

# Conservative and positivity-preserving semi-Lagrangian kinetic schemes with spectrally accurate phase-space resolution

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# Maxwell-Boltzmann system

Maxwell's equations:

$$\begin{aligned}\nabla \times \mathbf{E} &= -\frac{\partial \mathbf{B}}{\partial t} & \nabla \cdot \mathbf{E} &= \frac{\rho}{\varepsilon_0} \\ \nabla \times \mathbf{B} &= \mu_0 \mathbf{J} + \mu_0 \varepsilon_0 \frac{\partial \mathbf{E}}{\partial t} & \nabla \cdot \mathbf{B} &= 0\end{aligned}$$

Sources: charge and current density:

$$\rho(\mathbf{r}, t) = \sum_{\alpha} q_{\alpha} n_{\alpha}(\mathbf{r}, t), \quad \mathbf{J}(\mathbf{r}, t) = \sum_{\alpha} q_{\alpha} n_{\alpha}(\mathbf{r}, t) \mathbf{u}_{\alpha}(\mathbf{r}, t).$$

Number density and mean velocity of each species:

$$n_{\alpha}(\mathbf{r}, t) = \int_{\mathbb{R}^3} f_{\alpha}(\mathbf{r}, \mathbf{v}, t) d\mathbf{v}, \quad \mathbf{u}_{\alpha}(\mathbf{r}, t) = \frac{1}{n_{\alpha}(\mathbf{r}, t)} \int_{\mathbb{R}^3} \mathbf{v} f_{\alpha}(\mathbf{r}, \mathbf{v}, t) d\mathbf{v}.$$

Boltzmann's equation for each species:

$$\frac{\partial f_{\alpha}}{\partial t} + \mathbf{v} \cdot \nabla f_{\alpha} + \frac{q_{\alpha}}{m_{\alpha}} (\mathbf{E} + \mathbf{v} \times \mathbf{B}) \cdot \nabla_{\mathbf{v}} f_{\alpha} = \sum_{\beta} Q_{\alpha} (f_{\alpha}, f_{\beta})(\mathbf{r}, \mathbf{v}, t)$$



# Boltzmann's equation

**Eulerian formulation:**  $(t, \mathbf{x}, \mathbf{v})$  independent variables

$$\frac{\partial f_\alpha}{\partial t} + \mathbf{v} \cdot \nabla f_\alpha + \frac{\mathbf{F}_\alpha}{m_\alpha} \cdot \nabla_{\mathbf{v}} f_\alpha = \left. \frac{\partial f_\alpha}{\partial t} \right|_{\text{coll}}$$

**Lagrangian formulation:** follow trajectory  $(\mathbf{x}(t), \mathbf{v}(t))$  in phase space

$$\frac{d\mathbf{x}}{dt} = \mathbf{v}(t), \quad \frac{d\mathbf{v}}{dt} = \frac{1}{m_\alpha} \mathbf{F}_\alpha(t, \mathbf{x}(t), \mathbf{v}(t))$$

- Substituting into Boltzmann's equation:  $\frac{Df_\alpha}{Dt} = \left. \frac{\partial f_\alpha}{\partial t} \right|_{\text{coll}}$
- Time rate of change of  $f_\alpha(t, \mathbf{x}(t), \mathbf{v}(t))$  along phase-space trajectory only determined by collision operator
- Without collisions,  $f_\alpha$  constant along phase-space trajectory: fluid motion in phase-space is **incompressible**

**Semi-Lagrangian method:**

- $f_\alpha(t, \mathbf{x}, \mathbf{v})$  lies on Eulerian mesh
- Evolution within time step uses Lagrangian formulation (*method of characteristics*)



# Modeling challenges

## WEAKLY COLLISIONAL PLASMA:

- Electrons can be far from equilibrium and involved in strongly non-linear processes (e.g. ionization near threshold)
- Multiple species: electrons, multiple ions, neutrals;
- Multiple time and spatial scales;
- Complex geometries, different boundary conditions (perfect/real conductors, dielectrics, absorbing), often time varying and coupled to domain (plasma feedbacks into circuit);
- Complex collisional processes: elastic, inelastic (excitation, ionization, recombination, attachment, dissociation etc.);
- External magnetic fields: electrons may be strongly magnetized, possibly ions too;
- Other important processes: radiation transport, gas-phase chemical reactions, plasma-surface interaction, aggregates (dusty plasmas).

## CHALLENGES FOR LOW-ORDER EULERIAN CODES:

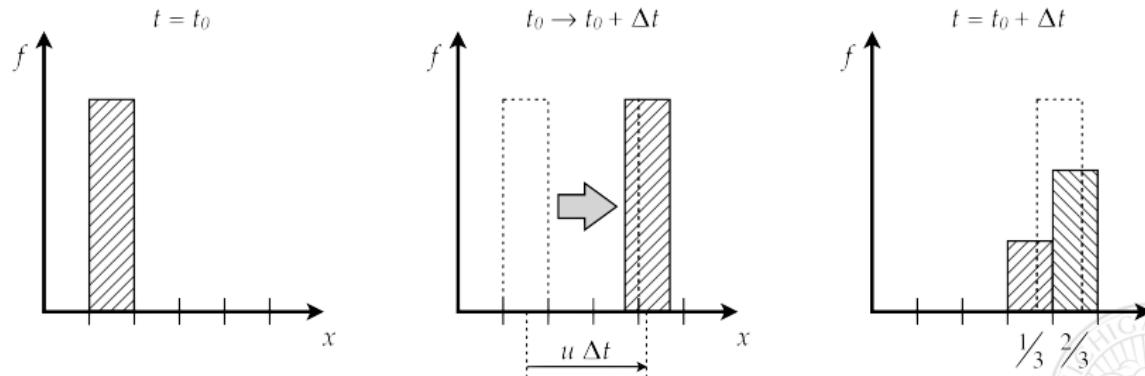
- For electrons, need high resolution over large velocity mesh
- Impressive memory requirement in multiple dimensions
- Explicit time-stepping imposes non-physical time-step restriction (CFL limit)
- Method of lines (MOL): multistep and multi-stage methods require additional storage



# Convected Scheme

The Convected Scheme [<sup>a</sup>] is a **forward semi-Lagrangian** method for Boltzmann's equation.  
Employs operator splitting:

1. **Collision operator** is local in configuration space, solves  $\frac{\partial f_\alpha}{\partial t} = \left. \frac{\partial f_\alpha}{\partial t} \right|_{\text{coll}}$
2. **Ballistic operator** advects  $f_\alpha(t, \mathbf{x}, \mathbf{v})$  along characteristic trajectories in phase space according to  $\frac{Df_\alpha}{Dt} = 0$ , integrated over a *moving cell* (MC).



$f_\alpha(t, \mathbf{x}, \mathbf{v})$  assumed uniform over MC, allowing for 'area remapping rule'

<sup>a</sup> W.N.G. HITCHON, D. KOCH, AND J. ADAMS. **An efficient scheme for convection-dominated transport.** *Journal of Computational Physics*, 83(1): 79-95, 1989.

# Convected Scheme

PROs:

- Preserves positivity (good as  $f_\alpha > 0$ )
- No CFL restriction on  $\Delta t$
- Very simple implementation
- Can enforce total energy conservation for stationary electric field

CONs:

- Difficult to handle boundary conditions
- **Numerical diffusion:** local remapping error  $O(\Delta x^2)$

## Reduced Numerical Diffusion

Numerical diffusion mitigated by reducing remapping frequency  $\Rightarrow$  “long-lived moving cells” [<sup>a</sup>]. Recently [<sup>b</sup>], we devised a high-order version of the Convected Scheme, for **neutral gas** kinetics:

**Model equation:** uniform velocity advection:  $n_t + u_0 n_x = 0$

**Basic idea:** compensating remapping error by applying small corrections to final position of moving cells prior to remapping  $\Rightarrow$  **antidiffusive velocity field**

**Tool:** modified equation analysis, perturbation analysis

<sup>a</sup> A.J. CHRISTLIEB, W.N.G. HITCHON AND E.R. KEITER. **A computational investigation of the effects of varying discharge geometry for an inductively coupled plasma.** *IEEE T. Plasma Sci.*, 28(6): 2214-2231, 2000.

<sup>b</sup> Y. GÜÇLÜ AND W.N.G. HITCHON. **A high order cell-centered semi-Lagrangian scheme for multi-dimensional kinetic simulations of neutral gas flows.** *Journal of Computational Physics*, 231(8): 3289-3316, Apr 2012.



# High-order semi-Lagrangian solution of the Vlasov-Poisson system

## PROBLEM:

Difficult to construct high-order semi-Lagrangian ballistic operator when mean force is present  
(no straight trajectories)

## SOLUTION:

- Further split ballistic operator into separate constant advection operators along  $x$  and  $v$  [<sup>a</sup>]
- Apply favorite high-order semi-Lagrangian solver to each operator
- Combine operators to high-order in time using Runge-Kutta-Nyström methods [<sup>b,c</sup>]  
**(symplectic**  $\Rightarrow$  energy stable)

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<sup>a</sup>C.Z. CHENG AND G. KNORR. **The integration of the Vlasov equation in configuration space.** *J. Comput. Phys.*, 22: 330-351, 1976.

<sup>b</sup>J.A. ROSSMANITH AND D.C. SEAL. **A positivity-preserving high-order semi-Lagrangian discontinuous Galerkin scheme for the Vlasov-Poisson equations.** *J. Comput. Phys.*, 227: 9527-9553, 2011.

<sup>c</sup>N. CROUSEILLES, E. FAOU AND M. MEHRENBERGER. **High order Runge-Kutta-Nyström splitting methods for the Vlasov-Poisson equation.** *INRIA-00633934*, 2011.

# Arbitrarily High-Order Convected Scheme (1)

## 1D CONSTANT ADVECTION EQUATION

$$\left( \frac{\partial}{\partial t} + u \frac{\partial}{\partial x} \right) n(x, t) = 0$$

- Exact solution (method of characteristics):  $n(x, t + \Delta t) \equiv n(x - u\Delta t, t)$
- Courant parameter:  $\alpha := u \Delta t / \Delta x$

## CONVECTED SCHEME UPDATE

- Discretize time (arbitrary  $\Delta t$ ) and space (uniform  $\Delta x$ ):  $n_i^k \approx n(x_i, t_k)$
- Because of uniform  $\Delta x$ , solution can be shifted exactly by integer number of cells
- Without loss of generality, assume  $0 \leq \alpha \leq 1$  (this is **not** a CFL limit)
- Under these assumptions, CS update is

$$n_i^{k+1} = U_{i-1}^k n_{i-1}^k + (1 - U_i^k) n_i^k$$

- As long as  $0 \leq U_i^k \leq 1$ , CS is **mass** and **positivity preserving**
- With no high-order corrections,  $U(x, t) \equiv \alpha \Rightarrow$  1st-order Upwind scheme
- With high-order corrections,  $U(x, t) = [u + \tilde{u}(x, t)] \Delta t / \Delta x = \alpha + \tilde{\alpha}(x, t)$
- $\tilde{\alpha}(x, t)$  is **anti-diffusive Courant parameter**



## Arbitrarily High-Order Convected Scheme (2)

## LOCAL TRUNCATION ERROR (LTE)

- Exact solution, Taylor expand in space (smooth initial conditions):

$$n(x, t + \Delta t) = n(x, t) + \left( \sum_{p=1}^{N-1} (-\alpha)^p \frac{(\Delta x)^p}{p!} \frac{\partial^p}{\partial x^p} \right) n(x, t) + O(\Delta x^N),$$

- CS solution, Taylor expand in space about  $(x, t) = (x_i, t^k)$ :

$$n_{\text{CS}}(x, t + \Delta t) = n(x, t) + \left( \sum_{p=1}^{N-1} (-1)^p \frac{(\Delta x)^p}{p!} \frac{\partial^p}{\partial x^p} \right) U(x, t) n(x, t) + O((\Delta x)^N)$$

- We want the local truncation error  $\mathcal{E}(x, t, \Delta t) := n(x, t + \Delta t) - n_{\text{CS}}(x, t + \Delta t) = O(\Delta x^N)$ , hence we find  $\tilde{\alpha}(x, t)$  by imposing the **order condition**

$$\sum_{p=1}^{N-1} (-\alpha)^p \frac{(\Delta x)^p}{p!} \frac{\partial^p n}{\partial x^p} - \sum_{p=1}^{N-1} (-1)^p \frac{(\Delta x)^p}{p!} \frac{\partial^p (Un)}{\partial x^p} = O(\Delta x^N)$$

# Arbitrarily High-Order Convected Scheme (3)

## HIGH-ORDER CORRECTIONS [<sup>a</sup>]

- Make the polynomial ansatz  $Un(x, t) = \sum_{q=0}^{N-2} (-1)^q \beta_q(\alpha) (\Delta x)^q \frac{\partial^q n(x, t)}{\partial x^q}$ ,

and solve for the unknown polynomials  $\beta_q(\alpha)$ ;

- Substitute in order condition to find (after algebraic manipulations)

$$[\tilde{\alpha}n]_i^k = \sum_{q=1}^{N-2} (-1)^q \frac{B_{q+1}(\alpha) - B_{q+1}(0)}{(q+1)!} (\Delta x)^q \left. \frac{\partial^q n(x, t)}{\partial x^q} \right|_i^k,$$

where  $B_q(\cdot)$  are Bernoulli polynomials;

- Approximate products  $(\Delta x)^q \left. \frac{\partial^q n(x, t)}{\partial x^q} \right|_i^k$  with error no larger than  $O(\Delta x^{N-1})$ , e.g.:
  1. linear polynomial interpolation,
  2. weighted essentially non-oscillatory (WENO) interpolation,
  3. fast Fourier transform (FFT).

<sup>a</sup>Y. GÜÇLÜ, A.J. CHRISTLIEB AND W.N.G. HITCHON. **Arbitrarily high order Convected Scheme solution of the Vlasov-Poisson system.** *In preparation.*

# Arbitrarily High-Order Convected Scheme (4)

## NUMERICAL IMPLEMENTATION [<sup>a</sup>]

- 6th-order finite difference scheme

$$\Delta x \left. \frac{\partial n}{\partial x} \right|_i^k \approx \frac{n_{i-2}^k - 8n_{i-1}^k + 8n_{i+1}^k - n_{i+2}^k}{12} + O(\Delta x^5),$$

$$(\Delta x)^2 \left. \frac{\partial^2 n}{\partial x^2} \right|_i^k \approx \frac{-n_{i-2}^k + 16n_{i-1}^k - 30n_i^k + 16n_{i+1}^k - n_{i+2}^k}{12} + O(\Delta x^6),$$

$$(\Delta x)^3 \left. \frac{\partial^3 n}{\partial x^3} \right|_i^k \approx \frac{-n_{i-2}^k + 2n_{i-1}^k - 2n_{i+1}^k - n_{i+2}^k}{2} + O(\Delta x^5),$$

$$(\Delta x)^4 \left. \frac{\partial^4 n}{\partial x^4} \right|_i^k \approx n_{i-2}^k - 4n_{i-1}^k + 6n_i^k - 4n_{i+1}^k + n_{i+2}^k + O(\Delta x^6),$$

- 22nd-order pseudo-spectral scheme

$$Un(x) = \mathcal{F}^{-1} \left[ \sum_{q=0}^{N-2} (-j)^q \beta_q(\alpha) (\xi \Delta x)^q \cdot \mathcal{F}[n](\xi) \right] (x),$$



<sup>a</sup>Y. GÜÇLÜ, A.J. CHRISTLIEB AND W.N.G. HITCHON. **Arbitrarily high order Convected Scheme solution of the Vlasov-Poisson system.** *In preparation.*

# Spectrally Accurate CS: Filtering

## PROBLEM:

- Stable scheme for nonlinear Vlasov-Poisson must dissipate sub-cell features;
- Pseudo-spectral method has no dissipation mechanism, and preserves discrete L2-norm  $\Rightarrow$  aliasing, Gibbs.

## SOLUTION:

- Apply Fourier windowing in  $k$ -space [<sup>a</sup>]  $\Rightarrow$  equivalent to Gaussian-regularized Sinc convolution in  $x$  (Gaussian has width  $W\Delta x$ );
- Adaptive filtering investigated in [<sup>a</sup>], but in our experience, need very accurate indicator to avoid excessive dissipation;
- We use constant  $W = 4$  in all numerical examples  $\Rightarrow$  low filter strength preserves spectral accuracy.

<sup>a</sup>Y. SUN, Y. ZHOU, S.-G. LI AND G. WEI. A windowed Fourier pseudospectral method for hyperbolic conservation laws. *J. of Comput. Phys.*, 214(2): 466-490, 2006.



# Spectrally Accurate CS: 1D Refinement Study

1D constant advection equation (normalized):

$$\frac{\partial n}{\partial t} + \frac{\partial n}{\partial x} = 0, \quad x \in [-0.5, 0.5], \quad t \in [0, 1],$$

Smooth initial conditions:

$$n_0(0, x) = 0.5 e^{-(\frac{x+0.2}{0.03})^2} + e^{-(\frac{x}{0.06})^2} + 0.5 e^{-(\frac{x-0.2}{0.03})^2}.$$

Refinement study:

<b>N<sub>x</sub></b>	<b>P4</b>	<b>Order</b>	<b>P6</b>	<b>Order</b>	<b>F22</b>	<b>Order</b>
32	$1.41 \times 10^{-01}$	—	$7.68 \times 10^{-02}$	—	$2.47 \times 10^{-02}$	—
64	$5.99 \times 10^{-02}$	1.24	$2.55 \times 10^{-02}$	1.59	$1.84 \times 10^{-04}$	7.07
128	$2.28 \times 10^{-02}$	1.39	$4.45 \times 10^{-03}$	2.52	$7.55 \times 10^{-11}$	21.22
256	$5.44 \times 10^{-03}$	2.07	$2.14 \times 10^{-04}$	4.37	$1.02 \times 10^{-13}$	9.53
512	$7.94 \times 10^{-04}$	2.78	$7.03 \times 10^{-06}$	4.93	$1.20 \times 10^{-13}$	m.p.
1024	$1.02 \times 10^{-04}$	2.96	$2.21 \times 10^{-07}$	4.99	$4.44 \times 10^{-13}$	m.p.
2048	$1.28 \times 10^{-05}$	2.99	$6.93 \times 10^{-09}$	5.00	$4.20 \times 10^{-13}$	m.p.

# Operator Splitting: ODEs

Separable ODE

$$\dot{u}(t) = A(u) + B(u)$$

Assume we can construct the exact solution to the subproblems

$$\dot{u}(t) = A(u) \quad \dot{u}(t) = B(u)$$

in the form of the Lie operators

$$u_A(t + \Delta t) = e^{A\Delta t} u(t) \quad u_B(t + \Delta t) = e^{B\Delta t} u(t)$$

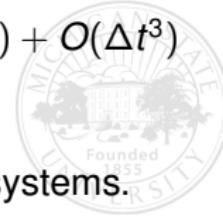
If the operators do not commute, composition introduces splitting error. E.g.

Lie-Trotter  $u(t + \Delta t) = e^{A\Delta t} e^{B\Delta t} u(t) + O(\Delta t^2)$

Leapfrog/Störmer-Verlet/Strang  $u(t + \Delta t) = e^{A\Delta t/2} e^{B\Delta t} e^{A\Delta t/2} u(t) + O(\Delta t^3)$

**NOTE 1:** Composition of low-order schemes to get higher order.

**NOTE 2:** Most popular as symplectic integrators for Hamiltonian systems.



# Operator Splitting: Linear Transport PDEs

Separable PDE: let  $u = u(t, x)$ , and  $A(\cdot)$  and  $B(\cdot)$  be integro-differential operators:

$$\frac{\partial u}{\partial t} = A(u) + B(u)$$

After discretization in  $x$  we have the approximate  $e^{\tilde{A}\Delta t}$  and  $e^{\tilde{B}\Delta t}$  where some error  $O(\Delta x^p)$  was introduced.

Dimensional splitting for linear equations

$$\frac{\partial u}{\partial t} = a(x_2) \frac{\partial u}{\partial x_1} + b(x_1) \frac{\partial u}{\partial x_2} \quad \text{with } u = u(t, x_1, x_2)$$

leads to 2 families of **constant advection equations**.

Hence, we can use our semi-Lagrangian solver for:

- 2D rotating advection
- linear Vlasov equation



# Operator Splitting: Vlasov-Poisson (1)

## Complication:

Time-varying electric field may reduce accuracy of splitting method!?

## CHENG & KNORR'S ALGORITHM [a]

Based on Strang splitting, uses previous value of  $\mathbf{E}$ , 2nd-order:

1.  $\Delta t/2$  step on  $(\partial_t + \mathbf{v} \cdot \nabla_{\mathbf{x}}) f_e = 0$ ;
2. Compute  $n_e = \int_{\mathbb{R}^3} f_e d\mathbf{v}$ , solve  $\nabla_{\mathbf{x}}^2 \phi = \frac{q_e}{\epsilon_0} (n_e - n_0)$ , and evaluate  $\mathbf{E} = -\nabla_{\mathbf{x}} \phi$ ;
3.  $\Delta t$  step on  $(\partial_t + \frac{q_e}{m_e} \mathbf{E} \cdot \nabla_{\mathbf{v}}) f_e = 0$ ;
4.  $\Delta t/2$  step on  $(\partial_t + \mathbf{v} \cdot \nabla_{\mathbf{x}}) f_e = 0$ .

**NOTE:** In [a] semi-Lagrangian advection (1D-1V, cubic splines in  $x$ , trigonometric interpolation in  $v$ )



<sup>a</sup> C. CHENG, AND G. KNORR. The integration of the Vlasov equation in configuration space. *J. Comput. Phys.*, 22(3): 330–351, 1976.

# Operator Splitting: Vlasov-Poisson (2)

**Problem:** how to go higher order in time?

- High-order extrapolation of  $\mathbf{E}$  and time averaging [<sup>a</sup>]  
(Lax-Wendroff / Cauchy-Kovalesky procedure)
- Use high-order symplectic integrators [<sup>b,c,d</sup>] based on Hamiltonian splitting  $H = T + V$ , with  $s$  stages:

$$\begin{cases} f_k^*(\mathbf{x}, \mathbf{v}) = f_{k-1}\left(\mathbf{x} - (a_k \Delta t) \mathbf{v}, \mathbf{v}\right) \\ f_k(\mathbf{x}, \mathbf{v}) = f_k^*\left(\mathbf{x}, \mathbf{v} - (b_k \Delta t) \frac{q}{m} \mathbf{E}[f_k^*](\mathbf{x})\right) \end{cases} \quad (k = 1, 2, \dots, s),$$

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<sup>a</sup>J.A. ROSSMANITH AND D.C. SEAL. **A positivity-preserving high-order semi-Lagrangian discontinuous Galerkin scheme for the Vlasov-Poisson equations.** *J. Comput. Phys.*, 227: 9527-9553, 2011.

<sup>b</sup>T. WATANABE AND H. SUGAMA. **Vlasov and drift-kinetic simulation methods based on the symplectic integrator.** *Trans. Th. Stat. Phys.* 34: 287-309, 2005.

<sup>c</sup>E. POHN, M. SHOUCRU AND G. KAMELANDER. **Eulerian Vlasov codes.** *Comput. Phys. Comm.*, 166(2): 81-93, 2005.

<sup>d</sup>N. CROUSEILLES, E. FAOU AND M. MEHRENBERGER. **High order Runge-Kutta-Nyström splitting methods for the Vlasov-Poisson equation.** INRIA-00633934, 2011.

# Operator Splitting: Vlasov-Poisson (3)

Methods compared:

Label	Description	Order	Stages	Refs.
LF2	Leap-frog / Strang / Störmer-Verlet	2	1	[ <sup>a</sup> ]
Y4	Triple-jump composition of LF2	4	3	[ <sup>b</sup> ],[ <sup>c</sup> ],[ <sup>d</sup> ]
O6-4	4th-order RKN, optimized	4	6	[ <sup>e</sup> ]
O11-6	6th-order RKN, optimized	6	11	[ <sup>e</sup> ]
O14-6	6th-order RKN, optimized	6	14	[ <sup>e</sup> ]

RKN = Runge-Kutta-Nyström

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<sup>a</sup> G. STRANG. **On the Construction and Comparison of Difference Schemes.** *SIAM J. Numer. Anal.* 5(3): 506–517, 1968.

<sup>b</sup> H. YOSHIDA. **Construction of higher order symplectic integrators.** *Phys. Lett. A*, 150(5-7): 262-268, 1990.

<sup>c</sup> E. FOREST AND R.D.RUTH. **Fourth-order symplectic integration.** *Physica D: Nonlinear Phenomena*, 43(1): 105-117, 1990.

<sup>d</sup> J. CANDY AND W. ROZMUS. **A symplectic integration algorithm for separable Hamiltonian functions.** *J. Comput. Phys.* 92(1): 230-256, 1991.

<sup>e</sup> S. BLANES AND P. MOAN. **Practical symplectic partitioned Runge–Kutta and Runge–Kutta–Nyström methods.** *J. Comp. Appl. Math.* 142(2): 313-330, 2002.

# 2D Rotating Advection

**Goal:** characterize splitting error (closed orbits? phase error?)

- Model equation:  $\frac{\partial n}{\partial t} - (2\pi y) \frac{\partial n}{\partial x} + (2\pi x) \frac{\partial n}{\partial y} = 0$
- Square domain:  $(x, y) \in [-1, 1] \times [-1, 1]$
- Boundary conditions: periodic
- Initial conditions:  
superposition of two  
( $\cos$ )<sup>22</sup> bells with  
elliptical cross-section

$$n(0, x, y) = 0.5 B(r_1(x, y)) + 0.5 B(r_2(x, y)),$$

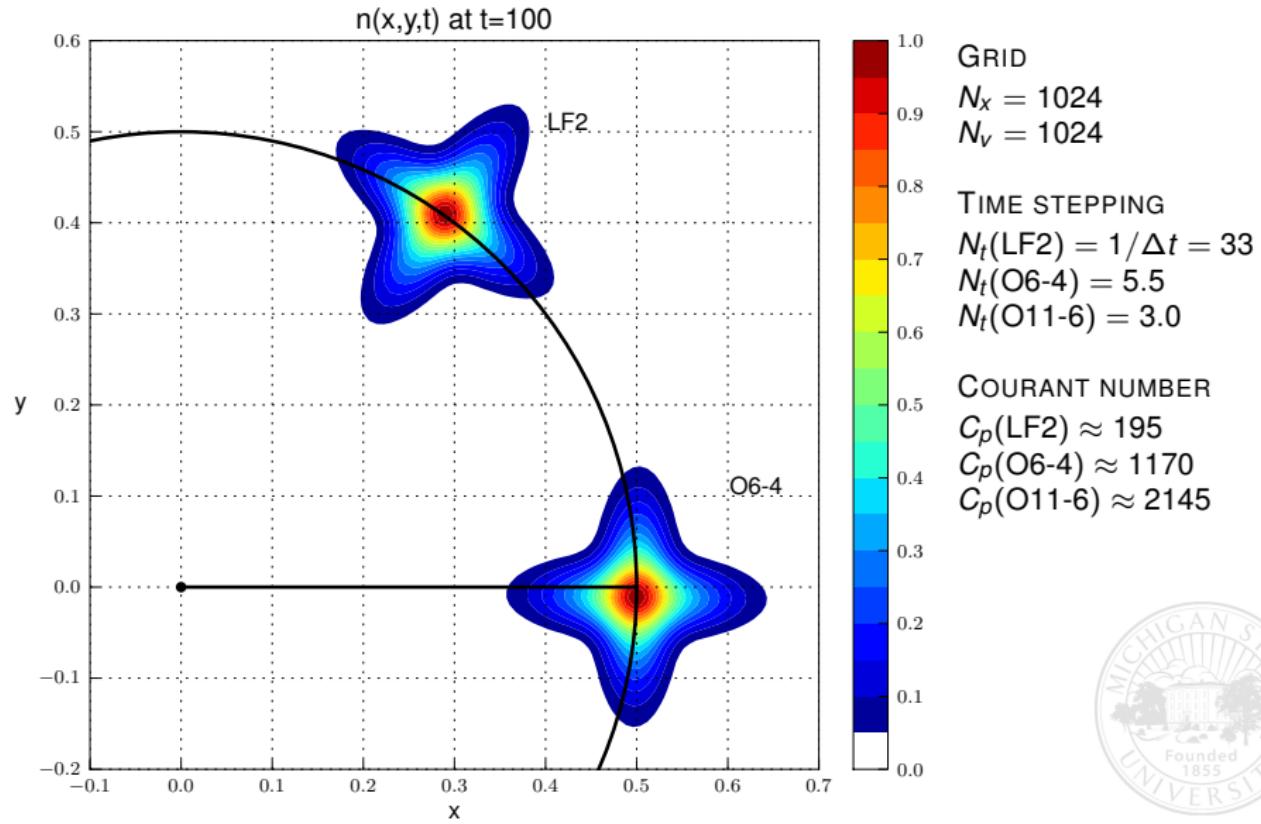
$$B(r) = \begin{cases} \cos\left(\frac{\pi r}{2a}\right)^{22} & \text{if } r \leq a, \\ 0 & \text{otherwise,} \end{cases}$$

$$r_1(x, y) = \sqrt{(x - x_c)^2 + 8(y - y_c)^2},$$

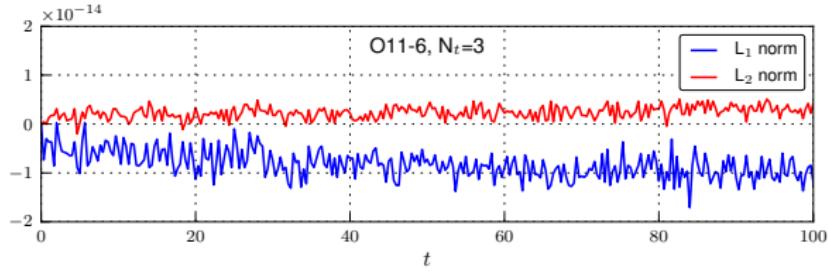
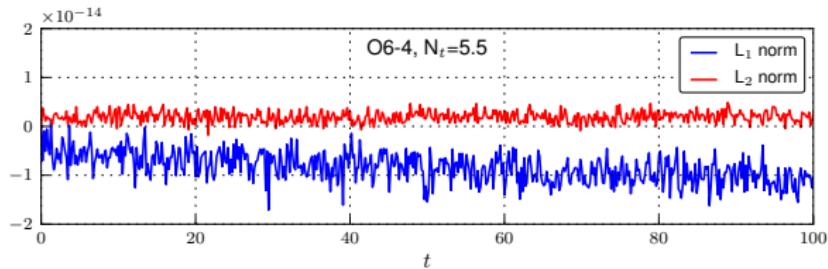
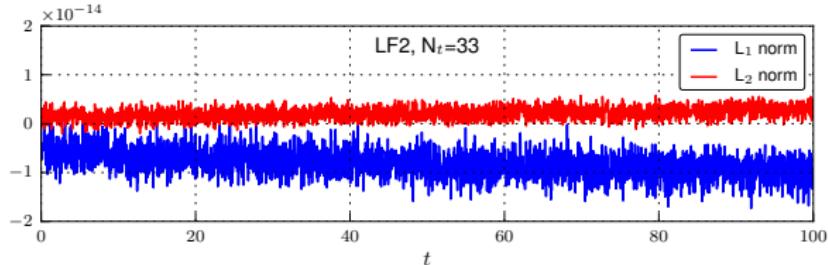
$$r_2(x, y) = \sqrt{8(x - x_c)^2 + (y - y_c)^2},$$



# 2D Rotating Advection: phase-error



# 2D Rotating Advection: invariants / area preservation



GRID

$$N_x = 1024$$

$$N_v = 1024$$

TIME STEPPING

$$N_t(\text{LF2}) = 1/\Delta t = 33$$

$$N_t(\text{O6-4}) = 5.5$$

$$N_t(\text{O11-6}) = 3.0$$

COURANT NUMBER

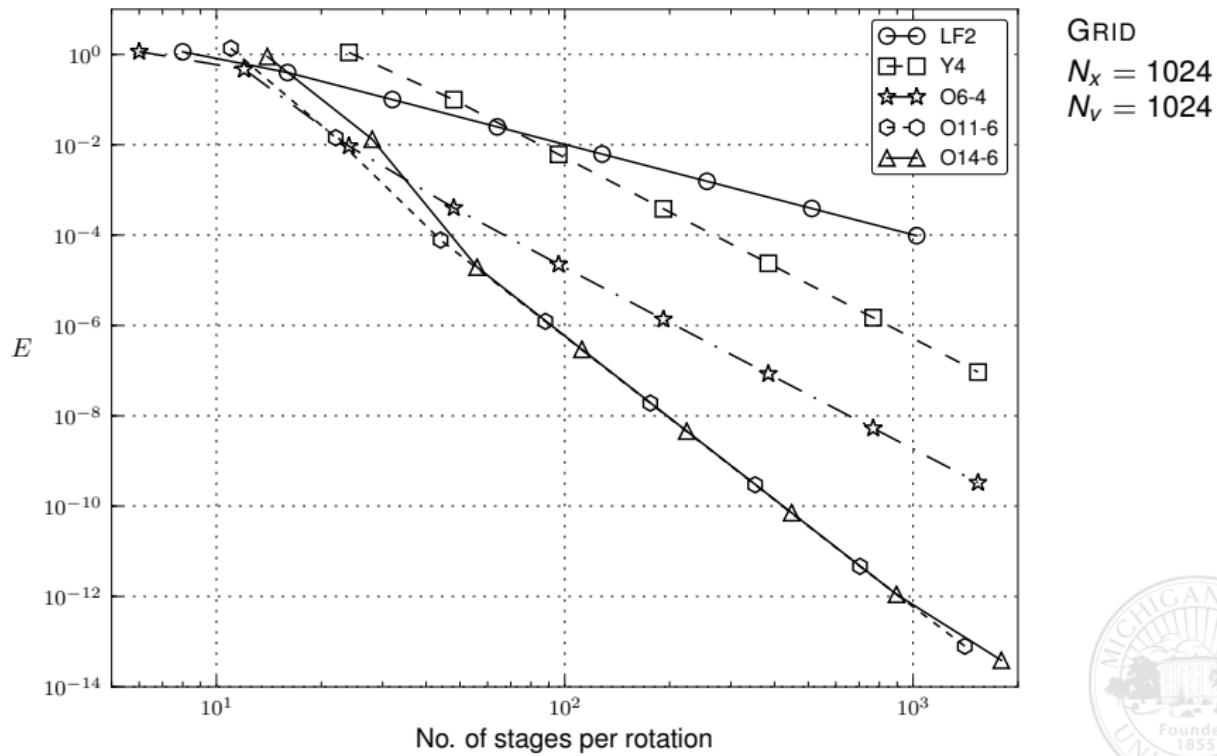
$$C_p(\text{LF2}) \approx 195$$

$$C_p(\text{O6-4}) \approx 1170$$

$$C_p(\text{O11-6}) \approx 2145$$



# 2D Rotating Advection: efficiency



# Linear Vlasov Equation with Stationary Field

**Goal:** simulate trapping of electrons in stationary field

**Challenges:** fast filamentation, long-time confinement

- Model equation (normalized)

$$\frac{\partial f}{\partial t} + v \frac{\partial f}{\partial x} - E(x) \frac{\partial f}{\partial v} = 0$$

- Electric field  $E(x)$  given, not self-consistent
- Square domain:  $(x, v) \in [-1, 1] \times [-1, 1]$
- Periodic boundary conditions in  $x$  and  $v$  (assume compact support)
- Special case of 2D advection equation in  $(x, v)$  coordinates
- Phase-space flow is incompressible:  $\partial v / \partial x - \partial E / \partial v = 0$



# Linear Vlasov: phase-space vorticity

- Prescribed stationary potential:

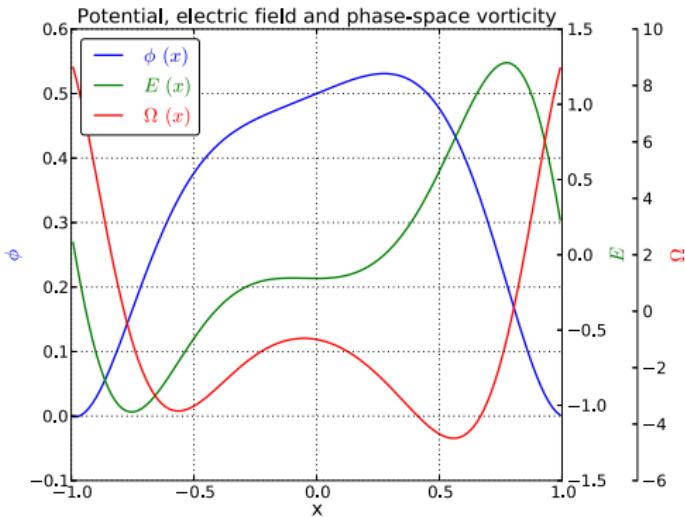
$$\phi(x) = \frac{1 + \cos(\pi x^2)}{4} + \frac{\sin(\pi x)}{20}$$

- 2D flow field:

$$\mathbf{u}(x) = \begin{bmatrix} u_1(x_1, x_2) \\ u_2(x_1, x_2) \end{bmatrix}$$

- 2D vorticity:

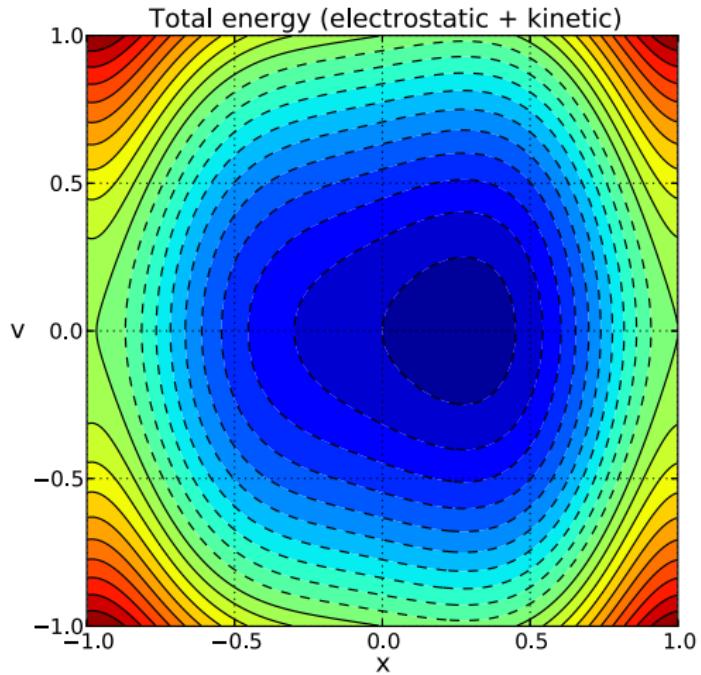
$$\nabla \times \mathbf{u} = \left[ \frac{\partial u_2}{\partial x_1} - \frac{\partial u_1}{\partial x_2} \right] \hat{\mathbf{n}}_3 = \Omega(x_1, x_2) \hat{\mathbf{n}}_3$$



For 1D-1V Vlasov,  $(x_1, x_2) = (x, v)$  and  $(u_1, u_2) = (v, -E(x))$ .  
Hence the **phase-space vorticity** depends on  $x$  only:

$$\Omega(x, v) = -\frac{\partial E}{\partial x} - \frac{\partial v}{\partial v} = \frac{\partial^2 \phi}{\partial x^2} - 1 = \Omega(x)$$

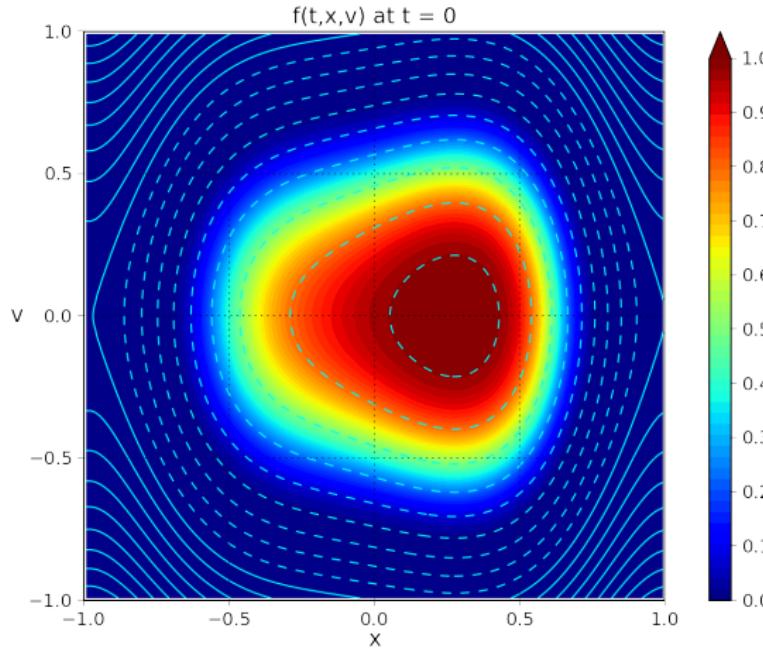
# Linear Vlasov: Hamiltonian



- One-particle Hamiltonian:  
$$H(x, v) = \frac{v^2}{2} - \phi(x)$$
- Hamiltonian is constant of motion:  $H(x(t), v(t)) \equiv c$
- $H(x, v) > 0$  (solid lines): open trajectories
- $H(x, v) \leq 0$  (dashed lines): closed trajectories **(electrostatic confinement)**

If Vlasov solver is **energy-stable**, confined particles will remain confined.

# Linear Vlasov: steady-state preservation



INITIAL CONDITIONS

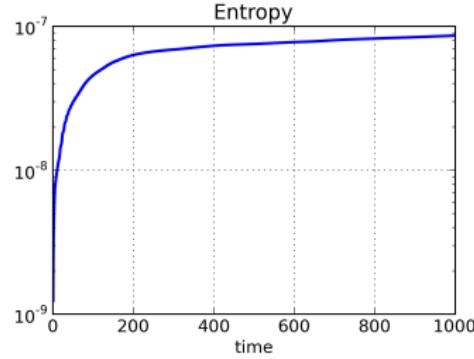
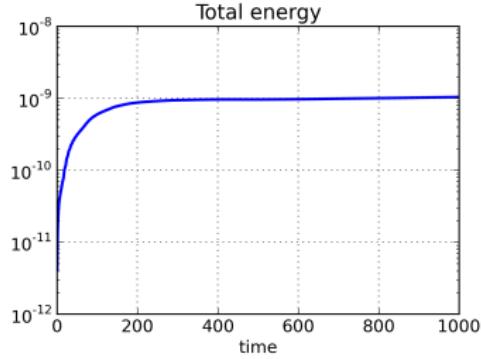
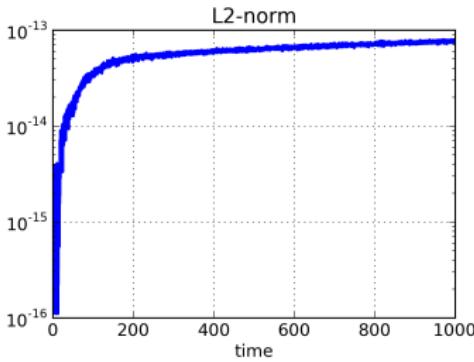
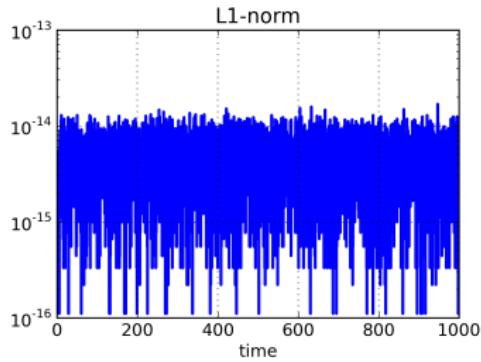
$$f(0, x, v) = g(H(x, v) - \min(H))$$

$$g(h) = \begin{cases} \cos(\pi h)^6 & \text{if } h < 0.5 \\ 0.0 & \text{otherwise} \end{cases}$$



# Linear Vlasov: steady-state preservation

Relative errors in conserved quantities



GRID

$N_x = 128$

$N_v = 128$

TIME STEPPING

$\Delta t = 0.25$   
4000 steps

COURANT NO.

$C_x \approx 16$

$C_v \approx 20$



# Linear Vlasov: filamentation (1)

## GRID

$N_x = 512$

$N_v = 512$

## TIME STEPPING

$\Delta t = 0.5$

100+100 steps

## COURANT NUMBER

$C_x \approx 130$

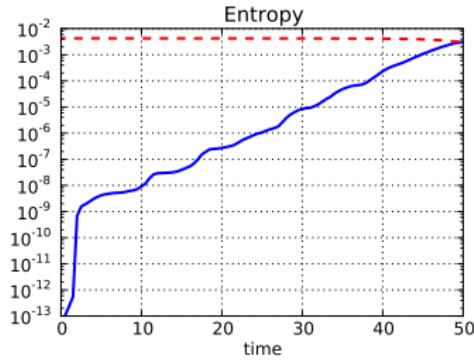
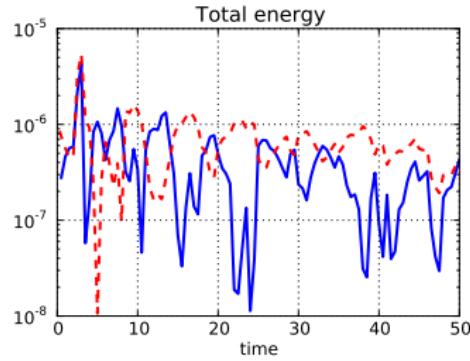
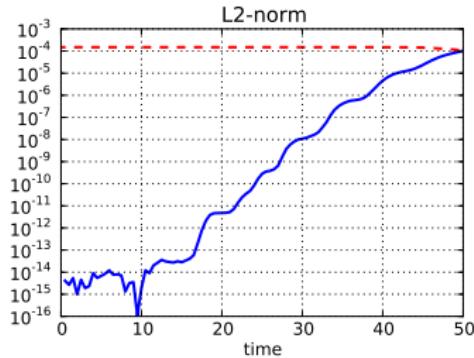
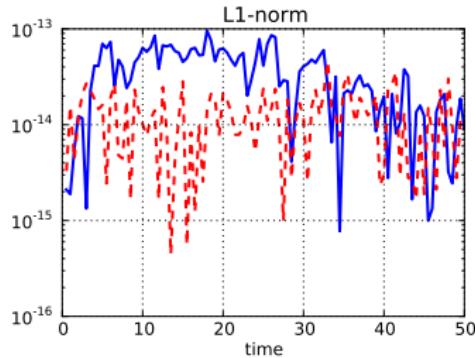
$C_v \approx 160$

(Loading Filamentation\_coarse\_animation.mp4)



# Linear Vlasov: filamentation (2)

Relative errors in conserved quantities



GRID  
 $N_x = 512$   
 $N_v = 512$

TIME STEPPING  
 $\Delta t = 0.5$   
100+100 steps

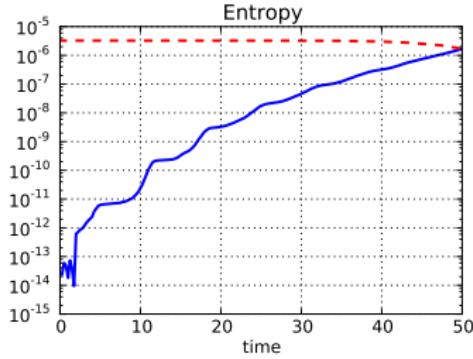
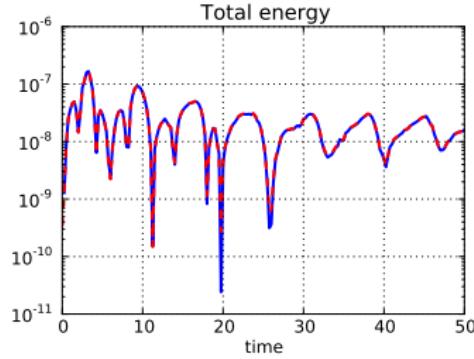
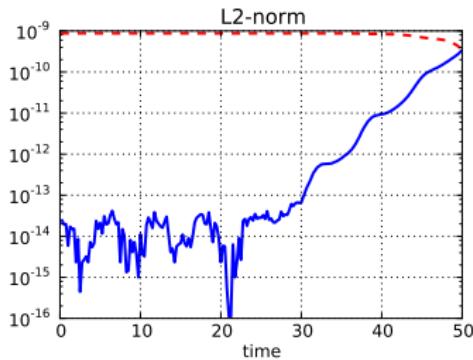
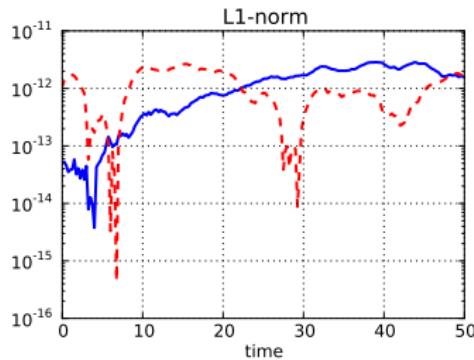
COURANT NO.  
 $C_x \approx 130$   
 $C_v \approx 160$

Solid blue line:  
**forward** evolution  
Dashed red line:  
**backward** evolut.

Founded  
1855  
UNIVERSITY

# Linear Vlasov: filamentation (3)

Relative errors in conserved quantities



GRID  
 $N_x = 1024$   
 $N_v = 1024$

TIME STEPPING  
 $\Delta t = 0.25$   
200+200 steps

COURANT NO.  
 $C_x \approx 130$   
 $C_v \approx 160$

Solid blue line:  
**forward** evolution  
Dashed red line:  
**backward** evolut.

Founded  
1855

# Vlasov-Poisson: bump-on-tail instability (1)

GRID

$N_x = 256$

$N_v = 512$

TIME STEPPING

$\Delta t = 0.5$

44 steps

O6-4

(Loading BumpOnTail\_animation.mp4)

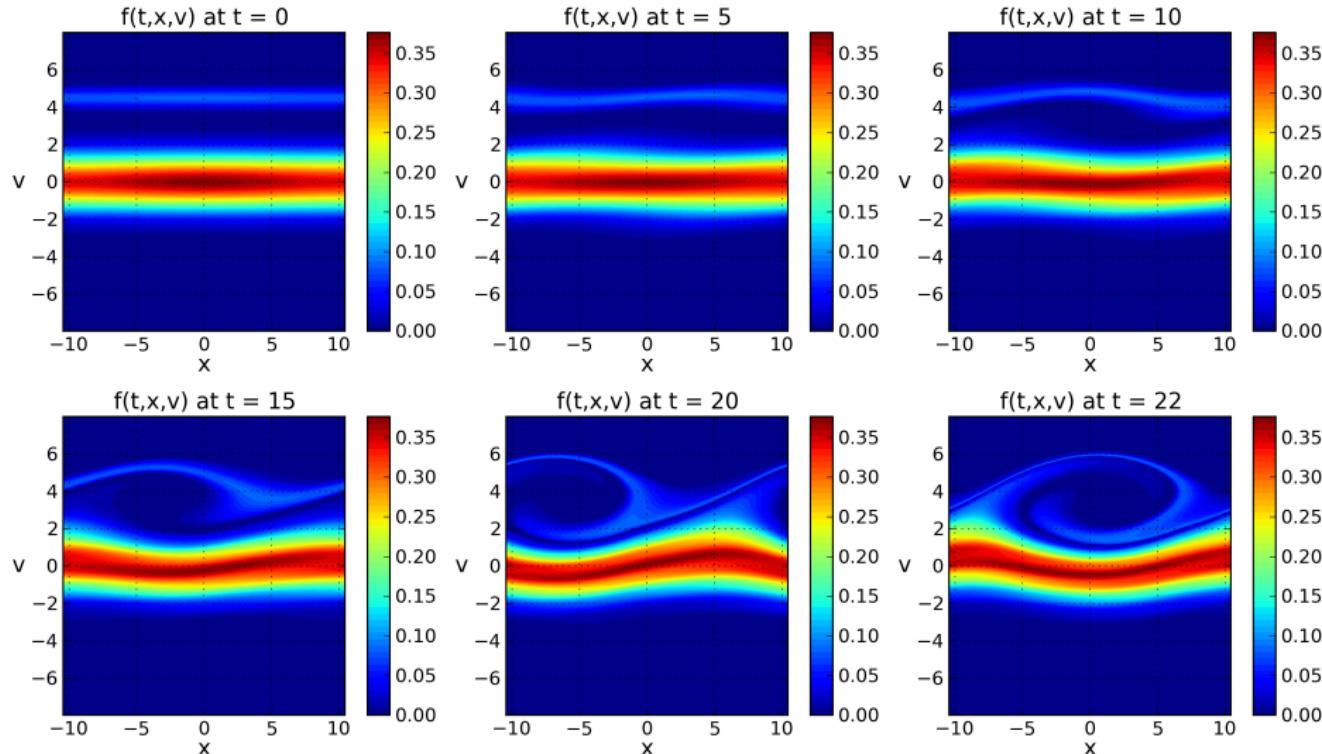
COURANT NUMBER

$C_x \approx 49$

$C_v \approx 9.4$



# Vlasov-Poisson: bump-on-tail instability (2)



$$N_x = 256$$

$$N_v = 512$$

$$\Delta t = 0.5$$

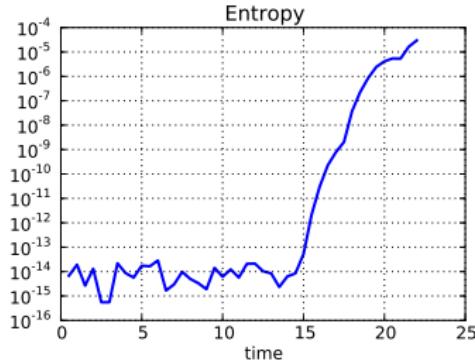
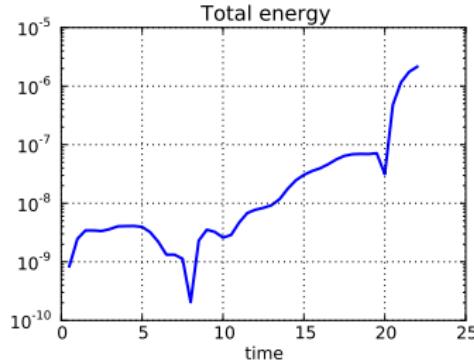
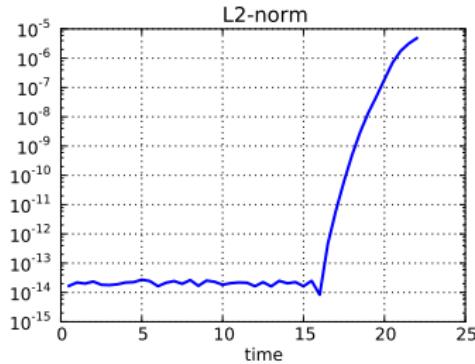
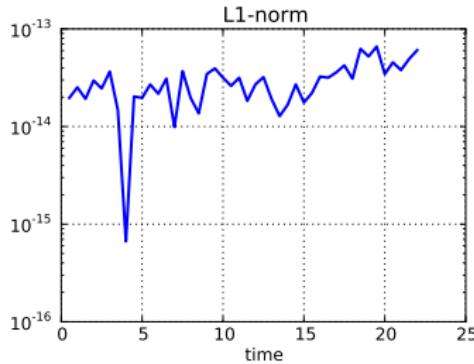
44 steps

$$[C_x \approx 49 \quad C_v \approx 9.4]$$

$$O6-4$$

# Vlasov-Poisson: bump-on-tail instability (3)

Relative errors in conserved quantities



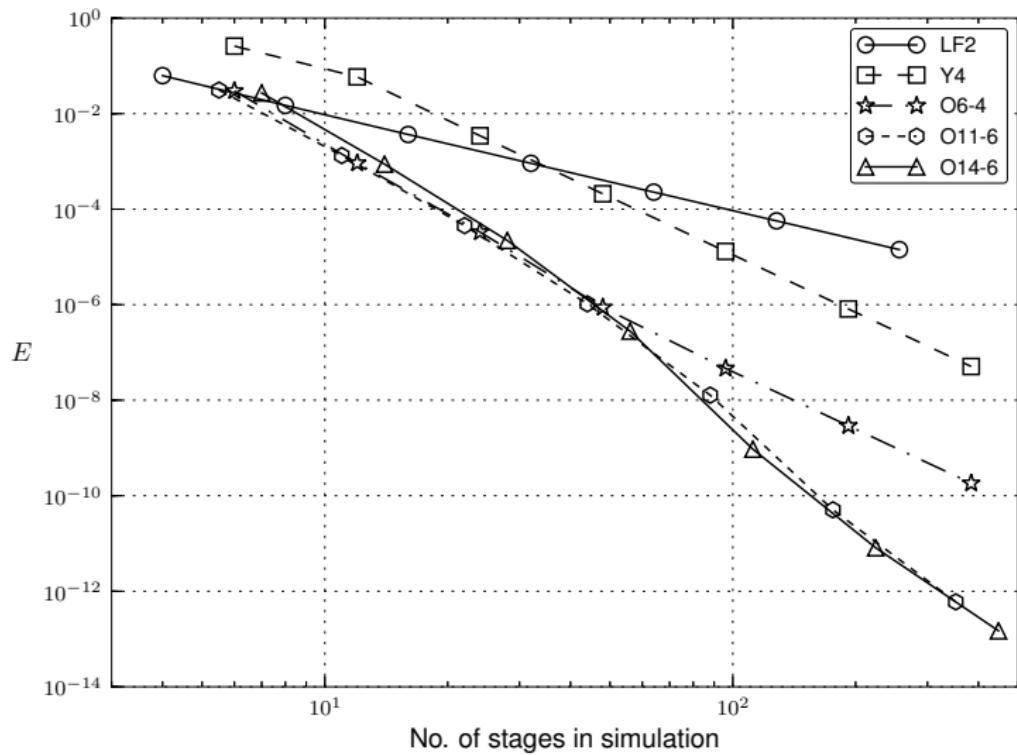
GRID  
 $N_x = 256$   
 $N_v = 512$

TIME STEPPING  
 $\Delta t = 0.5$   
44 steps  
O6-4

COURANT NO.  
 $C_x \approx 49$   
 $C_v \approx 9.4$



# Vlasov-Poisson: bump-on-tail instability (4)



FINAL TIME  
 $T = 16$

GRID  
 $N_x = 1024$   
 $N_v = 1024$

REFERENCE  
 $N_x = 1024$   
 $N_v = 1024$   
 $\Delta t = 0.05$   
O11-6



# Vlasov-Poisson: linear Landau damping (1)

GRID

$N_x = 16$

$N_v = 256$

TIME STEPPING

$\Delta t = 0.5$

120 steps

O6-4

(Loading LinearLandau\_animation.mp4)

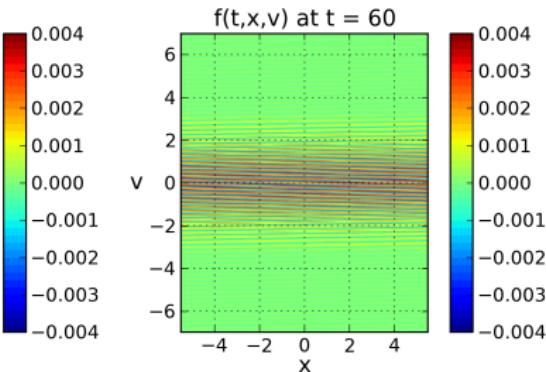
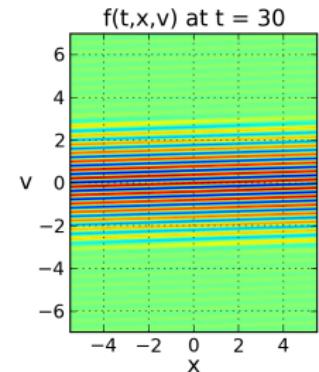
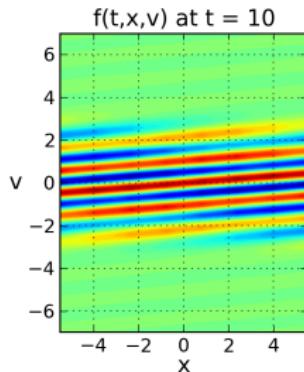
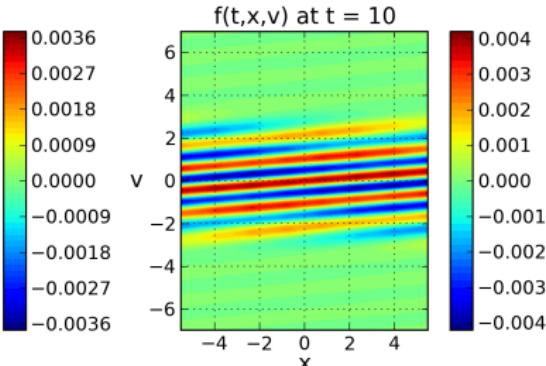
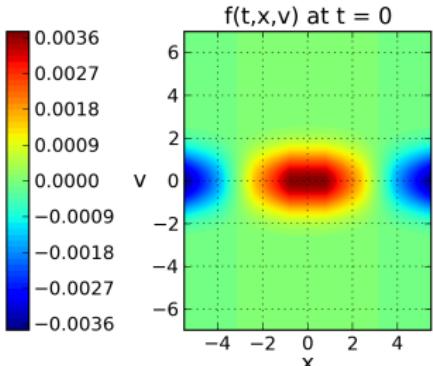
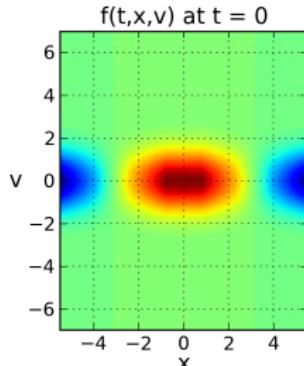
COURANT NUMBER

$C_x \approx 4.0$

$C_v \approx 0.2$



# Vlasov-Poisson: linear Landau damping (2)



$N_x = 8$

$N_v = 256$

$\Delta t = 1.0$

60 steps

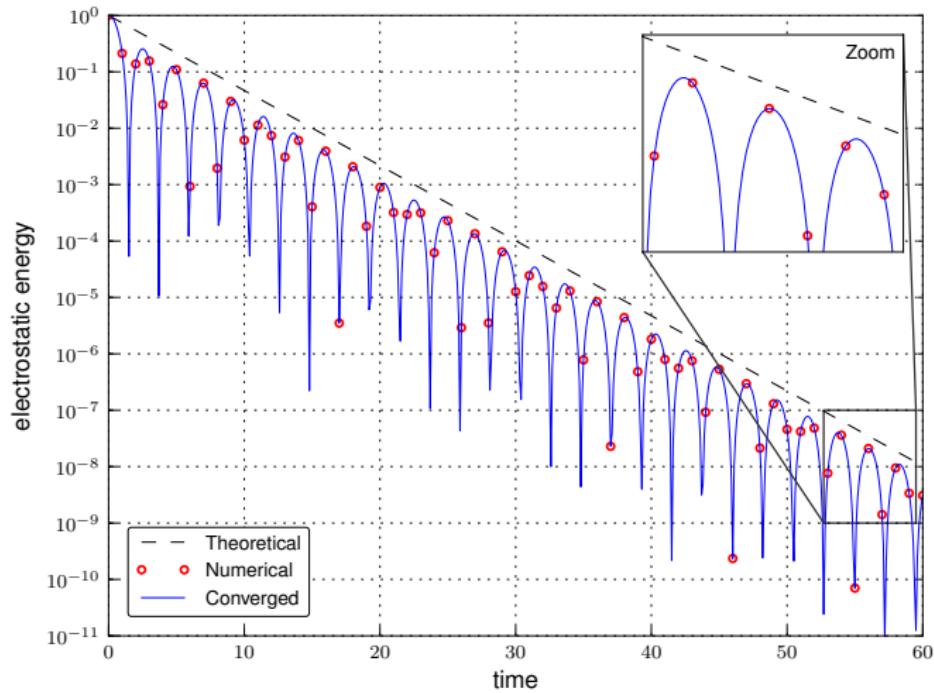
$[C_x \approx 4.0]$

$C_v \approx 0.4]$

O6-4



# Vlasov-Poisson: linear Landau damping (3)



GRID

$N_x = 8$

$N_v = 256$

TIME STEPPING

$\Delta t = 1.0$

60 steps

O6-4

COURANT NUMBER

$C_x \approx 4.0$

$C_v \approx 0.4$

REFERENCE

$N_x = 16$

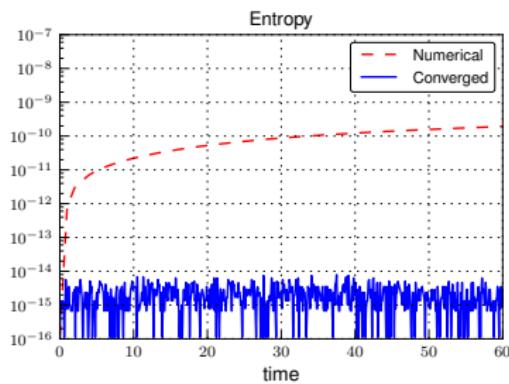
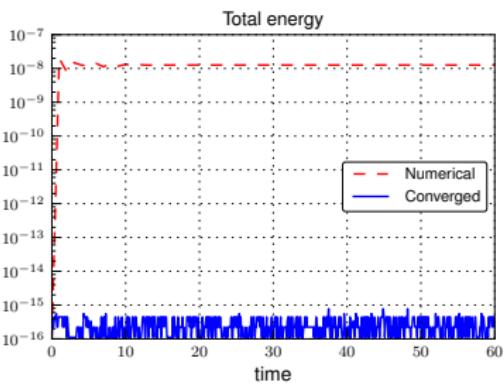
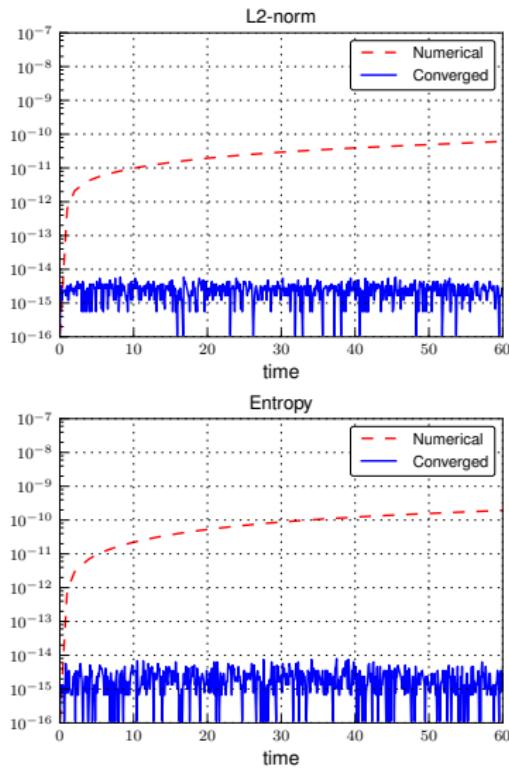
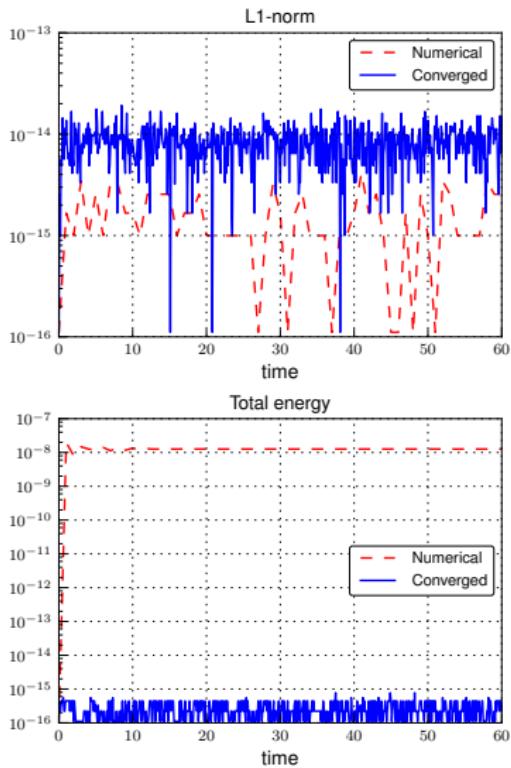
$N_v = 512$

$\Delta t = 0.1$

O11-6



# Vlasov-Poisson: linear Landau damping (4)



GRID

$N_x = 8$   
 $N_v = 256$

TIME STEPPING

$\Delta t = 1.0$   
60 steps  
O6-4

COURANT NO.

$C_x \approx 4.0$   
 $C_v \approx 0.4$

REFERENCE

$N_x = 16$   
 $N_v = 512$   
 $\Delta t = 0.1$   
O11-6



# Vlasov-Poisson: non-linear Landau damping (1)

GRID

$N_x = 256$

$N_v = 512$

TIME STEPPING

$\Delta t = 0.5$

120 steps

(Loading LinearLandau\_animation.mp4)

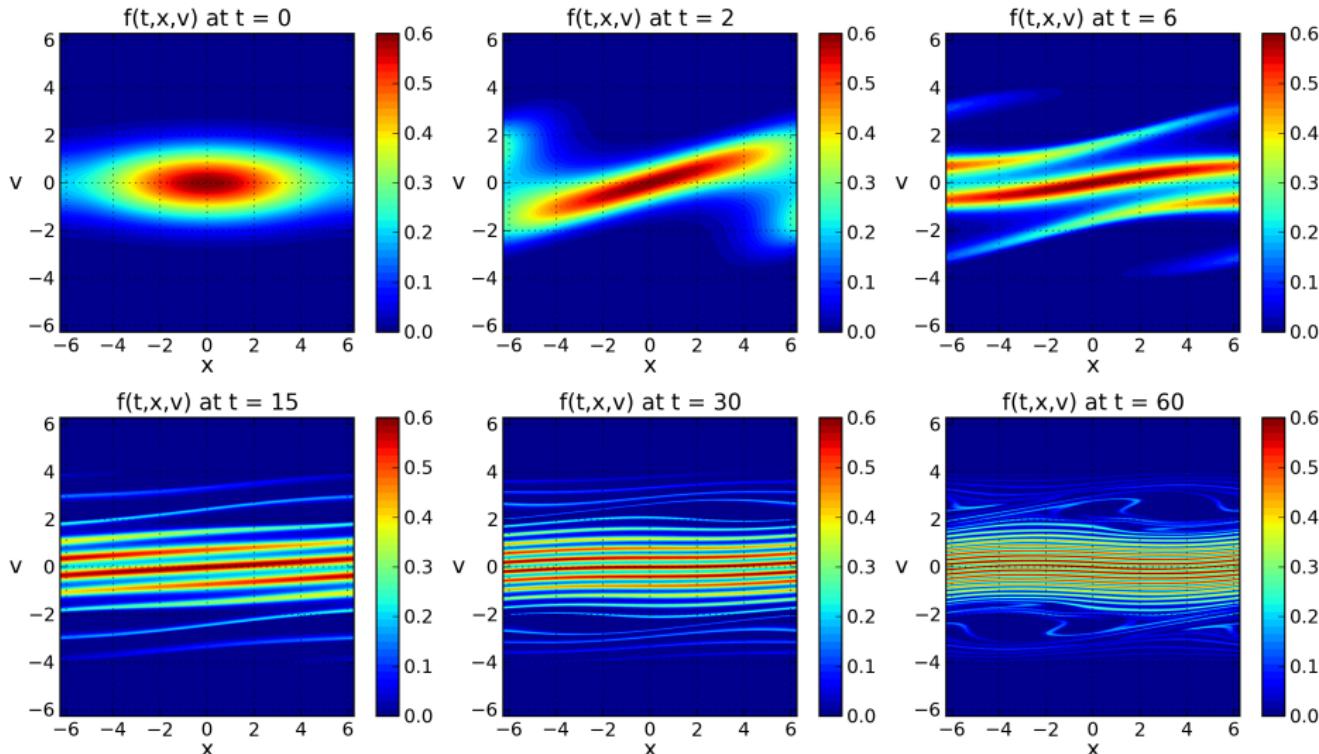
COURANT NUMBER

$C_x \approx 64$

$C_v \approx 20$



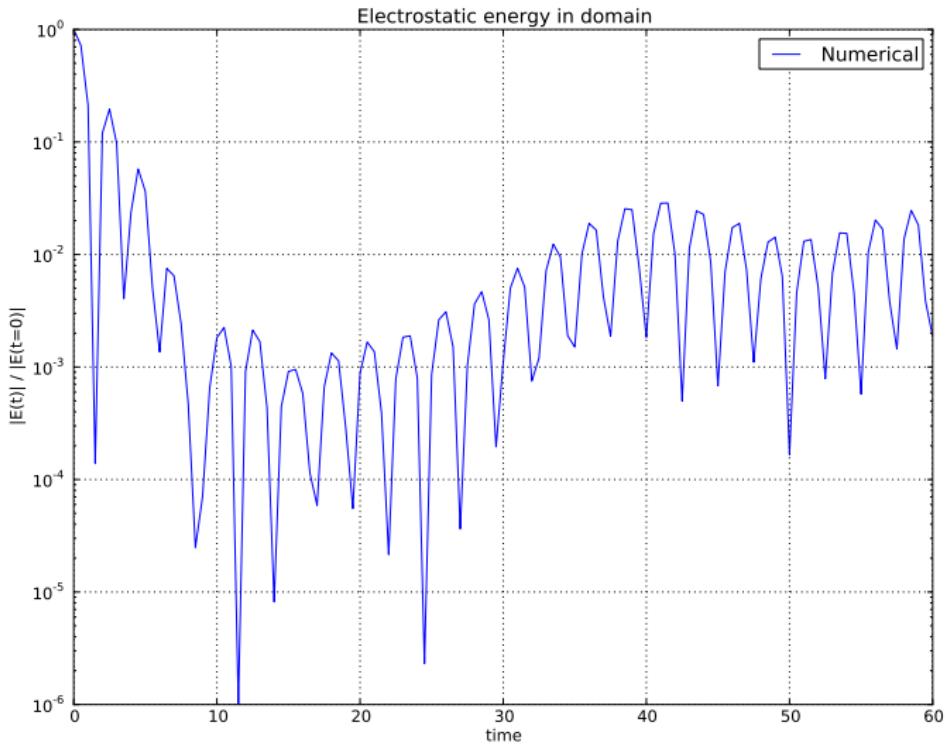
# Vlasov-Poisson: non-linear Landau damping (2)



$N_x = 256 \quad N_v = 512 \quad \Delta t = 0.5 \quad 120$  steps  $[C_x \approx 64 \quad C_v \approx 20]$



# Vlasov-Poisson: non-linear Landau damping (3)



GRID

$N_x = 256$

$N_v = 512$

TIME STEPPING

$\Delta t = 0.5$

120 steps

COURANT NUMBER

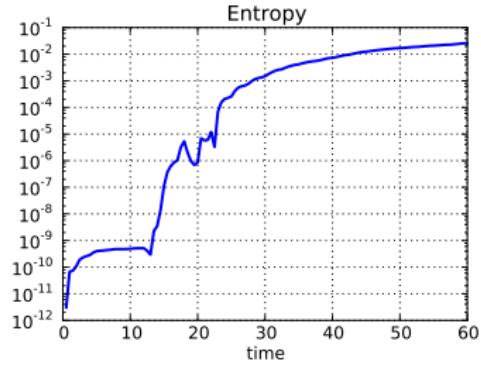
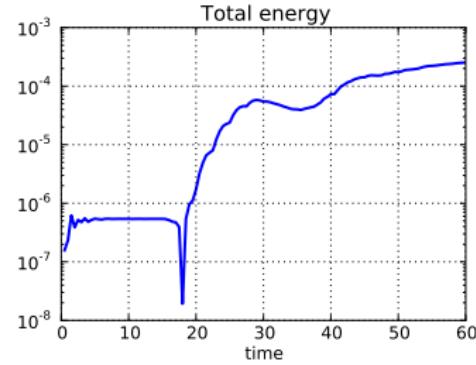
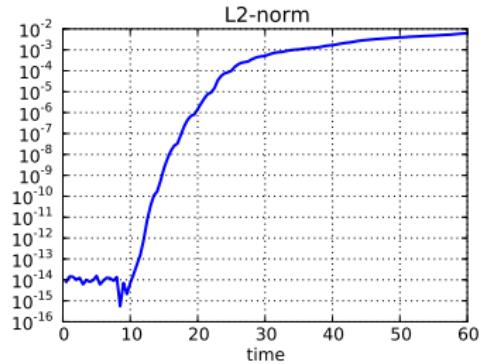
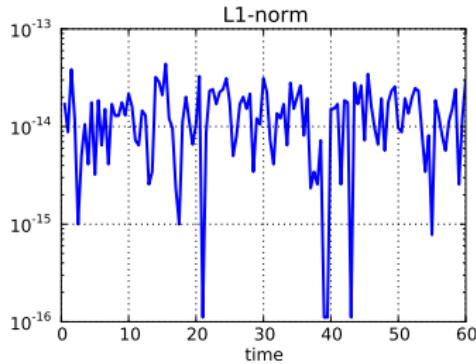
$C_x \approx 64$

$C_v \approx 20$



# Vlasov-Poisson: non-linear Landau damping (4)

Relative errors in conserved quantities



GRID  
 $N_x = 256$   
 $N_v = 512$

TIME STEPPING  
 $\Delta t = 0.5$   
120 steps

COURANT NO.  
 $C_x \approx 64$   
 $C_v \approx 20$



# Vlasov-Poisson: two-stream instability (1)

GRID

$N_x = 256$

$N_v = 512$

TIME STEPPING

$\Delta t = 0.5$

90 steps

(Loading TwoStream\_animation.mp4)

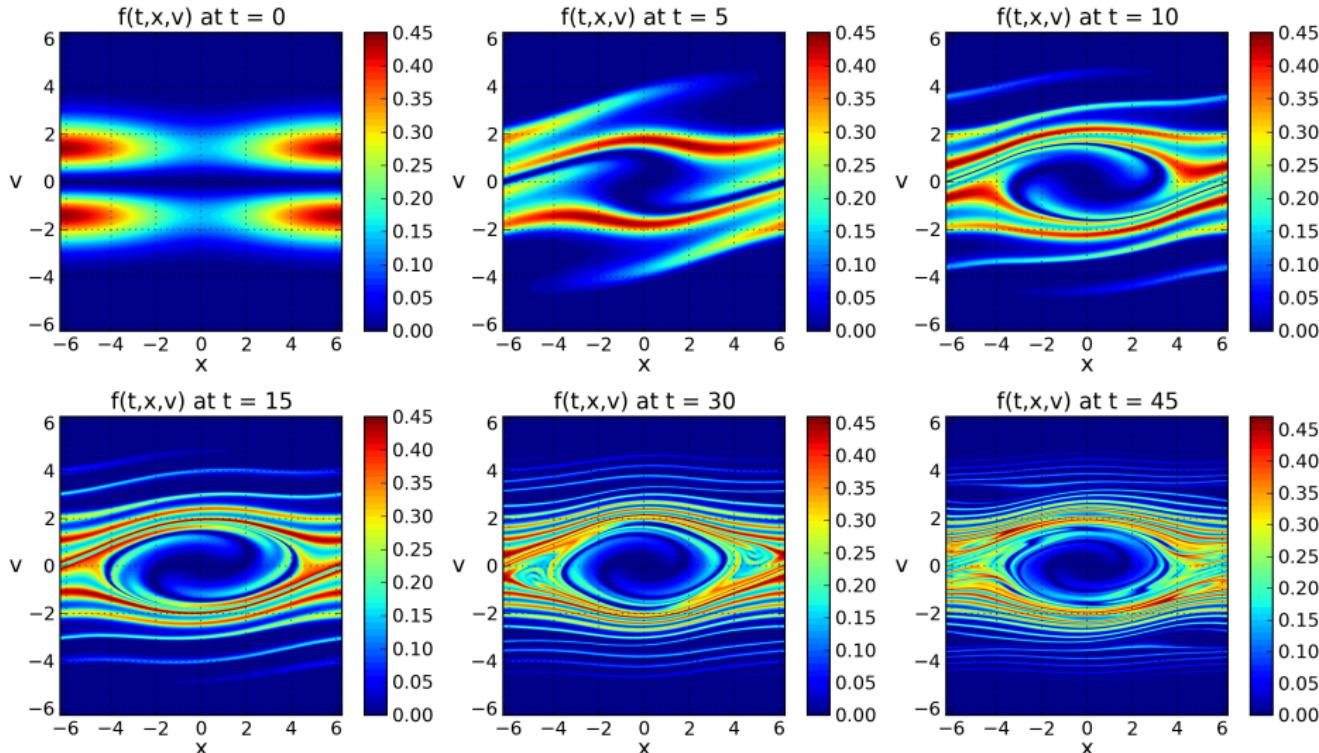
COURANT NUMBER

$C_x \approx 64$

$C_v \approx 20$



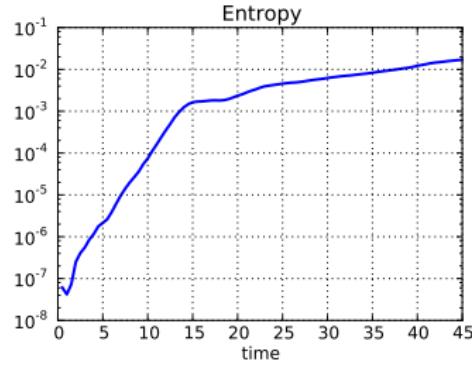
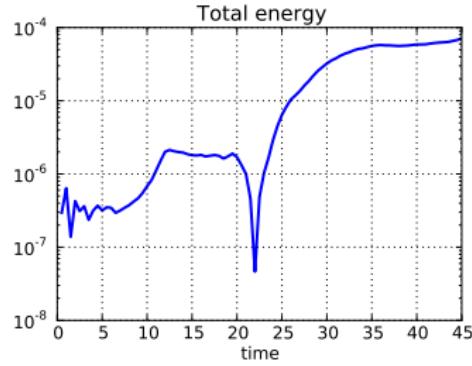
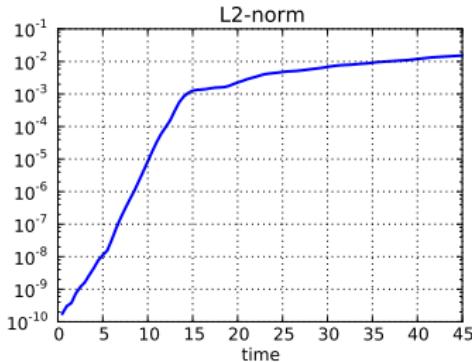
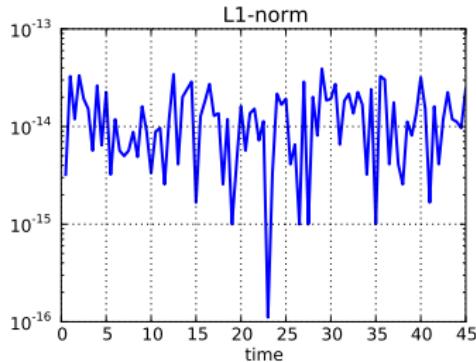
# Vlasov-Poisson: two-stream instability (2)



$N_x = 256$     $N_v = 512$     $\Delta t = 0.5$    90 steps    $[C_x \approx 64 \quad C_v \approx 20]$

# Vlasov-Poisson: two-stream instability (3)

Relative errors in conserved quantities



GRID  
 $N_x = 256$   
 $N_v = 512$

TIME STEPPING  
 $\Delta t = 0.5$   
90 steps

COURANT NO.  
 $C_x \approx 64$   
 $C_v \approx 20$



# Conclusions and Outlook

## RECALL:

- Multi-dimensional mesh-based solution of Boltzmann's eq. for weakly collisional plasmas;
- In standard Eulerian codes, memory requirement may be too large (large mesh + RK storage) and time steps may be too small (CFL restriction);
- Splitting ballistic and collision operators allows semi-Lagrangian algorithms (no CFL limit);
- Splitting configuration-advection and velocity-advection permits one to use very accurate constant advection solvers (coarser mesh, no more RK storage).

## SUMMARY:

- Convected Scheme (CS) is semi-Lagrangian algorithm, mass and positivity preserving;
- Constant advection CS extended to arbitrarily high order (22nd-order version with FFTs);
- High order (4th/6th) Runge-Kutta-Nyström operator splitting guarantees energy stability;
- Tested with standard benchmarks for 1D-1V Vlasov-Poisson system;
- Error in total energy conservation bounded until solution is spatially resolved.

## WORK IN PROGRESS:

- Implement absorbing boundary conditions (wall recombination);
- Couple to simple collision operator (electron scattering on neutrals);
- Extend to higher dimensions (1D-2V, 2D-3V).

## Acknowledgments

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- MICHIGAN STATE UNIVERSITY FOUNDATION  
SPG-RG100059
  - AIR FORCE OFFICE OF SCIENTIFIC RESEARCH (AFOSR)  
FA9550-11-1-0281, FA9550-12-1-0343, FA9550-12-1-0455
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DMS-1115709