

A splitting approach for freezing waves

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Abstract We present a numerical method which is able to approximate traveling waves (e.g. viscous profiles) in systems with hyperbolic and parabolic parts by a direct long-time forward simulation. A difficulty with long-time simulations of traveling waves is that the solution leaves any bounded computational domain in finite time. To handle this problem one should go into a suitable co-moving frame. Since the velocity of the wave is typically unknown, we use the method of freezing [2], see also [1], which transforms the original partial differential equation (PDE) into a partial differential algebraic equation (PDAE) and calculates a suitable co-moving frame on the fly. The efficient numerical approximation of this freezing PDAE is a challenging problem and we introduce a new numerical discretization, suitable for problems that consist of hyperbolic conservation laws which are (nonlinearly) coupled to parabolic equations. The idea of our scheme is to use the operator splitting approach. The benefit of splitting methods in this context lies in the possibility to solve hyperbolic and parabolic parts with different numerical algorithms. We test our method at the (viscous) Burgers' equation. Numerical experiments show linear and quadratic convergence rates for the approximation of the numerical steady state obtained by a long-time simulation for Lie- and Strang-splitting respectively. Due to these affirmative results we expect our scheme to be suitable also for freezing waves in hyperbolic-parabolic coupled PDEs.

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1 Introduction

Many partial differential equations from applications consist of different parts,

$$u_t = Au_{xx} + f(u)_x + g(u) =: F(u). \quad (1)$$

Sometimes, one part is parabolic while another part is hyperbolic and these parts are nonlinearly coupled. Examples of such hyperbolic-parabolic coupled PDEs are hyperbolic models of chemosensitive movement or reaction-diffusion equations for which not all components diffuse.

One is often interested in special solutions, which arise as (time-)asymptotic limits of solutions to the Cauchy problem for (1). An important class of such solutions are traveling waves. They describe how mass (or information) travels through the domain. From this interpretation, it is obvious, that one is often interested not only in the shape but also the velocity of the traveling wave.

Traveling waves are solutions of the PDE (1) of the form

$$u(x, t) = \bar{v}(x - \bar{\mu}t), \quad x \in \mathbb{R}, \quad t \in \mathbb{R},$$

where $\bar{v} : \mathbb{R} \rightarrow \mathbb{R}$ is the non-constant profile and $\bar{\mu} \in \mathbb{R}$ the velocity of the wave. For Burgers' equation there is a family of traveling wave solutions,

$$\begin{aligned} u(x, t) &= \varphi(x - \bar{\mu}t) + \frac{1}{2}(b + c) = \bar{v}(x - \bar{\mu}t), \\ \varphi(x) &= a \frac{1 - e^{ax}}{1 + e^{ax}}, \quad a = \frac{1}{2}(b - c), \quad \bar{\mu} = \frac{1}{2}(b + c), \end{aligned} \quad (2)$$

parametrized by the asymptotic states $\lim_{x \rightarrow -\infty} u(x, t) = b > c = \lim_{x \rightarrow \infty} u(x, t)$.

As a toy example we consider the Cauchy problem for the viscous Burgers' equation

$$u_t + \left(\frac{1}{2}u^2\right)_x = u_{xx} \quad \text{on } \mathbb{R} \times [0, \infty), \quad u(\cdot, 0) = u_0 \quad \text{on } \mathbb{R}. \quad (3)$$

For the numerical approximation of (3) one has to truncate the unbounded spatial domain to a finite interval. This leads to the problem that every traveling wave solution with non-zero speed eventually leaves the computational domain. The simplest remedy is to use periodic boundary conditions on a very large domain, but this is only reasonable in the case of pulses. Instead, we use the freezing method from [2]. The idea is to move the spatial frame with the speed of the traveling wave. We make the **ansatz** that the solution of (1) is of the form

$$u(x, t) = v(x - \gamma(t), t), \quad \gamma(t) \in \mathbb{R}, \quad (4)$$

where $\gamma(t)$ denotes a time dependent position. Then, $\mu(t) := \gamma'(t)$ can be interpreted as the velocity of the wave at time t . Plugging (4) into (1) yields

$$v_t = F(v) + \mu(t)v_x, \quad (5)$$

where both v and μ are unknown. Due to the additional unknown μ one has to complement (5) with an additional algebraic equation, called phase condition in [2], to retain well-posedness. In Burgers' case this transforms (3) into the PDAE

$$\begin{cases} v_t = v_{xx} - (\frac{1}{2}v^2)_x + \mu v_x, \\ 0 = \Psi(v, \mu), \\ \gamma_t = \mu(t), \end{cases} \quad \begin{aligned} v(\cdot, 0) &= u_0, \\ \gamma(0) &= 0. \end{aligned} \quad (6)$$

This is called the *freezing method* in [2]. We restrict to two standard choices for the phase condition, the *orthogonal phase condition* given by

$$0 = \Psi(v, \mu) := \langle v_t \mid v_x \rangle_{L^2} = \langle v_{xx} - (\frac{1}{2}v^2)_x + \mu v_x \mid v_x \rangle_{L^2} \quad (7)$$

and the *fixed phase condition* given by

$$0 = \Psi(v, \mu) := \langle v - \hat{v} \mid \hat{v}_x \rangle_{L^2} \quad (8)$$

with \hat{v} an appropriately chosen reference function.

For the numerical approximation of (6) we use splitting methods, which we briefly recall for convenience. Assume that a solution to an initial value problem of the form

$$u_t = A(u) + B(u) \quad (9)$$

is sought. Let $\Phi_A^t(u_0)$ and $\Phi_B^t(u_0)$ denote the solution operators for $u_t = A(u)$ and $u_t = B(u)$ with initial value u_0 , respectively. The *Lie-Trotter splitting*,

$$u^{n+1} = \Phi_B^{\Delta t} \circ \Phi_A^{\Delta t}(u^n), \quad (10)$$

typically converges linearly to the exact solution for $\Delta t \rightarrow 0$. A splitting method that typically leads to second order convergence is *Strang splitting*,

$$u^{n+1} = \Phi_A^{\Delta t/2} \circ \Phi_B^{\Delta t} \circ \Phi_A^{\Delta t/2}(u^n). \quad (11)$$

For example in [4], the authors show that this scheme is second order convergent for the viscous Burgers' PDE.

To apply this approach to the freezing PDAE (6), we split the equation into two parts to separate the hyperbolic and parabolic problem. Then we solve each part with a method which is particularly adapted to the respective sub-problem. Namely we solve the hyperbolic problem with an explicit scheme from Kurganov and Tadmor [5]. The parabolic sub-problem is solved by an implicit second order finite-difference approximation, due to the restrictive CFL condition.

Our main focus in this article is on approximating the limits of the time evolution and, unlike [4], not on the finite time convergence properties of the scheme. In particular, we aim to understand the preservation of steady states and their stability by our schemes.

In the case of ordinary differential equations there is a well-established theory for numerical steady states. For example in [8] there are results which state that one-step

methods preserve fixed points and their stability in a Δt^r -shrinking neighborhood under Lipschitz assumptions, where r denotes the order of consistency of the one-step method. An analogous result holds for the Strang splitting:

Theorem 1 ([3]). *Let $A, B \in C^3(\mathbb{R}^m, \mathbb{R}^m)$ and assume that \hat{u} is a hyperbolic fixed point of (9). Let φ_A, φ_B be one-step methods approximating Φ_A, Φ_B respectively. If φ_A, φ_b are second order Runge-Kutta methods then there exist $\Delta t_0, K > 0$, such that the numerical Strang splitting, $U^{n+1} = \varphi^{\Delta t}(U^n) = \varphi_B^{\Delta t/2} \circ \varphi_A^{\Delta t} \circ \varphi_B^{\Delta t/2}(U^n)$, has a fixed point \hat{U} which is unique in the ball $B(\hat{u}; K\Delta t^2)$ for all $0 < \Delta t \leq \Delta t_0$. Furthermore, \hat{U} is a stable (resp. unstable) fixed point of $\varphi^{\Delta t}$ if \hat{u} is a stable (resp. unstable) steady state of (9).*

For the freezing method there are several results available, where it is shown that the (continuous) method provides good approximations including preservation of asymptotic stability of traveling waves for certain problem classes in the continuous and semi-discrete case, see [6, 1, 9]. But the time-asymptotic behavior of a discretization with a splitting approach was never considered before.

A different approach to apply adapted schemes for different parts of the freezing PDAE appears in [7], where the freezing method is used to capture similarity solutions of the multidimensional Burgers' equation. There an IMEX-Runge-Kutta approach is used and second order convergence for the time dependent problem is shown on finite time intervals.

2 The splitting scheme

We now explicitly state our numerical scheme. We split (6) into two sub-problems as follows: Let $\Phi_A^t : (z_0, \gamma_0, \mu_0) \mapsto (z(t), \gamma(t), \mu(t))$ be the solution operator to the parabolic problem

$$\begin{cases} z_t = z_{xx}, & z(\cdot, 0) = z_0, \\ \gamma_t = 0, & \gamma(0) = \gamma_0, \\ \mu_t = 0, & \mu(0) = \mu_0, \end{cases} \quad (\text{A})$$

let $\Phi_B^t : (w_0, \gamma_0, \mu_0) \mapsto (w(t), \gamma(t), \mu(t))$ be the solution operator to

$$\begin{cases} w_t = -(\frac{1}{2}w^2)_x + \mu w_x, & w(\cdot, 0) = w_0, \\ 0 = \Psi(w, \mu), & \gamma(0) = \gamma_0. \\ \gamma_t = \mu(t), \end{cases} \quad (\text{B})$$

Here Ψ is one of the phase conditions (7) or (8). Note that the initial value μ_0 is ignored for this operator (B), because it is uniquely determined by the constraints. Since the splitting approach now iterates both solution operators consecutively, the question when and how to solve the algebraic constraint arises. For the orthogonal phase condition we chose an explicit and for the fixed phase condition we use a half-explicit approach. Thus we calculate the speed μ prior to solving the nonlinear PDE,

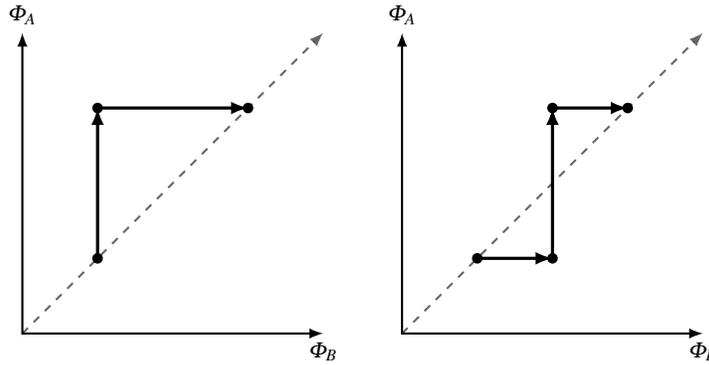


Fig. 1 Diagram of the Lie-Trotter splitting on the left and the Strang splitting on the right.

convergence	order 1		order 2	
full problem	$u_t + (\frac{1}{2}u^2)_x = u_{xx}$			
freezing method	orthogonal or fixed p.c.		fixed p.c.	
sub-problem	$w_t = -(\frac{1}{2}w^2)_x$	$z_t = z_{xx}$	$w_t = -(\frac{1}{2}w^2)_x$	$z_t = z_{xx}$
semi-discrete formulation	Rusanov Scheme	discrete Laplacian	Kurganov-Tadmor	discrete Laplacian
time discretization	forward Euler	backward Euler	Heun's method	Crank-Nicolson
splitting method	Lie		Strang	

Fig. 2 Overview of the applied numerical schemes for the presented schemes which offer a numerical steady state.

the μw_x part is then discretized by using finite differences. Lie and Strang splitting are illustrated by classical diagrams in **Fig. 1**. A step in the vertical direction in Fig. 1 amounts in numerically solving the Cauchy problem for the heat equation (A), whereas a step in the horizontal direction amounts to solve the hyperbolic PDAE (B). Only states on the dashed diagonal line might be considered as approximations to solutions to the original problem. In addition, the order of the sub-problems (A), (B) in the splitting approach is chosen such that the phase condition is satisfied at the end of a full time step. More details about how to calculate the speed with the algebraic constraint can be found in the description of the schemes, see (12), (13), (16) and (17). A schematic overview of the schemes is given in Fig. 2.

2.1 First order scheme.

We first present a first order scheme. For this we use a method of lines (MOL) approach for (B): We choose a finite interval $[L_-, L_+]$ and choose a spatial grid with uniform step size Δx and spatially discretize with the semi-discrete version of the Rusanov scheme. We use Dirichlet boundary conditions. It is worth mentioning here that for example the LxF or NT scheme do not offer a semi-discrete version. The **Rusanov scheme (RS)** in its semi-discrete form is given by

$$\begin{aligned} \frac{d}{dt} w_j(t) &= -\frac{f(w_{j+1}(t)) - f(w_{j-1}(t))}{2\Delta x} \\ &\quad + \frac{\kappa}{2\Delta x} [w_{j+1}(t) - 2w_j(t) + w_{j-1}(t)] \\ &= -\partial_0 f(w(t))_j + \kappa \frac{\Delta x}{2} \partial_0^2 w(t)_j \\ &=: \text{RS}^{\Delta x}(w(t)), \end{aligned}$$

where ∂_0 is the central difference quotient, $\partial_0 w_j = \frac{1}{2\Delta x}(w_{j+1} - w_{j-1})$, ∂_0^2 the discrete Laplacian, both with Dirichlet boundary conditions and $\kappa = \max_j u(j\Delta x, 0)$ is the maximum over the initial value evaluated at all grid points.

This scheme is in a simplified form: Since the local maximal speeds, used in the Kurganov-Tadmor scheme, ensure that all information of the Riemann fans stay in each cell of the discretized problem, they can be replaced by an upper bound. In the case of the Burgers' nonlinearity this upper bound is given by the maximal absolute value of the solution, which, in turn, is given by the maximal absolute value κ of the initial function u_0 due to the maximums principle.

The time discretization is done with a uniform step size Δt , for the first order version we use the forward Euler method. The numerical approximation of $\Phi_B^{\Delta t}$ will be denoted by $\phi_{B, \text{RS}\Delta x}^{\Delta t}$ and $\varphi_{B, \text{RS}\Delta x}^{\Delta t}$ for the two different phase conditions (7) and (8), respectively. The operator $\phi_{B, \text{RS}\Delta x}^{\Delta t}$ is given as the function which maps w_0, γ_0, μ_0 to the solution w^1, γ^1, μ^1 of the system

$$\begin{cases} w^1 = w^0 + \Delta t \text{RS}^{\Delta x}(w^0) + \Delta t \mu^* \partial_0 w^0, \\ \mu^* = -\frac{\partial_0 w^{0\top} (\partial_0^2 w^0 - w^0 \partial_0 w^0)}{\partial_0 w^{0\top} \partial_0 w^0}, \\ \gamma^1 = \gamma_0 + \Delta t \mu^1, \\ \mu^1 = -\frac{\partial_0 w^{1\top} (\partial_0^2 w^1 - w^1 \partial_0 w^1)}{\partial_0 w^{1\top} \partial_0 w^1}, \end{cases} \quad \begin{cases} w^0 = w_0, \\ \gamma^0 = \gamma_0, \end{cases} \quad (12)$$

where we use a discrete version of the orthogonal phase condition (7). For the fixed phase condition (8) the operator $\varphi_{B, \text{RS}\Delta x}^{\Delta t}$ is given as the mapping, which maps w_0, γ_0, μ_0 to the solution w^1, γ^1, μ^1 of the system

$$\begin{cases} w^1 = w^0 + \Delta t \text{RS}^{\Delta x}(w^0) + \Delta t \mu^1 \partial_0 w^0, \\ \mu^1 = -\frac{\partial_0 \hat{v}^\top (w^0 + \Delta t \text{RS}^{\Delta x}(w^0) - \hat{v})}{\Delta t \partial_0 \hat{v}^\top \partial_0 w}, \\ \gamma^1 = \gamma_0 + \Delta t \mu^1, \end{cases} \quad \begin{cases} w^0 = w_0, \\ \gamma^0 = \gamma_0. \end{cases} \quad (13)$$

Also for the sub-problem (A) we use a MOL approach, namely we spatially discretize (A) by finite differences, i.e. the discrete Laplacian with Dirichlet boundary conditions, ∂_0^2 , is used to approximate the second spatial derivative,

$$\frac{d}{dt} z_j = \partial_0^2 z_j, \quad z_j(0) = z_j^0.$$

For the time discretization we use backward Euler, because implicit methods have better stability properties for this type of equation. Using the linearity of ∂_0^2 , this leads to $\phi_{A, \text{BE} \Delta x}^{\Delta t} : (z_0, \gamma_0, \mu_0) \mapsto (z^1, \gamma^1, \mu^1)$ where

$$\begin{cases} z^1 = (I - \Delta t \partial_0^2)^{-1} z^0, & z^0 = z_0, \\ \gamma^1 = \gamma^0, & \gamma^0 = \gamma_0, \\ \mu^1 = \mu^0, & \mu^0 = \mu_0, \end{cases} \quad (14)$$

such that $\phi_{A, \text{BE} \Delta x}^{\Delta t} \approx \Phi_A^{\Delta t}$.

By using the Lie splitting (10), the full scheme for the freezing PDAE (6) is given by

$$\begin{pmatrix} v^{n+1} \\ \gamma^{n+1} \\ \mu^{n+1} \end{pmatrix} := \phi_{B, \text{RS} \Delta x}^{\Delta t} \circ \phi_{A, \text{BE} \Delta x}^{\Delta t} \begin{pmatrix} v^n \\ \gamma^n \\ \mu^n \end{pmatrix} \quad (\text{LO})$$

for the orthogonal phase condition and by

$$\begin{pmatrix} v^{n+1} \\ \gamma^{n+1} \\ \mu^{n+1} \end{pmatrix} := \phi_{B, \text{RS} \Delta x}^{\Delta t} \circ \phi_{A, \text{BE} \Delta x}^{\Delta t} \begin{pmatrix} v^n \\ \gamma^n \\ \mu^n \end{pmatrix} \quad (\text{LF})$$

for the fixed phase condition.

2.2 Second order scheme.

To construct a scheme with quadratic convergence in time and space we have to replace our numerical solution operators by suitable second order schemes and use Strang splitting instead of Lie splitting. For the nonlinear hyperbolic part we use the second order semi-discrete scheme from [5]. It is given as

$$\begin{aligned}
\frac{d}{dt}u_j(t) &= -\frac{1}{2\Delta x} \left(f(u_{j+1/2}^+(t)) + f(u_{j+1/2}^-(t)) - f(u_{j-1/2}^+(t)) - f(u_{j-1/2}^-(t)) \right) \\
&\quad + \frac{\kappa}{2\Delta x} \left(u_{j+1/2}^+(t) - u_{j+1/2}^-(t) - u_{j-1/2}^+(t) + u_{j-1/2}^-(t) \right) \\
&=: \text{KT}^{\Delta x}(u(t)),
\end{aligned} \tag{15}$$

where

$$u_{j+\frac{1}{2}}^{\pm}(t) := u_{j+\frac{1}{2}\pm\frac{1}{2}}(t) \mp \frac{\Delta x}{2}(u_x)_{j+\frac{1}{2}\pm\frac{1}{2}}(t)$$

for $j = -M, \dots, M$ with $u(t) \in \mathbb{R}^{2M+1}$ and $u_j(t) \in \mathbb{R}$ its j -th element. The slopes are approximated using the minmod limiter

$$(u_x)_j^n = \text{minmod} \left(\frac{u_j^n - u_{j-1}^n}{\Delta x}, \frac{u_{j+1}^n - u_j^n}{\Delta x} \right),$$

where $\text{minmod}(a, b) := \frac{1}{2}[\text{sgn}(a) + \text{sgn}(b)] \cdot \min(|a|, |b|)$. For the time integration we use Heun's method. In the case of (7), $\phi_{B, \text{KT}^{\Delta x}}^{\Delta t}$ is the mapping $\phi_{B, \text{KT}^{\Delta x}}^{\Delta t} : (w_0, \gamma_0, \mu_0) \mapsto (w^1, \gamma^1, \mu^1)$ given by the solution of

$$\begin{cases}
w^* = w^0 + \Delta t \text{KT}^{\Delta x}(w^0) + \Delta t \mu^* \partial_0 w^0, \\
w^1 = \frac{1}{2}w^0 + \frac{1}{2}(w^* + \Delta t \text{KT}^{\Delta x}(w^*) + \Delta t \mu^* \partial_0 w^*), \\
\mu^* = -\frac{\partial_0 w^{0\top} (\partial_0^2 w^0 - w^0 \partial_0 w^0)}{\partial_0 w^{0\top} \partial_0 w^0}, \\
\gamma^1 = \gamma_0 + \Delta t \mu^1, \\
\mu^1 = -\frac{\partial_0 w^{1\top} (\partial_0^2 w^1 - w^1 \partial_0 w^1)}{\partial_0 w^{1\top} \partial_0 w^1},
\end{cases} \quad \begin{aligned} w^0 &= w_0, \\ \gamma^0 &= \gamma_0. \end{aligned} \tag{16}$$

For the fixed phase condition (8) we define $\phi_{B, \text{KT}^{\Delta x}}^{\Delta t}$ as the mapping $(w_0, \gamma_0, \mu_0) \mapsto (w^1, \gamma^1, \mu^1)$

$$\begin{cases}
w^* = w^0 + \Delta t \text{KT}^{\Delta x}(w^0) + \Delta t \mu^1 \partial_0 w^0, \\
w^1 = \frac{1}{2}w^0 + \frac{1}{2}(w^* + \Delta t \text{KT}^{\Delta x}(w^*) + \Delta t \mu^1 \partial_0 w^*), \\
\mu^1 = -\frac{\partial_0 \hat{v}^\top (w^0 + \Delta t \text{KT}^{\Delta x}(w^0) - \hat{v})}{\partial_0 \hat{v}^\top \partial_0 w}, \\
\gamma^1 = \gamma_0 + \Delta t \mu^1,
\end{cases} \quad \begin{aligned} w^0 &= w_0, \\ \gamma^0 &= \gamma_0. \end{aligned} \tag{17}$$

For the heat equation, we use the Crank-Nicolson¹ method to discretize in time and, as in the first order version, the discrete Laplacian with Dirichlet boundary conditions, ∂_0^2 , is used in space. The solution operator $\phi_{A, \text{CN}^{\Delta x}}^{\Delta t}$ is given by the mapping $(z_0, \gamma_0, \mu_0) \mapsto (z^1, \gamma^1, \mu^1)$ of

¹ The Crank-Nicolson method used here is only the discretization in time by combining the forward and backward Euler method.

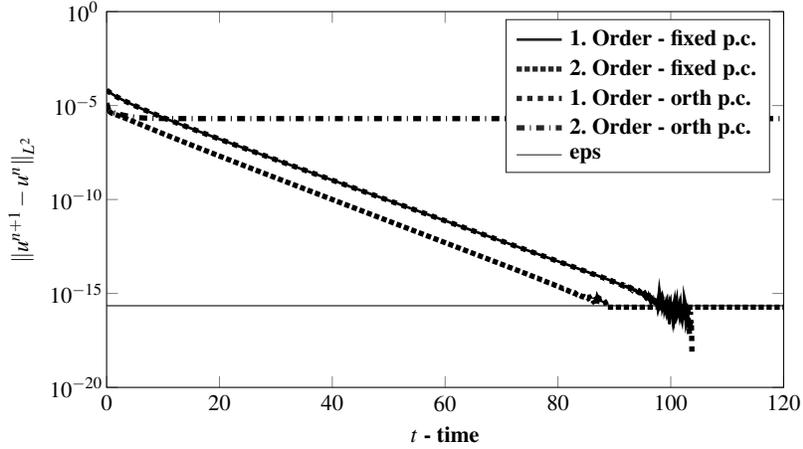


Fig. 3 Convergence to a numerical steady state except for the second order scheme with orthogonal phase condition.

$$\begin{cases} z^1 = (I - \frac{\Delta t}{2} \partial_0^2)^{-1} (I + \frac{\Delta t}{2} \partial_0^2) z^0, & z^0 = z_0, \\ \gamma^1 = \gamma^0, & \gamma^0 = \gamma_0, \\ \mu^1 = \mu^0, & \mu^0 = \mu_0. \end{cases}$$

These methods were chosen, because they offer quadratic convergence for the individual problems and thus we can hope for quadratic convergence of the full problem with Strang splitting. Strang splitting (11) leads to our second order scheme given by

$$\begin{pmatrix} v^{n+1} \\ \gamma^{n+1} \\ \mu^{n+1} \end{pmatrix} = \phi_{B,KT\Delta x}^{\Delta t/2} \circ \phi_{A,CN\Delta x}^{\Delta t} \circ \phi_{B,KT\Delta x}^{\Delta t/2} \begin{pmatrix} v^n \\ \gamma^n \\ \mu^n \end{pmatrix} \quad (\text{SO})$$

for the orthogonal phase condition and by

$$\begin{pmatrix} v^{n+1} \\ \gamma^{n+1} \\ \mu^{n+1} \end{pmatrix} = \phi_{B,KT\Delta x}^{\Delta t/2} \circ \phi_{A,CN\Delta x}^{\Delta t} \circ \phi_{B,KT\Delta x}^{\Delta t/2} \begin{pmatrix} v^n \\ \gamma^n \\ \mu^n \end{pmatrix} \quad (\text{SF})$$

for the fixed phase condition.

3 Numerical Results

The purpose of our schemes is to calculate viscous profiles by a simple forward simulation and thus we are interested in the quality of those profiles obtained at the end of a long-time simulation. Note that we do not consider the initial convergence

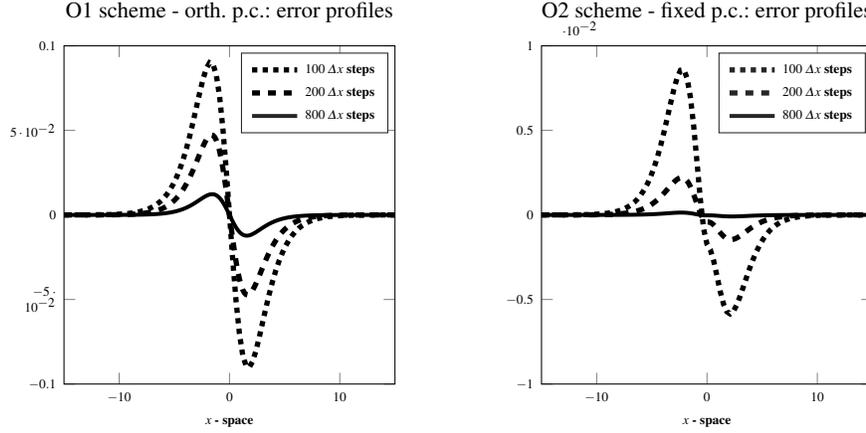


Fig. 4 Different numerical steady states for different $dt = \frac{\Delta x}{10}$. Note that the errors dominate where the profile varies the most and not at the boundary.

order on finite intervals. For all following computations we use the finite interval $[-15, 15]$ and Dirichlet boundary conditions. The reference function is given by (2) using $b = 1.5$, $c = -0.5$, which is also used as initial value with $t = 0$. This leads to a speed of 0.5 for the traveling wave. Since we are looking for numerical steady states in the co-moving frame, we have to check if our numerical schemes yield steady states. A steady state has the property $\frac{d}{dt}u(t) = 0$, which translates in the numerical case to $u^{n+1} = u^n$. In Fig. 3 we plot the time against the discrete L^2 distance $\|u^{n+1} - u^n\|_{L^2}$ and see that our schemes yield steady states at around $t \approx 100$ for (LO), (LF) and (SF) since $\|u^{n+1} - u^n\|_{L^2} \approx$ machine precision. For the Strang splitting scheme with orthogonal phase condition (SO) we see that $\|u^{n+1} - u^n\|_{L^2}$ does not converge to zero and the scheme does not offer a steady state. Solutions for this scheme leave the co-moving frame because the approximation of the speed is incorrect in this case. For these computations we use 300 steps in space and $\Delta t = \frac{\Delta x}{10}$.

Next, we consider the error profiles of the calculated steady states with different step sizes. This result is shown in Fig. 4. Obviously, we get different numerical steady states for different $dt = \frac{\Delta x}{10}$, which approximates the exact steady state better for smaller steps sizes. In addition, we observe that the dominant error occurs in the profile and there is more-or-less no error at the boundary.

The most interesting observation in our case is the convergence of our numerical steady states to the exact one. For this we plot the discrete L^2 error of our states to the exact one for different step sizes in Fig. 5. Here we see linear convergence of our numerical steady states to the exact one for our first order scheme. For the second order version we see quadratic convergence.

Finally, we note that usually the exact solution of the traveling wave is unknown. Therefore one has to guess some suitable reference function. In Fig. 6 we see that a

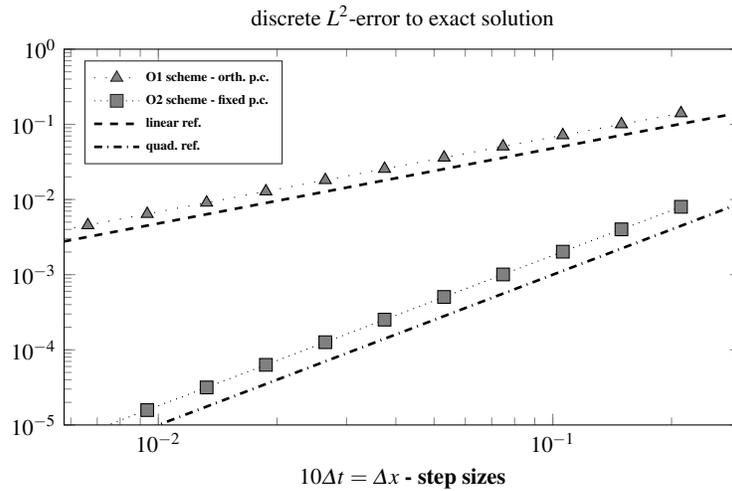


Fig. 5 Convergence rates of the numerical steady states to the exact one. The scheme (LF) was omitted because it produces the same results as (LO), whereas the scheme (SO) was ignored because it does not offer steady states.

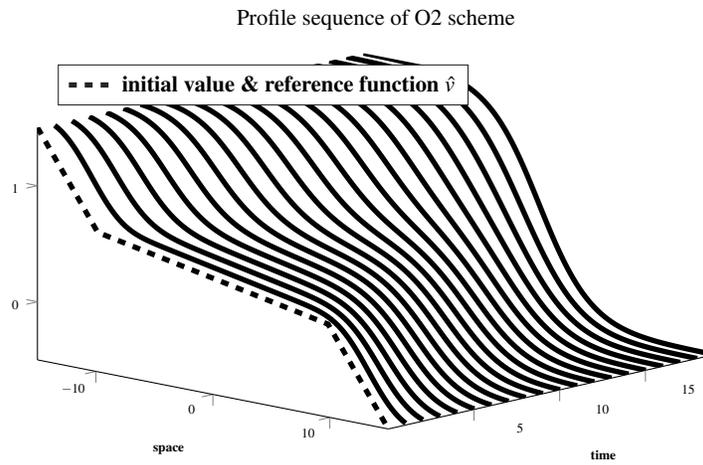


Fig. 6 Solution using initial value and reference function which only covers the rough behavior of the solution.

rough guess is sufficient for the initial value as well as for the reference function \hat{v} . The forward simulation approximates the traveling wave as before.

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