Exponential methods for solving hyperbolic problems with application to kinetic equations

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Outline

- Motivation for Vlasov-Poisson equations
- 2 Linear analysis
 - Lawson methods
 - Exponential Runge-Kutta methods
- 3 Numerical simulation: Vlasov-Poisson equations
- 4 Numerical simulation: drift-kinetic equations
- Conclusion

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Vlasov-Poisson equations $1D \times 1D$

Our model: a non-linear transport in $(x, v) \in \Omega \times \mathbb{R}$ of an electron density distribution f = f(t, x, v):

$$\begin{cases} \partial_t f + v \partial_x f + E \partial_v f = 0 \\ \partial_x E = \int_{\mathbb{R}} f \, dv - 1 \end{cases}$$

Motivation:

- We want high order methods in (x, v)
- We want high order methods in time t:
 - Splitting methods: could have a lot of steps
 - Runge-Kutta methods: stability constraints (CFL condition)
 - The most restrictive CFL condition is associated with the linear part $(\partial_t f + v \partial_x f = 0)$
- →We want to propose a compromise: exponential integrators.

Vlasov-Poisson equations $1D \times 1D$

Fourier transform in x direction of Vlasov, amenable to exponential integrators:

$$\partial_t \hat{f} + ikv\hat{f} + \widehat{E\partial_v f} = 0$$

Vlasov is of the form:

$$\dot{u} = iau + F(u)$$

Variation of constant: $\partial_t(e^{-iat}u)=e^{-iat}F(u)$. No more CFL in x of the form $\Delta t \leq \sigma \frac{\Delta x}{v_{\max}}$ with $[-v_{\max},v_{\max}] \equiv \mathbb{R}$. Time integration:

$$u(t_n + \Delta t) = \exp(ia\Delta t)u(t_n) + \int_0^{\Delta t} \exp(ia(\Delta t - s))F(u(t_n + s)) ds$$

with $\Delta t > 0$, $t_n = n\Delta t$ with $n \in \mathbb{N}$ Linear part is exact!

Idea of exponential integrators

2 classes of methods:

exponential Runge-Kutta: solve exactly what we can, and interpolate the rest. For example first order exponential Euler method:

$$u(t_n+\Delta t)pprox u^{n+1}=e^{-ia\Delta t}u^n+\Delta t arphi_1(ia\Delta t) F(u^n)$$
 where $arphi_1(z)=rac{e^z-1}{z}$



Hochbruck and Ostermann (2010)

Lawson: Change of variable: $v(t) = e^{-iat}u(t)$, we solve with a RK method: $\dot{v} = \tilde{F}(t, v) = e^{-iat} F(e^{iat} v(t))$ For example, Lawson Euler method:

$$v(t_n + \Delta t) \approx v^{n+1} = v^n + \Delta t e^{-iat_n} F(e^{iat_n} v^n)$$

or as an expression of u:

$$u^{n+1} = e^{-ia\Delta t}u^n + \Delta t e^{ia\Delta t}F(u^n)$$



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Reminder of stability tools

If we want to study stability of:

$$\partial_t u + \partial_x u = 0$$

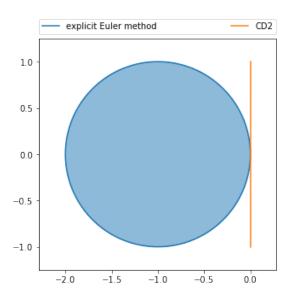
with centered scheme (CD2) $(\partial_X u)_j \approx \frac{1}{2\Delta_X}(u_{j+1} - u_{j-1})$. After a Fourier transform (von Neumann analysis):

$$\dot{u} + i \frac{\sin(k\Delta x)}{\Delta x} u = 0$$

Explicit Euler method in time: we have to stretch eigenvalues (or Fourier symbol) of CD2 into explicit Euler stability domain.

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Reminder of stability tools



From linear Vlasov equation to toy model

Linear Vlasov equation:

$$\partial_t f + a \partial_x f + b \partial_v f = 0$$

Fourier transform in x, CD2 in v plus a Fourier transform in v, formally:

$$\frac{\mathrm{d}f}{\mathrm{d}t} + iakf + b\frac{i\sin(\varphi)}{\Delta x}f = 0$$

Toy model:

$$\dot{u} + iau + \lambda u = 0$$

with $a \in \mathbb{R}$, $\lambda \in \mathbb{C}$ (diffusive scheme for example).

 λ is the Fourier symbol (or eigenvalues) of FD method to approximate $\partial_{\nu} f$.

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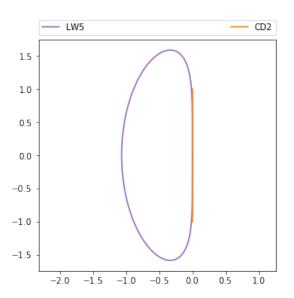
In v direction we use a FD method:

- CD2 (centered difference of order 2): $(\partial_{\nu} f)(v_j) \approx \frac{f_{j+1} f_{j-1}}{2\Delta \nu}$
- WENO5 (weighted essentially non-oscillatory of order 5):
 - WENO5: non linear scheme: Von Neumann analysis
 - LW5 (linearized WENO5): linear scheme (this is Lagrange interpolation of order 5)

$$(\partial_{\nu}f)(\nu_{j}) \approx \frac{1}{\Delta\nu} \left(-\frac{1}{30}f_{j-3} + \frac{1}{4}f_{j-2} - f_{j-1} + \frac{1}{3}f_{j} + \frac{1}{2}f_{j+1} - \frac{1}{20}f_{j+2} \right)$$

- Wang and Spiteri (2007)
- Motamed, Macdonald, and Ruuth (2010)

Fourier symbols



Lawson methods stability domain

For our toy model:

$$\dot{u} = iau + \lambda(u)$$

Change of variable: $v(t) = e^{-iat}u(t)$

$$\dot{v} = e^{-iat} \lambda e^{iat} v$$

Apply a Runge-Kutta method to compute stability function of Lawson method:

$$v^{n+1} = \underbrace{p(\lambda \Delta t)}_{\text{stability function of RK}} v^n$$

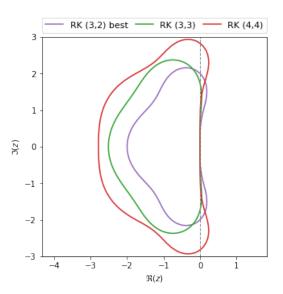
i.e.:

stability function of Lawson

$$u^{n+1} = \overbrace{p(\lambda \Delta t)e^{-ia\Delta t}} u^n$$

Stability domain: $\mathcal{D}=\{z\in\mathbb{C}, |p(z)|\leq 1\}$ of Lawson method is **the same** as the underlying Runge-Kutta method **because** $ia\in i\mathbb{R}$

Considered Lawson(RK(s, p)) methods



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Lawson methods – CD2

For stability between a Lawson method and CD2, we solve:

$$|p(iy)| = 1, \quad y \in \mathbb{R}$$

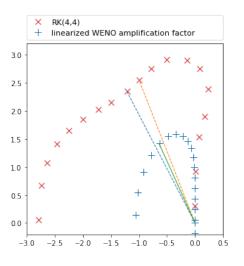
Methods	Lawson($RK(3,2)$ best)	Lawson $(RK(3,3))$	Lawson $(RK(4,4))$
y _{max}	2	$\sqrt{3}$	$2\sqrt{2}$

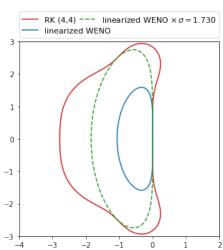
Table: CFL number for some Lawson schemes



Baldauf (2008)

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Lawson methods – LW5: CFL estimates

Methods	Lawson($RK(3,2)$ best)	Lawson $(RK(3,3))$	Lawson $(RK(4,4))$
σ	1.344	1.433	1.73

Table: CFL number for some Lawson schemes.



Motamed, Macdonald, and Ruuth (2010)



Lunet et al. (2017)

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Exponential Runge-Kutta methods

$$\dot{u} = iau + F(u)$$

Example on ExpRK(2,2):

$$u^{(1)} = e^{-ia\Delta t}u^n - \Delta t\varphi_1 F(u^n)$$

$$u^{n+1} = e^{-ia\Delta t}u^n - \Delta t \left[(\varphi_1 - \varphi_2)F(u^n) + \varphi_2 F(u^{(1)}) \right]$$

Stability function becomes:

$$p_{\text{ExpRK}(2,2)}(z) = \frac{1}{2}\varphi_1\varphi_{1,2}z^2 + (\varphi_1 + i\frac{\varphi_1\varphi_{1,2}}{2}a)z + 1 + i\varphi_1a$$

Stability domain depends of $a\Delta t...X$

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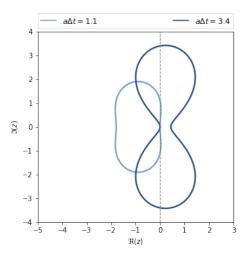


Figure: Stability domain of ExpRK(2,2) for $a\Delta t \in \{1.1, 3.4\}$

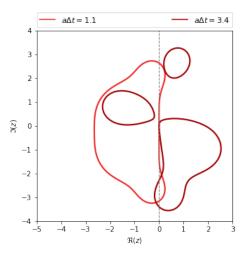


Figure: Stability domain of Cox-Matthews for $a\Delta t \in \{1.1, 3.4\}$

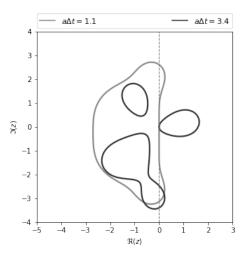


Figure: Stability domain of Krogstad for $a\Delta t \in \{1.1, 3.4\}$

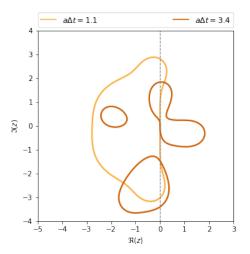


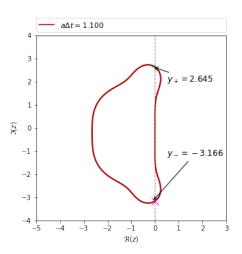
Figure: Stability domain of Hochbruck–Ostermann for $a\Delta t \in \{1.1, 3.4\}$

Stability domain informations

Fourier symbol must fit in the stability domain of ExpRK method **for all** values of $a\Delta t \in \mathbb{R}$.

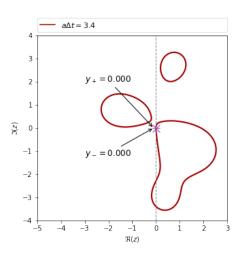
- Impossible with WENO5 (LW5 Fourier symbol)
 - \rightarrow Numerical test: unstable in very short time
- ✓ Singleton $\{0\}$ is alway in stability domain of ExpRK method for each values of $a\Delta t$
 - \rightarrow We can try to stabilize CD2
 - SPOILER: CFL is equal to zero

ExpRK – CD2



 $y_{
m max}^{
m exp}=\min(y_+,|y_-|)$ the largest value to stretch i[-1,1] into the stability domain at $a\Delta t$

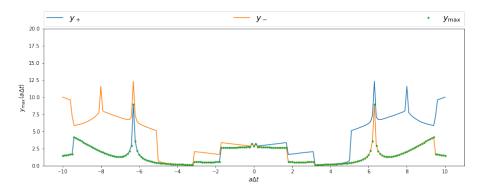
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ExpRK - CD2: CFL number

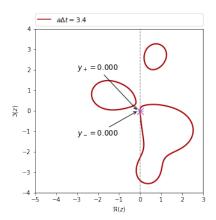


CFL $y_{\text{max}} = \min_{a \Delta t} y_{\text{max}}^{exp}$ is still 0...

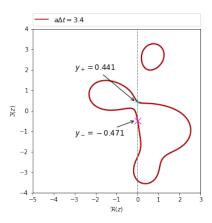
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ExpRK - CD2: Relaxed CFL condition

$$\mathcal{D}_{\varepsilon} = \{ z \in \mathbb{C}, |p(z)| \le 1 + \varepsilon \}$$

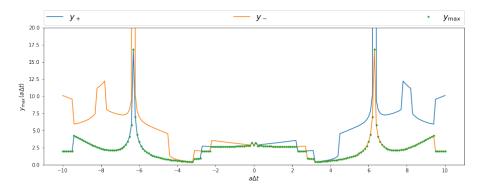


Cox-Matthews stability domain, relaxation $\varepsilon = 0$



Cox-Matthews stability domain, relaxation $\varepsilon = 10^{-2}$

ExpRK - CD2: Relaxed CFL condition



Relaxed CFL $y_{\rm max}(\varepsilon=10^{-2})\approx 0.450 \neq 0$! (but unstable in theory)

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ExpRK – CD2: Relaxed CFL estimates

Methods	ExpRK22	Krogstad	Cox-Matthews	Hochbruck
				–Ostermann
$y_{\sf max}(\varepsilon=10^{-3})$	0.300	0.100	0.150	0.250
$y_{max}(\varepsilon = 10^{-2})$	0.551	0.200	0.450	0.501
$y_{\sf max}(\varepsilon=10^{-1})$	1.001	0.601	1.351	1.702

Table: CFL number, assuming the relaxed stability constraint, for some exponential integrators.

It's unstable in theory, in practice, number of iterations is finished, so amplification is controlled.

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$$\begin{cases} \partial_t f + v \partial_x f + E \partial_v f = 0 \\ \partial_x E = \int_{\mathbb{R}} f \, dv - 1 \end{cases}$$

Numerical tools:

- FFT in x direction
- CD2 or WENO5 in v direction
- Lawson(RK(s, p)) or ExpRK method in time t

CFL:
$$\Delta t_n \leq \frac{C\Delta v}{||E^n||_{\infty}} \leq \frac{C\Delta v}{\max_n ||E^n||_{\infty}}$$
 where $C = y_{\max}$ or σ from the linear theory.

We can choose:
$$\Delta t = \min\left(0.1, \frac{C\Delta v}{\max_n ||E^n||_{\infty}}\right)$$

Landau damping

$$f(t=0,x,v)=f_0(x,v)=\frac{1}{\sqrt{2\pi}}e^{-\frac{v^2}{2}}(1+0.001\cos(0.5x))$$

 $x \in [0, 4\pi], \ v \in [-8, 8], \ N_x = 81, \ N_v = 128$

Because of damping:

$$\max_{n}||E^n||_{\infty}=||E^0||_{\infty}$$

So, we choose $\Delta t = 0.1$ (with $\Delta t = 100$ it is still stable!)

Landau damping: numerical results

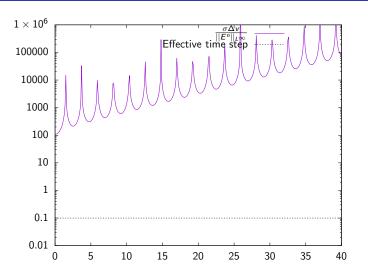


Figure: Landau damping test: time history of the CFL condition (semi-log scale).

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Landau damping: numerical results

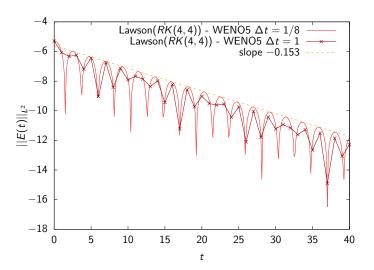


Figure: Landau damping test: time history of $||E(t)||_{L^2}$ (semi-log scale) obtained with Lawson(RK(4,4)) and WENO5 with $\Delta t = 1/8$ and $\Delta t = 1$.

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Bump on Tail (BoT)

$$f(t=0,x,v) = \left[\frac{0.9}{\sqrt{2\pi}}e^{-\frac{v^2}{2}} + \frac{0.2}{\sqrt{2\pi}}e^{-2(v-4.5)^2}\right](1+0.001\cos(0.5x))$$

 $x \in [0, 20\pi], v \in [-8, 8], N_x = 135, N_v = 256$ Numerical estimation of $\max_n ||E^n||_{\infty} \approx 0.6$, we choose $\Delta t = \frac{C\Delta v}{0.6}$

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BoT: numerical results

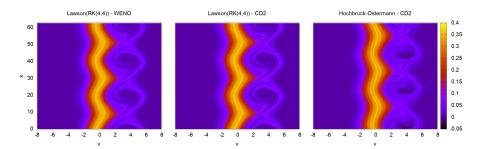


Figure: Distribution function at time t = 40 as a function of x and v for Lawson(RK(4,4)) + WENO5 (left), Lawson(RK(4,4)) + centered scheme (center), Hochbruck–Ostermann + centered scheme (right).

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BoT: numerical results

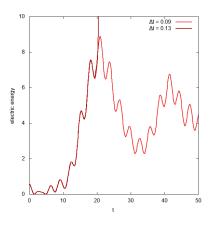


Figure: Illustration of the accuracy of the CFL estimate obtained from the linear theory. History of electric energy with Lawson(RK(4,4)) + WENO5

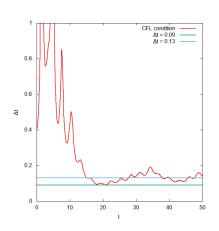
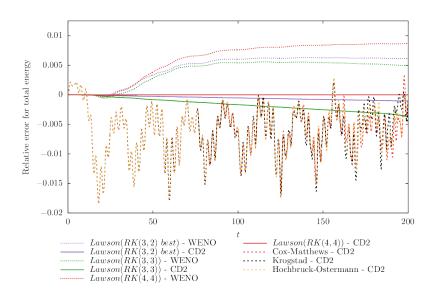


Figure: History of CFL condition for Lawson(RK(4,4)) + WENO5 case

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BoT: numerical results



$$\max_n ||E^n||_{\infty}$$
 is not accessible in practice.

To capture correctly the phenomena involved in the bump on tail test, we take the following time step size:

$$\Delta t_n = \min\left(0.1, \frac{C\Delta v}{||E^n||_{\infty}}\right)$$

with $C = y_{\text{max}}$ or σ from the linear theory.

→Good estimate in practice for Lawson methods.

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Drift-Kinetic equations

$$f = f(t, r, \theta, z, v)$$

$$\underbrace{\left\{ \underbrace{\partial_t f + v \partial_z f}_{\text{linear part}} \underbrace{-\frac{\partial_\theta \phi}{r} \partial_r f + \frac{\partial_r \phi}{r} \partial_\theta f - \partial_z \phi \partial_v f}_{\text{non linear part}} = 0 \right.}_{\text{non linear part}}$$

$$- \left[\partial_r^2 \phi + \left(\frac{1}{r} + \frac{\partial_r n_0(r)}{n_0(r)} \right) \partial_r \phi + \frac{1}{r^2} \partial_\theta^2 \phi \right] + \frac{1}{T_e(r)} (\phi - \langle \phi \rangle) = \frac{1}{n_0(r)} \int_{\mathbb{R}} f \, \mathrm{d}v - 1$$

$$(r, \theta, z, v) \in [0.1, 14.5] \times [0, 2\pi] \times [0, L] \times \mathbb{R}$$

After a Fourier transform in z, formally, the equation is still of the form of:

$$\partial_t f + ikvf + F(f) = 0$$

Compatible with all previous time integrators.

This is more complicated to use linear stability analysis, we use an other adaptive time step method.

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Adaptive time step size (error estimate)

For adaptive time step size with any time integrator φ :

$$f^{n+1} = \varphi_{\Delta t_n}(f^n)$$
 ; $\tilde{f}^{n+1} = \varphi_{\Delta t_n/2} \circ \varphi_{\Delta t_n/2}(f^n)$

Richardson extrapolated numerical solution of the method of order *p*:

$$f_R^{n+1} = \frac{2^{p+1}\tilde{f}^{n+1} - f^{n+1}}{2^{p+1} + 1}$$

estimate of the local error:

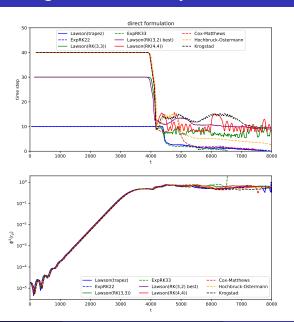
$$e_{n+1} = ||f_R^{n+1} - f^{n+1}||_{L^{\infty}} + \mathcal{O}(\Delta t_n^{p+2})$$

If $e_{n+1} > \text{tol}$: we reject the step and start again from time t_n . Else we determine the new time step size:

$$\Delta t_{new} = s \Delta t_n \left(rac{\mathsf{tol}}{e_{n+1}}
ight)^{1/(p+1)}$$

s = 0.8 is safety factor.

Ion temperature gradient instability: numerical results



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Summary

- Better understanding on stability of Lawson or ExpRK methods in transport equations
- Python script with sympy to compute estimates of CFL of Lawson –
 CD2, Lawson WENO (5 or 3) or ExpRK CD2 (with relaxing CFL)
- An adaptive time step size which works with any time integrators

Future works

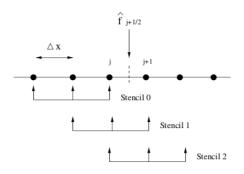
- We can improve method with an embedded Runge-Kutta method (Dormand-Prince method, used in ode45 of Matlab)
- Compare performance between exponential integrators and splitting methods (same stages/step, same order?)
- Use semi-Lagrangian method to remove dependency on periodic space (Fourier transform)

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Thank you for your attention

WENO5 method

Weighted **E**ssentially **N**on-**O**scillatory method of order 5: 3 estimates on 3 different stencils weighted with nonlinear weights.



3 steps:

- Indicator of smoothness
- Weights
- Flux

Indicator of smoothness β_i^{\pm}

To approximate $\partial_x f(u)$:

Split *f* as:

$$f(u) = f^{+}(u) + f^{-}(u)$$
 , $\frac{df^{+}}{du} \ge 0$ et $\frac{df^{-}}{du} \le 0$

Indicators of smoothness:

$$\beta_i^{\pm} \leftarrow (f_{[j-2,j+3]}^{\pm}) \ , \ i = 0, 1, 2$$

Approximations of derivatives of order 1 and 2 on 3 stencils.

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$$\begin{split} \beta_0^+ &= \frac{13}{12} \left(f_{j-2}^+ - 2 f_{j-1}^+ + f_j^+ \right)^2 + \frac{1}{4} \left(f_{j-2}^+ - 4 f_{j-1}^+ + 3 f_j^+ \right)^2 \\ \beta_1^+ &= \frac{13}{12} \left(f_{j-1}^+ - 2 f_j^+ + f_{j+1}^+ \right)^2 + \frac{1}{4} \left(f_{j-1}^+ - f_{j+1}^+ \right)^2 \\ \beta_2^+ &= \frac{13}{12} \left(f_j^+ - 2 f_{j+1}^+ + f_{j+2}^+ \right)^2 + \frac{1}{4} \left(3 f_j^+ - 4 f_{j+1}^+ + f_{j+2}^+ \right)^2 \\ \beta_0^- &= \frac{13}{12} \left(f_{j+1}^- - 2 f_{j-2}^- + f_{j+3}^- \right)^2 + \frac{1}{4} \left(3 f_{j-1}^- - 4 f_{j-2}^- + f_{j+3}^- \right)^2 \\ \beta_1^- &= \frac{13}{12} \left(f_j^- - 2 f_{j+1}^- + f_{j+2}^- \right)^2 + \frac{1}{4} \left(f_j^- - f_{j+2}^- \right)^2 \\ \beta_2^- &= \frac{13}{12} \left(f_{j-1}^- - 2 f_j^- + f_{j+1}^- \right)^2 + \frac{1}{4} \left(f_{j-1}^- - 4 f_j^- + 3 f_{j+1}^- \right)^2 \end{split}$$

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WENO5 method

Weights w_i^{\pm}

Unnormalized weights:

$$\alpha_i^{\pm} \leftarrow \frac{\gamma_i}{(\epsilon + \beta_i^{\pm})^2} , \ \gamma_i \in \mathbb{R}_+^* : \sum_k \gamma_k = 1$$

where $\gamma_0=\frac{1}{10}, \gamma_1=\frac{6}{10}, \gamma_2=\frac{3}{10}.$ Parameter $\epsilon=10^{-6}$ Linearized weights (LW5): $\alpha_i^\pm=\gamma_i+\mathcal{O}(\Delta x^2)$

Normalized weights:

$$w_i^{\pm} \leftarrow \frac{\alpha_i^{\pm}}{\sum_k \alpha_k^{\pm}}$$

Flux $f_{i+\frac{1}{2}}^{\pm}$

$$\begin{split} f^+_{j+\frac{1}{2}} \leftarrow w^+_0 \left(\frac{2}{6} f^+_{j-2} - \frac{7}{6} f^+_{j-1} + \frac{11}{6} f^+_j \right) + w^+_1 \left(-\frac{1}{6} f^+_{j-1} + \frac{5}{6} f^+_j + \frac{2}{6} f^+_{j+1} \right) \\ + w^+_2 \left(\frac{2}{6} f^+_j + \frac{5}{6} f^+_{j+1} - \frac{1}{6} f^+_{j+2} \right) \end{split}$$

$$\begin{split} f_{j+\frac{1}{2}}^{-} \leftarrow w_{2}^{-} \left(-\frac{1}{6} f_{j-1}^{-} + \frac{5}{6} f_{j}^{-} + \frac{2}{6} f_{j+1}^{-} \right) + w_{1}^{-} \left(\frac{2}{6} f_{j}^{-} + \frac{5}{6} f_{j+1}^{-} - \frac{1}{6} f_{j+2}^{-} \right) \\ + w_{0}^{-} \left(\frac{11}{6} f_{j+1}^{-} - \frac{7}{6} f_{j+2}^{-} + \frac{2}{6} f_{j+3}^{-} \right) \end{split}$$

$$\boxed{(\partial_x f(u))_j \approx \frac{1}{\Delta x} \left[(f_{j+\frac{1}{2}}^+ - f_{j-\frac{1}{2}}^+) + (f_{j+\frac{1}{2}}^- - f_{j-\frac{1}{2}}^-) \right]}$$

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More Lawson methods

We are interested in the numerical cost $\frac{\Delta t}{s}$ of RK(s,n). To compare each time integrator, we compute total energy in Vlasov-Poisson system:

$$H(t) = \int_{\Omega} \int_{\mathbb{R}} v^2 f \, dx dv + \int_{\Omega} E^2 \, dx$$

which is preserved in time. We propose to select the best method by considering:

$$h_{s,n}: \frac{\Delta t}{s} \mapsto \left| \left| \frac{H(t) - H(0)}{H(0)} \right| \right|_{\infty}$$

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More Lawson methods

