Tutorial on

Semi-Lagrangian schemes

Roberto Ferretti

BIRS workshop 11w5086



Dipartimento di Matematica Universita' di Roma Tre

- Some history
- Basic ideas and building blocks for SL schemes
- Convergence analysis for the linear problem
- Construction of Semi-Lagrangian schemes for convex HJ equations
- Convergence analysis for the nonlinear problem

Some history

- Semi–Lagrangian schemes: introduced as first–order schemes by Courant, Isaacson and Rees (CPAM, '52)
- Numerical Weather Prediction streamline: Wiin-Nielsen (Tellus, '59), Robert (Atmosphere-Ocean, '81), Staniforth, Côté, Smolarkiewicz...
- Plasma physics streamline: Cheng–Knorr ('76), Bertrand–Izzo, Besse–Mehrenberger,...

In the first developments it had not yet been realized that the possible advantage of SL schemes over conventional difference schemes was to be able to work at large Courant numbers.

This feature has become important in NWP problems, in which an orthogonal grid would have forced a conventional scheme to adopt prohibitively small time steps because of the singularity on the poles.

A further analysis shows that large Courant numbers cause the scheme to be less diffusive.

index

Basic ideas and building blocks for SL schemes

For simplicity, we will discuss SL schemes focusing on the model problem

$$\begin{cases} u_t(x,t) + f(x,t) \cdot Du(x,t) = 0, & (x,t) \in \mathbb{R}^d \times \mathbb{R} \\ u(x,0) = u_0(x) & x \in \mathbb{R}^d. \end{cases}$$

posed on the whole of \mathbb{R}^d .

Basic ideas and building blocks for SL schemes

For simplicity, we will discuss SL schemes focusing on the model problem

$$\begin{cases} u_t(x,t) + f(x,t) \cdot Du(x,t) = 0, & (x,t) \in \mathbb{R}^d \times \mathbb{R} \\ u(x,0) = u_0(x) & x \in \mathbb{R}^d. \end{cases}$$

posed on the whole of \mathbb{R}^d .

• We avoid the treatment of boundary conditions

Basic ideas and building blocks for SL schemes

For simplicity, we will discuss SL schemes focusing on the model problem

$$\begin{cases} u_t(x,t) + f(x,t) \cdot Du(x,t) = 0, & (x,t) \in \mathbb{R}^d \times \mathbb{R} \\ u(x,0) = u_0(x) & x \in \mathbb{R}^d. \end{cases}$$

posed on the whole of \mathbb{R}^d .

- We avoid the treatment of boundary conditions
- We treat separately and more explicitly the case of constant speed

Any large time-step technique (in particular, Semi-Lagrangian approximations) stem from the method of characteristics. Let a system of characteristic curves y(x, s; t) for the model equation be defined by:

$$\begin{cases} \frac{d}{dt}y(x,t;s) = f(y(x,t;s),s), \\ y(x,t;t) = x, \end{cases}$$

Then, the solution is constant along such trajectories, which means that the following representation formula

$$u(y(x,t;t+\tau),t+\tau) = u(x,t).$$

holds for the solution u.

Writing the representation formula at a node x_i and with $\tau = -\Delta t$, we have the time-discrete version

$$u(x_i,t) = u(y(x_i,t;t-\Delta t),t-\Delta t).$$

Its numerical discretization is obtained by combining:

Writing the representation formula at a node x_i and with $\tau = -\Delta t$, we have the time-discrete version

$$u(x_i,t) = u(y(x_i,t;t-\Delta t),t-\Delta t).$$

Its numerical discretization is obtained by combining:

• A numerical technique to integrate backwards the ODE of characteristics Writing the representation formula at a node x_i and with $\tau = -\Delta t$, we have the time-discrete version

$$u(x_i,t) = u(y(x_i,t;t-\Delta t),t-\Delta t).$$

Its numerical discretization is obtained by combining:

• A numerical technique to integrate backwards the ODE of characteristics

• A reconstruction to approximate the value $u(y(x_i, t; t - \Delta t), t - \Delta t)$, since in general the foot of the characteristic $y(x_i, t; t - \Delta t)$ does not coincide with any grid point.

$$v_i^{n+1} = I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n))$$

$$v_i^{n+1} = I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n))$$

• v_i^{n+1} is the numerical solution computed at (x_i, t_{n+1})

$$v_i^{n+1} = I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n))$$

- v_i^{n+1} is the numerical solution computed at (x_i, t_{n+1})
- $X^{\Delta}(x_i, t_{n+1}; t_n)$ approximates the characteristic through (x_i, t_{n+1})

$$v_i^{n+1} = I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n))$$

- v_i^{n+1} is the numerical solution computed at (x_i, t_{n+1})
- $X^{\Delta}(x_i, t_{n+1}; t_n)$ approximates the characteristic through (x_i, t_{n+1})
- $I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n)) = \sum_j v_j^n \psi_j(X^{\Delta}(x_i, t_{n+1}; t_n))$ is the interpolation computed at $(X^{\Delta}(x_i, t_{n+1}; t_n), t_n)$

Approximation of characteristics X^{Δ} : typically, by one-step or multisptep techniques

Approximation of characteristics X^{Δ} : typically, by one-step or multisptep techniques

• The advection field is known (in relevant problems) only at spacetime nodes **Approximation of characteristics** X^{Δ} : typically, by one-step or multisptep techniques

- The advection field is known (in relevant problems) only at spacetime nodes
- Need to avoid intermediate times, as well as to interpolate the advecting field among space grid nodes

1st example: the Euler scheme only needs informations at the time t.

$$y(x_i, t; t - \Delta t) \approx X^{\Delta}(x_i, t; t - \Delta t) = x_i - \Delta t f(x_i, t)$$

- This is the classical choice of the Courant–Isaacson–Rees scheme
- In general, it leads to a poor time approximation (1st order)

$$X^{\Delta}(x_i, t; t - \Delta t) = x_i - \frac{\Delta t}{2} \left[f(x_i, t) + \tilde{f}(x_i - \Delta t f(x_i, t), t - \Delta t) \right]$$

$$X^{\Delta}(x_i, t; t - \Delta t) = x_i - \frac{\Delta t}{2} \left[f(x_i, t) + \tilde{f}(x_i - \Delta t f(x_i, t), t - \Delta t) \right]$$

• $\tilde{f}(\xi, t - \Delta t)$ is an interpolate of the node values of $f(x_j, t - \Delta t)$, computed at the point ξ

$$X^{\Delta}(x_i, t; t - \Delta t) = x_i - \frac{\Delta t}{2} \left[f(x_i, t) + \tilde{f}(x_i - \Delta t f(x_i, t), t - \Delta t) \right]$$

• $\tilde{f}(\xi, t - \Delta t)$ is an interpolate of the node values of $f(x_j, t - \Delta t)$, computed at the point ξ

• No interpolation is needed if f is explicitly known

$$X^{\Delta}(x_i, t; t - \Delta t) = x_i - \frac{\Delta t}{2} \left[f(x_i, t) + \tilde{f}(x_i - \Delta t f(x_i, t), t - \Delta t) \right]$$

• $\tilde{f}(\xi, t - \Delta t)$ is an interpolate of the node values of $f(x_j, t - \Delta t)$, computed at the point ξ

- No interpolation is needed if f is explicitly known
- The approximation is second-order with respect to Δt

Numerical reconstruction of the value $u(y(x_j, t; t - \Delta t), t - \Delta t)$:

Linear:

- Symmetric Lagrange interpolation (most common)
- Finite Element interpolation, sparse grids, Chebyshev grids,...

Nonlinear:

• Non-Oscillatory (ENO/WENO) interpolation, monotone Hermite interpolations,...

Symmetric Lagrange interpolation is performed using a symmetric stencil of points around *x*:



stencils of interpolation (linear, cubic and quintic Lagrange)



region of interpolation (\mathbb{P}_3 finite elements and cubic Lagrange)

It might be convenient to consider a set of basis functions in which symmetric Lagrange interpolation may be expressed It might be convenient to consider a set of basis functions in which symmetric Lagrange interpolation may be expressed

• This basis function ψ_j is obtained by interpolating the sequences e_j , i.e., sequences which are everywhere zero except at the node x_j It might be convenient to consider a set of basis functions in which symmetric Lagrange interpolation may be expressed

• This basis function ψ_j is obtained by interpolating the sequences e_j , i.e., sequences which are everywhere zero except at the node x_j

• On a uniform grid, a basis function ψ_j can be written in terms of a reference basis function ψ :

$$\psi_j(\xi) = \psi\left(\frac{\xi}{\Delta x} - j\right)$$

(obtained reconstructing e_0 on a grid with $\Delta x = 1$)

When this procedure is applied to a Lagrange reconstruction of odd order r, the reference fasis function has the form:

$$\psi(\xi) = \begin{cases} \prod_{\substack{k \neq 0, k = -[r/2] \\ k \neq 0, k = -[r/2]}} \frac{\xi - k}{-k} & \text{if } 0 \le \xi \le 1 \\ \vdots & \vdots \\ \prod_{\substack{k=1 \\ k=1}}^{r} \frac{\xi - k}{-k} & \text{if } [r/2] \le \xi \le [r/2] + 1 \\ 0 & \text{if } \xi > [r/2] + 1 \end{cases}$$

and extended by symmetry for $\xi < 0$.

When this procedure is applied to a Lagrange reconstruction of odd order r, the reference fasis function has the form:

$$\psi(\xi) = \begin{cases} \prod_{\substack{k \neq 0, k = -[r/2] \\ k \neq 0, k = -[r/2]}} \frac{\xi - k}{-k} & \text{if } 0 \le \xi \le 1 \\ \vdots & \vdots \\ \prod_{\substack{k=1 \\ k=1}}^{r} \frac{\xi - k}{-k} & \text{if } [r/2] \le \xi \le [r/2] + 1 \\ 0 & \text{if } \xi > [r/2] + 1 \end{cases}$$

and extended by symmetry for $\xi < 0$.

• The interpolation error is $O(\Delta x^{r+1})$ for smooth functions



The reference basis functions ψ for \mathbb{P}_1 and cubic interpolation

index

Convergence analysis for the linear problem

To prove **consistency**, we need to compare the scheme:

$$v_i^{n+1} = I[V^n](X^{\Delta}(x_i, t_{n+1}; t_n))$$

with the representation formula:

$$u(x_i, t_{n+1}) = u(y(x_i, t_{n+1}; t_n), t_n).$$

assuming that u is a smooth solution and that $v_j^n = u(x_j, t_n)$.

We also assume to have a general approximation of order \boldsymbol{p} in time and \boldsymbol{r} in space

It turns out that the local truncation error is estimated as:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \left(\Delta t^p + \frac{\Delta x^{r+1}}{\Delta t} \right)$$

It turns out that the local truncation error is estimated as:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \left(\Delta t^p + \frac{\Delta x^{r+1}}{\Delta t} \right)$$

• The term Δt^p accounts for the error in the computation of characteristics

It turns out that the local truncation error is estimated as:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \left(\Delta t^p + \frac{\Delta x^{r+1}}{\Delta t} \right)$$

• The term Δt^p accounts for the error in the computation of characteristics

• The term $\frac{\Delta x^{r+1}}{\Delta t}$ accounts for the error generated by the accumulation of interpolation errors
It turns out that the local truncation error is estimated as:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \left(\Delta t^p + \frac{\Delta x^{r+1}}{\Delta t} \right)$$

- The term Δt^p accounts for the error in the computation of characteristics
- The term $\frac{\Delta x^{r+1}}{\Delta t}$ accounts for the error generated by the accumulation of interpolation errors
- There exists an optimal $\Delta x/\Delta t$ balance which maximizes the consistency rate

To prove **stability**, we restrict for simplicity to the equation in the constant coefficient form:

$$u_t + cu_x = 0.$$

Here, we have assumed that the advection has constant speed c, so that $X^{\Delta}(x_i, t_{n+1}; t_n) = x_i - c\Delta t$ and the SL scheme has the form

$$v_i^{n+1} = I[V^n](x_i - c\Delta t).$$

We are in the typical framework of Von Neumann analysis, and in fact it is possible to prove by Fourier analysis arguments that the scheme is stable.



Amplitude of the amplification factors λ for cubic interpolation

We will rather follow the line of proving stability by equivalence with a stable scheme, in this case the Lagrange–Galerkin scheme which has the form:

$$\int_{\mathbb{R}} v_{\Delta}^{n+1}(\xi) \phi_i(\xi) d\xi = \int_{\mathbb{R}} v_{\Delta}^n(\xi - c\Delta t) \phi_i(\xi) d\xi$$

that is, writing the numerical solution as $v_{\Delta}^k(x) = \sum_j v_j^k \phi_j(x)$,

$$\sum_{j} v_j^{n+1} \int_{\mathbb{R}} \phi_j(\xi) \phi_i(\xi) d\xi = \sum_{j} v_j^n \int_{\mathbb{R}} \phi_j(x_i - c\Delta t) \phi_i(\xi) d\xi$$

We will rather follow the line of proving stability by equivalence with a stable scheme, in this case the Lagrange–Galerkin scheme which has the form:

$$\int_{\mathbb{R}} v_{\Delta}^{n+1}(\xi) \phi_i(\xi) d\xi = \int_{\mathbb{R}} v_{\Delta}^n(\xi - c\Delta t) \phi_i(\xi) d\xi$$

that is, writing the numerical solution as $v_{\Delta}^k(x) = \sum_j v_j^k \phi_j(x)$,

$$\sum_{j} v_j^{n+1} \int_{\mathbb{R}} \phi_j(\xi) \phi_i(\xi) d\xi = \sum_{j} v_j^n \int_{\mathbb{R}} \phi_j(x_i - c\Delta t) \phi_i(\xi) d\xi$$

• In the LG scheme, interpolation is replaced by Galerkin projection

We will rather follow the line of proving stability by equivalence with a stable scheme, in this case the Lagrange–Galerkin scheme which has the form:

$$\int_{\mathbb{R}} v_{\Delta}^{n+1}(\xi) \phi_i(\xi) d\xi = \int_{\mathbb{R}} v_{\Delta}^n(\xi - c\Delta t) \phi_i(\xi) d\xi$$

that is, writing the numerical solution as $v_{\Delta}^k(x) = \sum_j v_j^k \phi_j(x)$,

$$\sum_{j} v_j^{n+1} \int_{\mathbb{R}} \phi_j(\xi) \phi_i(\xi) d\xi = \sum_{j} v_j^n \int_{\mathbb{R}} \phi_j(x_i - c\Delta t) \phi_i(\xi) d\xi$$

- In the LG scheme, interpolation is replaced by Galerkin projection
- As a consequence, $\|v_{\Delta}^{n+1}\|_2 \leq \|v_{\Delta}^n\|_2$ (i.e., the scheme is stable)

The Galerkin basis is supposed to have a structure similar to the SL basis:

$$\phi_j(\xi) = \frac{1}{\sqrt{\Delta x}} \phi\left(\frac{\xi}{\Delta x} - j\right)$$

where ϕ is the reference LG basis function, and the factor $\frac{1}{\sqrt{\Delta x}}$ gives the correct scaling in the integration

The Galerkin basis is supposed to have a structure similar to the SL basis:

$$\phi_j(\xi) = \frac{1}{\sqrt{\Delta x}} \phi\left(\frac{\xi}{\Delta x} - j\right)$$

where ϕ is the reference LG basis function, and the factor $\frac{1}{\sqrt{\Delta x}}$ gives the correct scaling in the integration

• The condition of equivalence between SL and LG schemes relates the reference functions ϕ and ψ with integral equation:

$$\int_{\mathbb{R}} \phi(\eta + t) \phi(\eta) d\eta = \psi(t)$$

that is, ϕ must have ψ as its autocorrelation

This problem has a solution (in general, nonunique) if and only if:

 \bullet The function ψ is positive definite, that is

$$\sum_{k=1}^n \sum_{j=1}^n a_k \psi(t_k - t_j) \bar{a}_j \ge 0$$

for any $t_k \in \mathbb{R}$, $a_k \in \mathbb{C}$ $(k = 1, \dots, n)$ and for all $n \in \mathbb{N}$

 \bullet Equivalently, the function ψ has a real positive Fourier transform $\hat{\psi}$

This problem has a solution (in general, nonunique) if and only if:

 \bullet The function ψ is positive definite, that is

$$\sum_{k=1}^n \sum_{j=1}^n a_k \psi(t_k - t_j) \bar{a}_j \ge 0$$

for any $t_k \in \mathbb{R}$, $a_k \in \mathbb{C}$ $(k = 1, \dots, n)$ and for all $n \in \mathbb{N}$

- \bullet Equivalently, the function ψ has a real positive Fourier transform $\hat{\psi}$
- The solution is given by $\phi(t) = \mathcal{F}^{-1} \left\{ \widehat{\psi}(\omega)^{1/2} \right\}$

This problem has a solution (in general, nonunique) if and only if:

 \bullet The function ψ is positive definite, that is

$$\sum_{k=1}^n \sum_{j=1}^n a_k \psi(t_k - t_j) \bar{a}_j \ge 0$$

for any $t_k \in \mathbb{R}$, $a_k \in \mathbb{C}$ (k = 1, ..., n) and for all $n \in \mathbb{N}$

- \bullet Equivalently, the function ψ has a real positive Fourier transform $\hat{\psi}$
- The solution is given by $\phi(t) = \mathcal{F}^{-1} \left\{ \widehat{\psi}(\omega)^{1/2} \right\}$
- Existence of a solution implies L^2 stability of SL schemes



The reference function ψ for \mathbb{P}_1 interpolation



The "obvious" LG counterpart ϕ for \mathbb{P}_1 interpolation



The minimal phase LG counterpart ϕ for \mathbb{P}_1 interpolation



SL and LG, cubic



SL and LG, quintic

Situations covered by this result:

• High–order Lagrange interpolations which can be shown to have a positive Fourier transform (tested for $n \leq 13$):

$$\widehat{\psi}^{(n)}(\omega) = p(\omega^2) \frac{\sin\left(\frac{\omega}{2}\right)^{n+1}}{\left(\frac{\omega}{2}\right)^{n+1}}$$

with $p(\omega^2)$ a polynomial of degree [n/2] with positive coefficients.

- Interpolatory wavelets, usually defined to be positive definite functions (e.g., in the case of the Shannon wavelet, $\hat{\psi}(\omega) = \mathbb{1}_{(-\pi,\pi)}(\omega)$).
- Cubic splines (no rigorous proof)

• In general, (as for the case of the \mathbb{P}_1 base) we expect to have multiple solutions to the problem: in fact, the relationship between $\hat{\phi}$ and $\hat{\psi}$,

$$|\widehat{\phi}(\omega)|^2 = \widehat{\psi}(\omega)$$

poses no constraint on the phase of $\widehat{\phi}$

• In general, (as for the case of the \mathbb{P}_1 base) we expect to have multiple solutions to the problem: in fact, the relationship between $\hat{\phi}$ and $\hat{\psi}$,

$$|\hat{\phi}(\omega)|^2 = \hat{\psi}(\omega)$$

poses no constraint on the phase of $\widehat{\phi}$

• The possibility to generate solutions with different phase terms is a tool to select a solution with prescribed decay and/or smoothness requirements (a key tool to treat the variable coefficient case)

index

Construction of Semi-Lagrangian schemes for convex HJ equations

Concerning HJ equations, we refer to the model problem:

$$\begin{cases} u_t(x,t) + H(Du(x,t)) = 0, & (x,t) \in \mathbb{R}^d \times [0,T] \\ u(x,0) = u_0(x) & x \in \mathbb{R}^d. \end{cases}$$

• Typical assumptions on H(p): smoothness, convexity, coercivity (e.g., a lower bound on H_{pp})

Construction of Semi-Lagrangian schemes for convex HJ equations

Concerning HJ equations, we refer to the model problem:

$$\begin{cases} u_t(x,t) + H(Du(x,t)) = 0, & (x,t) \in \mathbb{R}^d \times [0,T] \\ u(x,0) = u_0(x) & x \in \mathbb{R}^d. \end{cases}$$

- Typical assumptions on H(p): smoothness, convexity, coercivity (e.g., a lower bound on H_{pp})
- Various extensions (in particular, to Dynamic Programming Equations) are possible

The representation formula which parallels the formula of characteristics for HJ equations, is termed as the *Hopf–Lax formula*:

$$u(x,t+\tau) = \min_{a \in \mathbb{R}^d} [\tau H^*(a) + u(x-a\tau,t)]$$

where

$$H^*(a) = \sup_{p \in \mathbb{R}^d} [a \cdot p - H(p)]$$

is the Legendre transform of the Hamiltonian function H.

The representation formula which parallels the formula of characteristics for HJ equations, is termed as the *Hopf–Lax formula*:

$$u(x,t+\tau) = \min_{a \in \mathbb{R}^d} [\tau H^*(a) + u(x-a\tau,t)]$$

where

$$H^*(a) = \sup_{p \in \mathbb{R}^d} [a \cdot p - H(p)]$$

is the Legendre transform of the Hamiltonian function H.

Via the Hopf–Lax formula it can also be shown that the typical regularity achieved by the solution u is semiconcavity (roughly speaking, a unilateral upper bound on the second incremental ratio).

Semi–Lagrangian approximation for the convex HJ equation:

The Hopf–Lax representation formula is discretized as

$$v_i^{n+1} = \min_{\alpha \in \mathbb{R}^d} [\Delta t H^*(\alpha) + I[V^n](x_i - \alpha \Delta t)].$$

In addition to the reconstruction operator $I[V^n]$, two new ingredients are required:

Semi–Lagrangian approximation for the convex HJ equation:

The Hopf–Lax representation formula is discretized as

$$v_i^{n+1} = \min_{\alpha \in \mathbb{R}^d} [\Delta t H^*(\alpha) + I[V^n](x_i - \alpha \Delta t)].$$

In addition to the reconstruction operator $I[V^n]$, two new ingredients are required:

• A numerical approximation of the Legendre transform $H^*(\alpha)$ (whenever it cannot be explicitly computed)

Semi–Lagrangian approximation for the convex HJ equation:

The Hopf–Lax representation formula is discretized as

$$v_i^{n+1} = \min_{\alpha \in \mathbb{R}^d} [\Delta t H^*(\alpha) + I[V^n](x_i - \alpha \Delta t)].$$

In addition to the reconstruction operator $I[V^n]$, two new ingredients are required:

- A numerical approximation of the Legendre transform $H^*(\alpha)$ (whenever it cannot be explicitly computed)
- A derivative-free minimization procedure

index

Convergence analysis for the nonlinear problem

Beside consistency, two main concepts of stability are available for proving convergence in the nonlinear case:

Convergence analysis for the nonlinear problem

Beside consistency, two main concepts of stability are available for proving convergence in the nonlinear case:

• Barles–Souganidis theorem: the scheme should be invariant for the addition of constants, and *monotone up to a term* $o(\Delta t)$

Convergence analysis for the nonlinear problem

Beside consistency, two main concepts of stability are available for proving convergence in the nonlinear case:

- Barles–Souganidis theorem: the scheme should be invariant for the addition of constants, and *monotone up to a term* $o(\Delta t)$
- Lin–Tadmor theorem: the numerical solutions should be uniformly semiconcave

To prove **consistency**, at least in the sense of Barles–Souganidis, we compare again the scheme with the Hopf–Lax representation formula, assuming that u is a smooth solution and that $v_j^n = u(x_j, t_n)$. It results that the local truncation error has the estimate:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \frac{\Delta x^{r+1}}{\Delta t}$$

To prove **consistency**, at least in the sense of Barles–Souganidis, we compare again the scheme with the Hopf–Lax representation formula, assuming that u is a smooth solution and that $v_j^n = u(x_j, t_n)$. It results that the local truncation error has the estimate:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \frac{\Delta x^{r+1}}{\Delta t}$$

• A time discretization term $O(\Delta t^p)$ appears again as soon as characteristics are no longer straight lines To prove **consistency**, at least in the sense of Barles–Souganidis, we compare again the scheme with the Hopf–Lax representation formula, assuming that u is a smooth solution and that $v_j^n = u(x_j, t_n)$. It results that the local truncation error has the estimate:

$$L^{\Delta}(x_i, t_{n+1}) \Big| \le C \frac{\Delta x^{r+1}}{\Delta t}$$

- A time discretization term $O(\Delta t^p)$ appears again as soon as characteristics are no longer straight lines
- Consistency analysis is more technical in the Lin–Tadmor theory, although it comes to similar conclusions

Proving monotonicity up to an $o(\Delta t)$ is possible even for high-order reconstructions provided:

Proving **monotonicity up to an** $o(\Delta t)$ is possible even for high-order reconstructions provided:

• The numerical solutions are Lipschitz stable, so that the reconstruction satisfies monotonicity up to an $O(\Delta x)$ Proving **monotonicity up to an** $o(\Delta t)$ is possible even for high-order reconstructions provided:

• The numerical solutions are Lipschitz stable, so that the reconstruction satisfies monotonicity up to an $O(\Delta x)$

• The Courant number goes to infinity: $\Delta x = o(\Delta t)$ – here, the SL schemes have some more degrees of freedom in choosing the $\Delta t/\Delta x$ relationship

Lipschitz stability result: Consider the scheme in ${\mathbb R}$

$$v_i^{n+1} = \min_{\alpha \in \mathbb{R}} [\Delta t H^*(\alpha) + I_r[V^n](x_i - \alpha \Delta t)]$$

for a Hamiltonian function H(p) such that $H_{pp} \ge m_H$. Assume that, for some constant C < 1:

$$|I_r[V](x) - I_1[V](x)| \le C \max_{x_{j-1}, x_j, x_{j+1} \in \mathcal{S}(x)} |v_{j+1} - 2v_j + v_{j-1}|$$

(I_1 denoting the \mathbb{P}_1 interpolation, and $\mathcal{S}(x)$ denoting the reconstruction stencil at x) and that $\Delta x = O(\Delta t^2)$. Then, the family of numerical solutions V^n is Lipschitz stable.

Admissible reconstructions: the previous condition is satisfied for Lagrange reconstructions up to degree 5, provided the reconstruction stencil overlaps with the cell in which the reconstruction is performed.



admissible stencils
• Consequences: Lipschitz stability holds for ENO and finite element reconstructions up to degree 5

• Consequences: Lipschitz stability holds for ENO and finite element reconstructions up to degree 5

• Furthermore, the case of symmetric Lagrange or WENO reconstructions can be treated by proving that (linear) weights of WENO interpolation are nonnegative. This gives Lipschitz stability up to degree 5/9 for WENO and up to degree 9 for symmetric Lagrange. • Consequences: Lipschitz stability holds for ENO and finite element reconstructions up to degree 5

• Furthermore, the case of symmetric Lagrange or WENO reconstructions can be treated by proving that (linear) weights of WENO interpolation are nonnegative. This gives Lipschitz stability up to degree 5/9 for WENO and up to degree 9 for symmetric Lagrange.

• In the practical use of the SL scheme, the condition $\Delta x = O(\Delta t^2)$ seems overly restrictive.

index