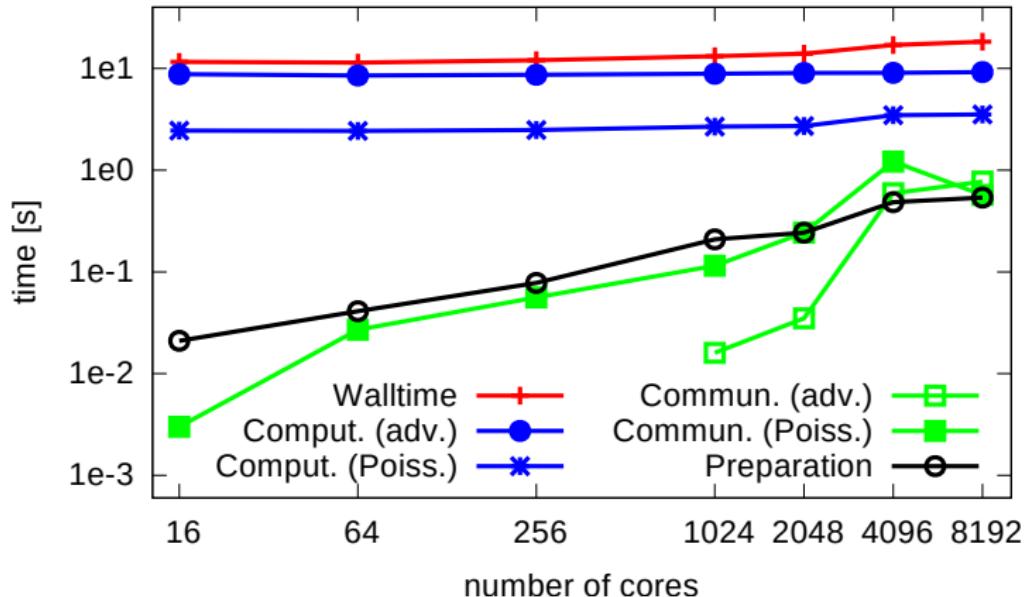


Figure: Weak scaling for the two-dimensional Vlasov–Poisson equations on the VSC-2.

What structure to preserve

Infinite number of conserved quantities for Vlasov–Poisson

- ▶ charge, all L^p norms, current, energy, entropy, ...

Focus here on the **physical relevant quantities**

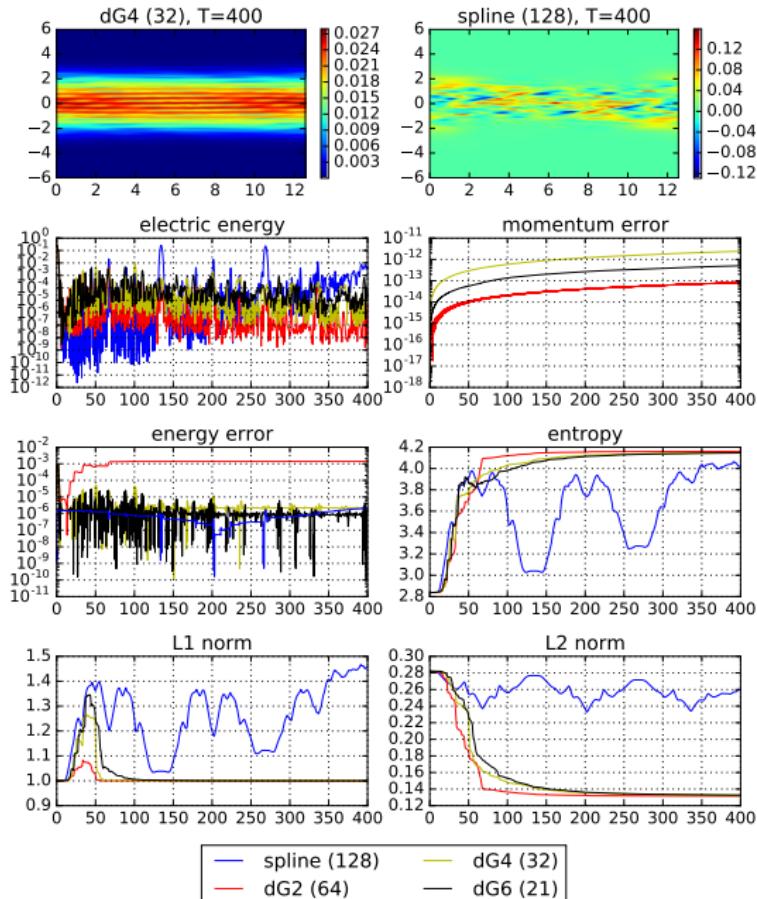
- ▶ charge, current, energy

and two measures of dissipation

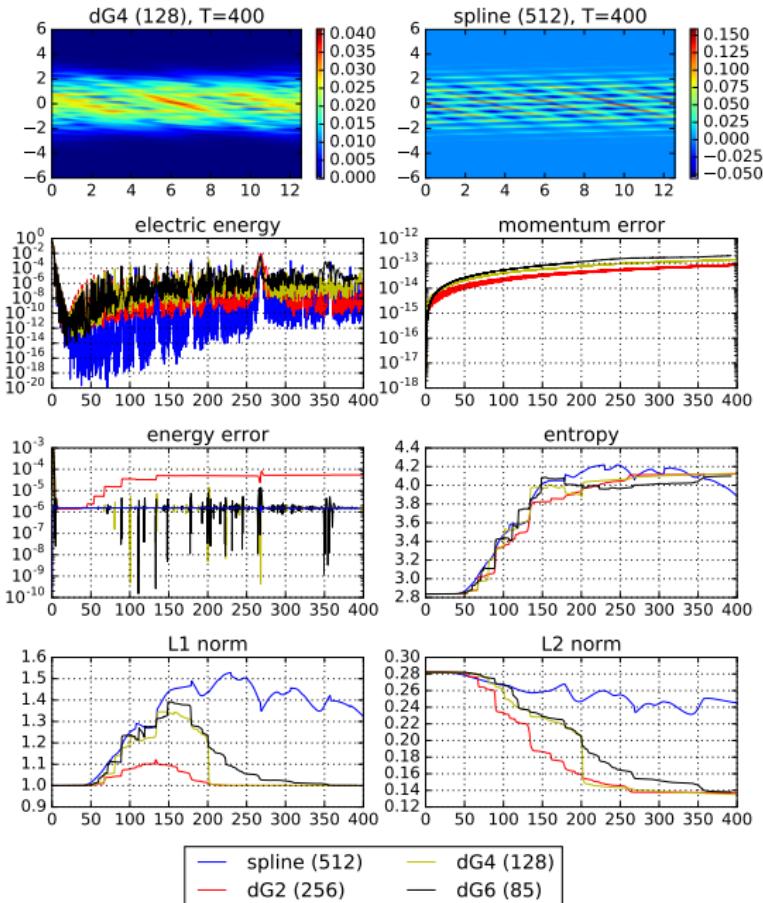
- ▶ $\|f\|_2^2$, entropy

conserved var.	splitting	dG	spline	remarks
charge	yes	yes	yes	
current	yes	yes [†]	yes [†]	additional term for $p = 0$
energy	no	no	no	additional term for $p = 0, p = 1$
entropy	yes	no	no	
L^1	yes	no*	no	
L^2	yes	no	no	

Expansion – low resolution



Expansion – high resolution



Dissipation

A common way to analyze the conservation of a numerical scheme is to consider a modified equation.

- ▶ Cubic spline interpolation **violates the second law of thermodynamics**
- ▶ and increases the L^2 norm.

Both of these **help to conserve the quantity in question** but are completely unphysical.

L^2 is strictly decreasing for dG

Theorem

The L^2 norm is a decreasing function of time for the semi-Lagrangian discontinuous Galerkin scheme.

Proof.

Let us consider a translated function $g(x)$. Expanding in a Legendre series gives

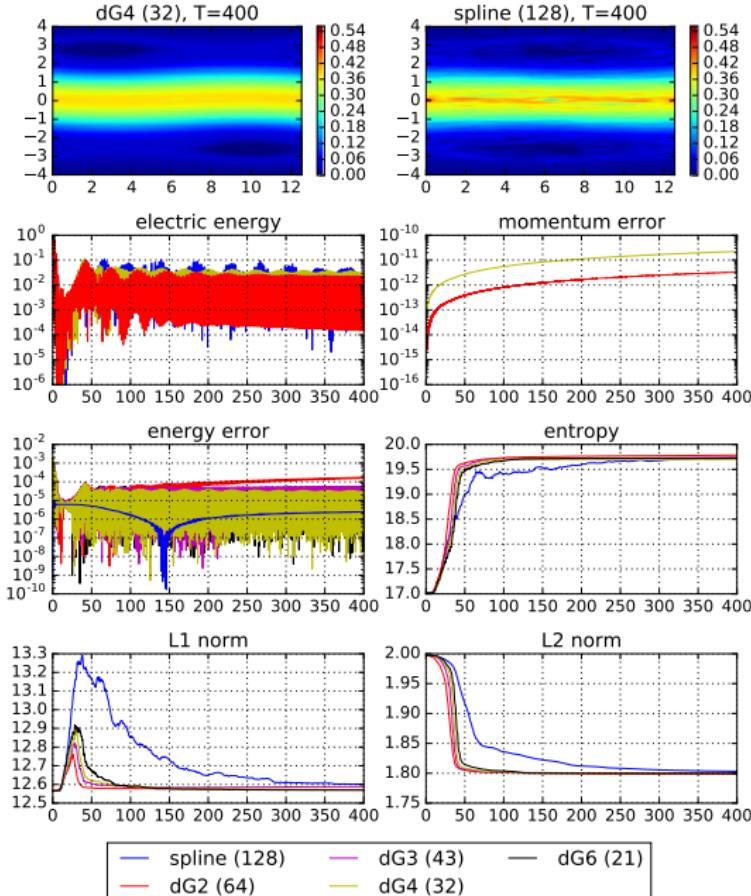
$$g(x) = \sum_{k=0}^{\infty} g_k p_k(x).$$

Therefore

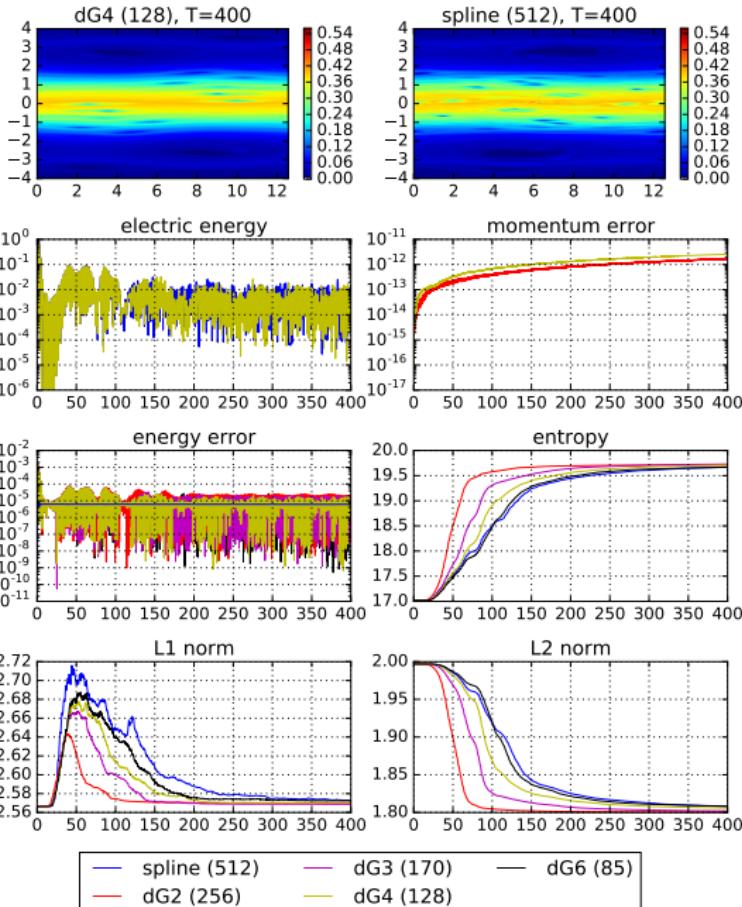
$$\|Pg\|_2^2 - \|g\|_2^2 = \sum_{k=0}^{\ell} g_k^2 - \sum_{k=0}^{\infty} g_k^2 \leq 0.$$



Nonlinear Landau damping – low resolution



Nonlinear Landau damping – high resolution



L^1 conservation

Theorem

A numerical algorithm that produces negative values and is mass conservative can not conserve the L^1 norm.

Proof.

$$\begin{aligned}\int |f(0, x, v)| d(x, v) &= \int f(0, x, v) d(x, v) \\ &= \int f(t, x, v) d(x, v) \\ &\leq \int |f(t, x, v)| d(x, v).\end{aligned}$$

The inequality becomes strict if the particle density function $f(t, x, v)$ is negative. □

To limit or not to limit

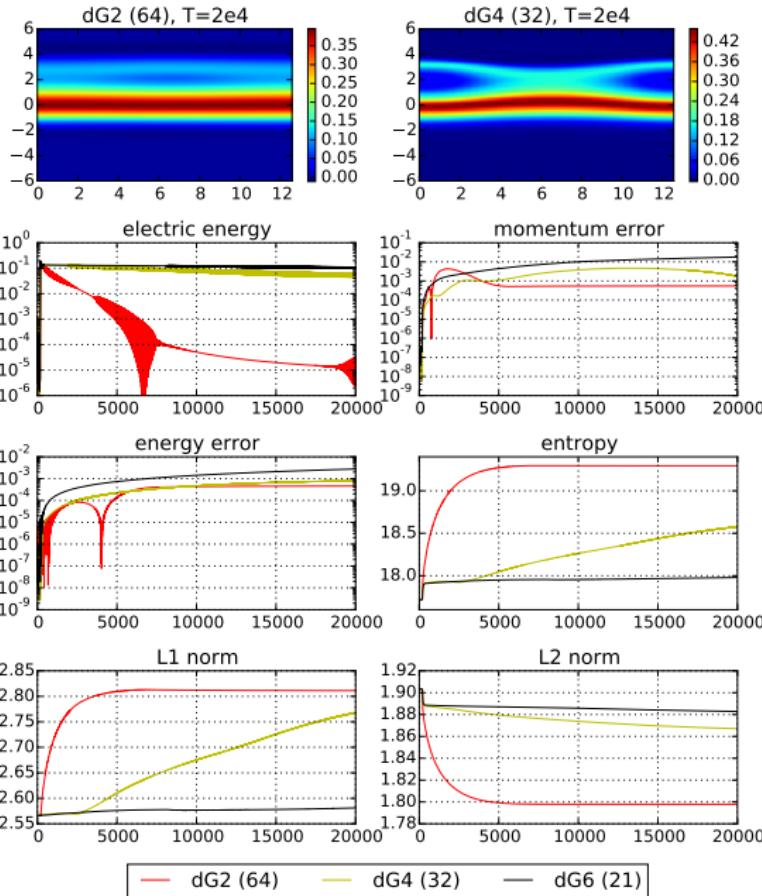
The semi-Lagrangian discontinuous Galerkin scheme can be made L^1 **conservative** by using **positivity limiters**.

But that necessarily introduces **additional dissipation** (or additional overshoot).

Our numerical simulations indicate that positivity is almost always a less severe problem than dissipation

- ▶ this is more of an issue for spline interpolation – positivity preservation is more difficult in that context

Long time behavior



Higher order methods

Suppose we start with

$$f(0, x, v) = e^{-v^2/2}(1 + \alpha \cos kx).$$

and solve $\partial_t f + v \partial_x f = 0$ we get

$$f(t, x, v) = e^{-v^2/2}(1 + \alpha \cos k(x - vt)).$$

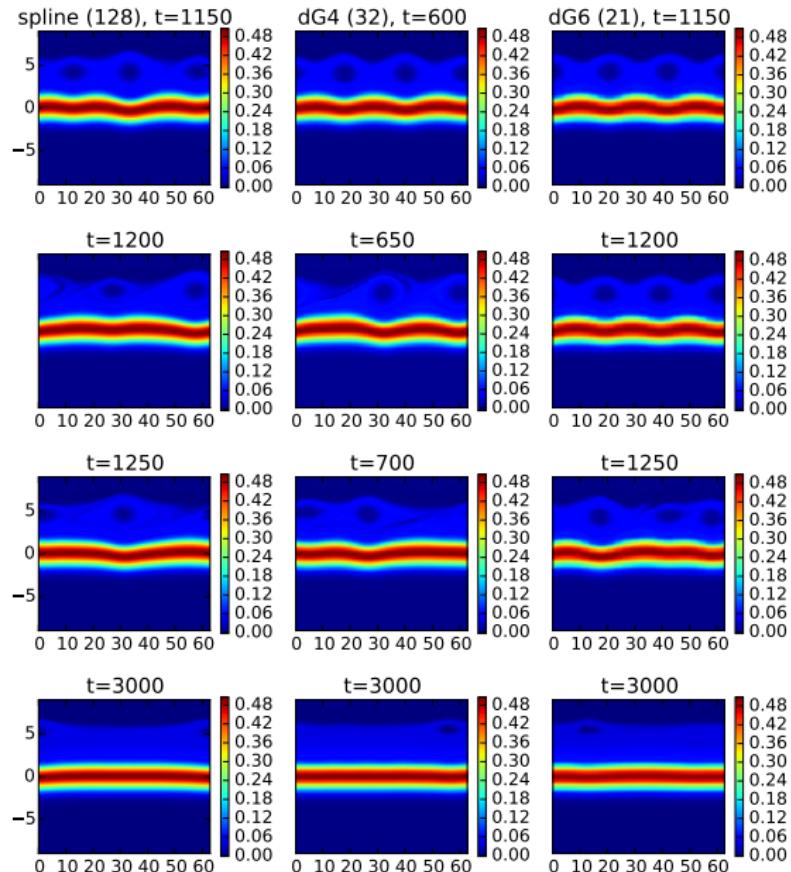
Thus,

$$\partial_v^m f(t, x, v) \propto (kvt)^m.$$

For all intents and purposes $\partial_v^6 f(t, x, v) = \infty$.

- ▶ But the **sixth order method outperforms** the lower order methods.

Bump-on-tail instability



Conservation in the exascale era

On future supercomputers we will most likely see

- ▶ **less memory per core**
- ▶ increased use of accelerators

Single precision would result in a reduction in memory and an increase in performance by **a factor of 2**.

- ▶ Conservation up to single precision?

Mixed-precision algorithm

In each cell we store c_0, \dots, c_p where

$$u(x) = \sum_{k=0}^p u_k p_k(x).$$

Store c_0 in double \Rightarrow Conservation of mass up to double

- ▶ Reduction in memory use by 1.6 (1D), 1.8 (2D), 1.97 (3D).

Intel Xeon Phi				
order	# double	bandwidth	speedup	memorydown
4	4	67.9 GB/s	–	–
4	1	76.3 GB/s	1.80	1.60
4	0	68.6 GB/s	2.02	2.00

NVIDIA K80				
order	# double	bandwidth	speedup	memorydown
4	4	137.9 GB/s	–	–
4	1	130.1 GB/s	1.51	1.60
4	0	142.5 GB/s	2.07	2.00

Conclusion

Thank you for your attention

<http://arxiv.org/abs/1601.02280>

<http://arxiv.org/abs/1603.07008>