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Numerical solution of nonlinear advection equations by the inflow-implicit/outflow-explicit finite volume method

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Abstract

In the first part in this work we present a scheme for solving nonlinear advection-diffusion equations. We treat the advection term by the inflow-implicit/outflow-explicit (IIOE) and we apply the Crank-Nicolson scheme for the diffusion term. We tested the performance of the scheme on the viscous Burgers' equation and the traffic flow equation by comparing numerical solutions with exact ones. In numerical experiments we can observe second order convergence.

In the second part of our work we deal with first order 1D scalar conservation laws. We show, that the IIOE scheme can be written in conservative form. It is shown by Taylor series analysis that the scheme is second order accurate for smooth solutions on a uniform grid in the 1D case. We test the scheme numerically for smooth solutions of conservation laws. For problems, where solutions can develop discontinuities, we present a possible flux limiting procedure, which yields the flux limited IIOE (FLIIOE) method. Numerical experiments are also presented for the flux limited method.

Keywords: viscous Burgers' equation, traffic flow, nonlinear conservation laws, finite volume method, semi-implicit scheme, flux limiter

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Chapter 1

Introduction

Our work deals with problems of computational fluid dynamics, which share similarities with the general equations describing flow phenomena. For example, the 1D Burgers' equation can be obtained through simplifications of the incompressible Navier-Stokes equations. Our intention is not to present detailed derivation of these equations. We only would like to discuss some concepts of fluid mechanics relevant to our work. For a more exhaustive explanation, see e.g. [3, 6, 9, 10, 12, 32].

Fluid motion is one of the most fascinating phenomenon in nature. It is of great theoretical and practical interest to better understand it. Mathematics has been proved to be very helpful throughout the history to tackle such problems. A long path and hard work of great minds lead to our current level of understanding, which is, however, far from complete. For historical details, see e.g. [6, 32] and references therein.

Although we know that fluids, liquids and gases, are made up of molecules, we can treat them as a continuum. This is a reasonable assumption assuming that the systems we want to analyze are large compared to the distance between molecules. This is the so called continuum hypothesis. This allows us to describe its properties as continuous scalar functions of space and time. There are two basic approaches to describe the motion of fluids: Lagrangian and Eulerian.

In the Lagrangian description we follow the particles as they move along trajectories. Every particle is somehow labeled, usually with its initial position. As one can imagine, it could be quiet cumbersome or even impossible to get a detailed information of the flow that way.

In the Eulerian viewpoint the properties of the fluid are represented by scalar func-

tions of a fixed position $\mathbf{x} = (x_1, x_2, x_3)$ in 3D space and time t. The scalar quantity usually means density, a velocity component, pressure, temperature, etc. Obviously we assume that at a given position there is only one particle at a given instant in time.

Our goal is to predict how these quantities change in the Eulerian description as time evolves. Suppose we are interested in the dynamics of fluid particles in an enclosed region in space, denoted as $\Omega(t)$. The region in general can be a function of time. The boundary of the region is denoted as $\partial\Omega(t)$. In particular we can assume that this region moves with the particles then it is usually referred to as a material volume [3]. A convenient way to describe changes of quantities in such region over time is to rely on the transport theorem, in fluid mechanics texts often named after Lord Osborne Reynolds, or after Gottfried Wilhelm von Leibniz. The transport theorem says

$$\frac{\mathrm{d}}{\mathrm{d}t} \int_{\Omega(t)} F(t, \mathbf{x}) \, \mathrm{d}\mathbf{x} = \int_{\Omega(t)} \frac{\partial}{\partial t} F(t, \mathbf{x}) \, \mathrm{d}\mathbf{x} + \int_{\partial \Omega(t)} (F(t, \mathbf{x}) \, \mathbf{v}(t, \mathbf{x})) \cdot \mathbf{n}(t, \mathbf{x}) \, \mathrm{d}S, \tag{1.1}$$

where $F(t, \mathbf{x})$ is the quantity of our interest per unit volume, $d\mathbf{x}$ denotes an infinitesimally small material volume, $\mathbf{v}(t, \mathbf{x}) = (v_1(t, \mathbf{x}), v_2(t, \mathbf{x}), v_3(t, \mathbf{x}))$ denotes velocity, $\mathbf{n}(t, \mathbf{x}) = (n_1(t, \mathbf{x}), n_2(t, \mathbf{x}), n_3(t, \mathbf{x}))$ is an outward unit normal vector, dS is an infinitesimally small surface element of the boundary. Using Gauss's theorem we can equivalently write (1.1) as

$$\frac{\mathrm{d}}{\mathrm{d}t} \int_{\Omega(t)} F(t, \mathbf{x}) \, \mathrm{d}\mathbf{x} = \int_{\Omega(t)} \frac{\partial}{\partial t} F(t, \mathbf{x}) + \nabla \cdot (F(t, \mathbf{x}) \, \mathbf{v}(t, \mathbf{x})) \, \mathrm{d}\mathbf{x}.$$

One of the basic laws of classical physics is the conservation of mass. Without the presence of any sources or sinks, the mass of a system of particles doesn't change over time. To express this mathematically we choose the scalar quantity to be the density, denoted as $\rho(t, \mathbf{x})$, which is mass per unit volume. Thus the conservation of mass can be expressed as

$$\frac{\mathrm{d}}{\mathrm{d}t} \int_{\Omega(t)} \rho(t, \mathbf{x}) \, \mathrm{d}\mathbf{x} = \int_{\Omega(t)} \frac{\partial}{\partial t} \rho(t, \mathbf{x}) + \nabla \cdot (\rho(t, \mathbf{x}) \, \mathbf{v}(t, \mathbf{x})) \, \mathrm{d}\mathbf{x} = 0.$$

The region can be arbitrarily chosen, which means that the expression under the integral must vanish everywhere. We end up with the conservation of mass in differential form

$$\frac{\partial}{\partial t}\rho(t, \mathbf{x}) + \nabla \cdot (\rho(t, \mathbf{x}) \mathbf{v}(t, \mathbf{x})) = 0,$$

which is often referred to as the continuity equation, since assuming that the continuum

hypothesis holds (without any sinks there won't be any voids in the fluid). If we further assume that the fluid under consideration is incompressible, in other words the variation of the density is, for our purposes, negligible, the continuity equation reduces to

$$\nabla \cdot \mathbf{v}(t, \mathbf{x}) = 0. \tag{1.2}$$

In describing the motion of fluids in a non-relativistic classical sense, we apply Newton's laws of motion on the material region, which says that

change of linear momentum of the system/time = net force acting on the system.

First we deal with the left hand side. The linear momentum of a point mass is defined as mass times velocity. We can extend this concept for fluids by defining the linear momentum/unit mass as $\rho(t, \mathbf{x}) \mathbf{v}(t, \mathbf{x})$, which is a vector with three components. Then according to the transport theorem we can express the time rate of change of the linear momentum as

$$\frac{\mathrm{d}}{\mathrm{d}t} \int_{\Omega(t)} \rho(t, \mathbf{x}) \, v_i(t, \mathbf{x}) \, \mathrm{d}\mathbf{x} = \int_{\Omega(t)} \frac{\partial}{\partial t} (\rho(t, \mathbf{x}) \, v_i(t, \mathbf{x})) + \nabla \cdot (\rho(t, \mathbf{x}) \, v_i(t, \mathbf{x}) \, \mathbf{v}(t, \mathbf{x})) \, \mathrm{d}\mathbf{x}.$$
for $i = 1, 2, 3$.

Now we focus our attention to forces. We can divide the forces to body forces (inertial forces, gravity, electromagnetic interactions, etc.), which are acting to every part of the material volume and surface forces, which are acting on the boundary of the region. Let $\mathbf{f}(t, \mathbf{x}) = (f_1(t, \mathbf{x}), f_2(t, \mathbf{x}), f_3(t, \mathbf{x}))$ denote the body forces per unit mass. The net body force acting on the material volume can be expressed as

$$\int_{\Omega(t)} \rho(t, \mathbf{x}) \, \mathbf{f}(t, \mathbf{x}) \, \mathrm{d}\mathbf{x}.$$

We can further divide the surface forces to normal and tangential components. These forces can be described by a second order stress tensor, denoted by τ_{ij} . The diagonal components are the normal stresses, which are composed of pressure and viscous stresses. Viscous stresses also act in the tangential direction, which are called shear stresses, making up the off-diagonal elements of the stress tensor. We can write the contribution of pressure acting in the direction of inward unit vector $-\mathbf{n}(t, \mathbf{x})$ as

$$-\int_{\partial\Omega(t)} p(t, \mathbf{x}) \mathbf{n}(t, \mathbf{x}) dS. \tag{1.3}$$

The contribution of the viscous stresses can be written as

$$\int_{\partial\Omega(t)} \tau_{ij}(t, \mathbf{x}) \cdot \mathbf{n}(t, \mathbf{x}) \, \mathrm{d}S. \tag{1.4}$$

We limit our discussion to Newtonian fluids, in which case the viscous stresses are linearly proportional to the strain rate. In that case the viscous stress tensor can be written as

$$\tau_{ij}(t, \mathbf{x}) = 2\mu \,\varepsilon_{ij}(t, \mathbf{x}),$$

where μ is the dynamic viscosity, which we assume is a constant in our case, ε_{ij} is the strain rate tensor, which is defined as

$$\varepsilon_{ij}(t, \mathbf{x}) = \frac{1}{2} \left(\frac{\partial v_i}{\partial x_j} + \frac{\partial v_j}{\partial x_i} \right).$$

Then using Gauss's theorem to rewrite the surface integrals (1.3), (1.4) to volume integrals, considering that the material volume can be chosen arbitrarily, it can be shown that Newton's second law can be written in differential form considering an incompressible Newtonian fluid as

$$\rho \frac{\partial}{\partial t} v_i(t, \mathbf{x}) + \rho \nabla \cdot (v_i(t, \mathbf{x}) \mathbf{v}(t, \mathbf{x})) = \rho f_i(t, \mathbf{x}) - \frac{\partial}{\partial x_i} p(t, \mathbf{x}) + \mu \Delta v_i(t, \mathbf{x}), \qquad (1.5)$$

for i = 1, 2, 3,

where $\rho(t, \mathbf{x}) = \rho = constant$ in case of an incompressible fluid, Δ is the Laplace operator.

From now on we will omit the independent variables in the notation. We can write the second term on the left hand side of the momentum equation (1.5) as

$$\rho \nabla \cdot (v_{i} \mathbf{v}) = \rho \left(\frac{\partial}{\partial x_{1}} (v_{i} v_{1}) + \frac{\partial}{\partial x_{2}} (v_{i} v_{2}) + \frac{\partial}{\partial x_{3}} (v_{i} v_{3}) \right) \\
= \rho \left(v_{1} \frac{\partial}{\partial x_{1}} v_{i} + v_{i} \frac{\partial}{\partial x_{1}} v_{1} + v_{2} \frac{\partial}{\partial x_{2}} v_{i} + v_{i} \frac{\partial}{\partial x_{2}} v_{2} + v_{3} \frac{\partial}{\partial x_{3}} v_{i} + v_{i} \frac{\partial}{\partial x_{3}} v_{3} \right) \\
= \rho \left(v_{i} \left(\underbrace{\frac{\partial}{\partial x_{1}} v_{1} + \frac{\partial}{\partial x_{2}} v_{2} + \frac{\partial}{\partial x_{3}} v_{3}}_{\nabla \cdot \mathbf{v} = 0} \right) + \underbrace{v_{1} \frac{\partial}{\partial x_{1}} v_{i} + v_{2} \frac{\partial}{\partial x_{2}} v_{i} + v_{3} \frac{\partial}{\partial x_{3}} v_{i}}_{\mathbf{v} \cdot \nabla v_{i}} \right). \tag{1.6}$$

Using the continuity equation for incompressible fluid (1.2) and dividing by ρ we arrive at the famous Navier-Stokes equations for incompressible flow:

$$\frac{\partial}{\partial t} v_i + \mathbf{v} \cdot \nabla v_i = f_i - \frac{1}{\rho} \frac{\partial}{\partial x_i} p + \frac{\mu}{\rho} \Delta v_i, \quad \text{for } i = 1, 2, 3,$$
 (1.7)

named after Claude-Louis Marie Henri Navier, and Sir George Gabriel Stokes, who are credited for the derivation. However, according to [29] and references therein, it is interesting to mention that Navier arrived at the correct equations despite making wrong assumptions about the underlying physics. He had no conception of shear stresses and only wanted to modify the inviscid equations, known as the Euler equations, to count with the molecular forces in the fluid. The first was Adhémar Jean Claude Barré de Saint-Venant who derived the equations correctly considering viscous stresses. Stokes also derived the equations correctly but two years after Saint-Venant. Interestingly, Saint-Venant's name has never become associated with these equations.

The nonlinear system (1.7) can be solved analytically only in a few simple cases. However, it is of great practical and theoretical interest to solve it in the general case. The trouble is mainly caused by the nonlinear advective term.

1.1 Viscous Burgers' equation

Considering the 1D case, the unknowns are functions of the x_1 coordinate. We denote x_1 by x, v_1 by u and f_1 by f. The equations (1.7) reduce to

$$u_t + u u_x = f - \frac{1}{\rho} p_x + \frac{\mu}{\rho} u_{xx}, \tag{1.8}$$

where the subscripts denote partial derivatives.

If we further neglect the pressure variations, in the absence of body forces, we get from (1.8) the one dimensional viscous Burgers' equation

$$u_t + u u_x = \sigma u_{xx},\tag{1.9}$$

where $\sigma = \mu/\rho$ is the diffusion coefficient (or kinematic viscosity in fluid mechanics context). The equation is named after a Dutch physicist Johannes Martinus Burgers, [5]. It is interesting to note that the equation was first studied by a British (later American) applied mathematician Harry Bateman [25, 2]. The equation captures some key features of the equations of fluid dynamics: nonlinear advection, arising from the advective terms in the linear momentum equations (1.7), and viscosity effects [25, 18]. The equation can be solved analytically by transforming it to the linear heat equation, known as the Cole-Hopf transformation [25, 33]. This fact allows us to compare numerical solutions with the exact ones. By studying the behaviour of its

numerical solutions, it can help us to predict the performance of a particular numerical scheme on the more complicated equations of fluid dynamics.

1.1.1 Cole-Hopf transformation

The nonlinear viscous Burgers' equation (1.9) can be converted to the linear heat equation, for which there exist different methods to solve explicitly. Interestingly, the transformation first appeared as an exercise in a 19th century textbook [25, 11]. Later the mathematicians Eberhard Hopf [15] and Julian Cole [7] rediscovered the method and now it is named in their honor [25]. Below we would like to present the key steps of the method.

We start with the function

$$v(t,x) = e^{\alpha \varphi(t,x)}, \tag{1.10}$$

where α is a nonzero constant. Then

$$\varphi(t, x) = \frac{1}{\alpha} \ln v(t, x),$$

which is a real function if v(t,x) > 0. We assume that v(t,x) solves the linear heat equation

$$v_t = \sigma \, v_{xx}. \tag{1.11}$$

We calculate the partial derivatives

$$v_t = \alpha \varphi_t e^{\alpha \varphi}, \quad v_x = \alpha \varphi_x e^{\alpha \varphi}, \quad v_{xx} = \alpha (\varphi_{xx} e^{\alpha \varphi} + \alpha \varphi_x^2 e^{\alpha \varphi}).$$

Substituting to the heat equation (1.11) we get

$$\alpha \varphi_t e^{\alpha \varphi} = \sigma \alpha (\varphi_{xx} e^{\alpha \varphi} + \alpha \varphi_x^2 e^{\alpha \varphi})$$

Cancelling the common term $\alpha e^{\alpha \varphi}$ the result is

$$\varphi_t = \sigma \, \varphi_{xx} + \sigma \, \alpha \, \varphi_x^2. \tag{1.12}$$

We differentiate with respect to x to end up with

$$\varphi_{tx} = \sigma \varphi_{xxx} + 2 \sigma \alpha \varphi_x \varphi_{xx}.$$

Notice that φ_x solves (1.9) for $\alpha = -1/(2\sigma)$.

We conclude that if v(t,x) > 0 is a positive solution to the heat equation then

$$u(t,x) = \frac{\partial}{\partial x} \left[-2\sigma \ln v(t,x) \right] = -2\sigma \frac{v_x(t,x)}{v(t,x)}$$
(1.13)

solves the viscous Burgers' equation (1.9).

Suppose we want to solve (1.9) with given initial conditions

$$u(0,x) = u^0(x), \quad \forall x \in \mathbb{R}.$$

First we choose a function $\varphi(t,x)$ that satisfies $\varphi_x(t,x) = u(t,x)$, e.g.

$$\varphi(t,x) = \int_0^x u(t,y) \, dy$$
, so $\varphi(0,x) = \int_0^x u^0(y) \, dy$.

According to (1.10) the corresponding initial condition for the heat equation

$$v(0,x) = e^{-\varphi(0,x)/(2\sigma)} = \exp\left(-\frac{1}{2\sigma} \int_{0}^{x} u^{0}(y) dy\right).$$

The lower limit of integration doesn't have to be zero, it can be arbitrarily chosen if it's needed. For example, in case of a delta function initial condition centered at the origin, which is one of our chosen examples, it would be problematic having 0 as a lower limit of integration. See, e.g. [25], where in case of the triangular wave solution, the lower limit was chosen to be $-\infty$. Notice that, in general, changing the lower limit has an effect of multiplying v by a constant so it doesn't change the solution u in (1.13).

The solution to the heat equation for t > 0 can be obtained as a convolution of the initial condition with the fundamental solution, see e.g. [25],

$$v(t,x) = \frac{1}{2\sqrt{\pi \sigma t}} \int_{-\infty}^{\infty} e^{-(x-s)^2/(4\sigma t)} v(0,s) ds$$

$$= \frac{1}{2\sqrt{\pi \sigma t}} \int_{-\infty}^{\infty} \exp\left(-\frac{(x-s)^2}{4\sigma t} - \frac{1}{2\sigma} \int_{0}^{s} u^0(r) dr\right) ds.$$
(1.14)

The solution u(t, x) is then obtained by (1.13).

1.1.2 Examples

Below we present some exact solutions to the viscous Burgers' equation (1.9), which can be obtained by the Cole-Hopf transformation. The examples are taken from [25].

Rarefaction wave. First we look for a solution of the Burgers' equation with an

initial condition in the form of a step function

$$u^{0}(x) = \begin{cases} u_{l}, & x \leq 0, \\ u_{r}, & x > 0, \end{cases}$$

 $u_l < u_r$ corresponds to a rarefaction wave.

Performing the steps of the Cole-Hopf transformation described previously one obtains for t>0

$$u(t,x) = u_l + \frac{u_r - u_l}{1 + \exp\left(\frac{u_r - u_l}{2\sigma}(x - st)\right) \operatorname{erfc}\left(\frac{x - u_l t}{2\sqrt{\sigma t}}\right) / \operatorname{erfc}\left(\frac{u_r t - x}{2\sqrt{\sigma t}}\right)}, \quad (1.15)$$

with $s = (u_l + u_r)/2$, and $\operatorname{erfc}(x) = 1 - \operatorname{erf}(x)$ denoting the complementary error function, where

$$\operatorname{erf}(x) = \frac{2}{\sqrt{\pi}} \int_{0}^{x} e^{-s^{2}} \, \mathrm{d}s.$$

The solution can be seen in Figure 2.5 on page 26.

Triangular wave. Another interesting example is the triangular-wave solution, which is a solution to (1.9), when the initial condition is the delta function $\delta(x)$ centered at the origin.

For t > 0 we get

$$u(t,x) = 2\sqrt{\frac{\sigma}{\pi t}} \frac{\exp(-x^2/4\sigma t)}{\coth(1/4\sigma) - \operatorname{erf}(x/2\sqrt{\sigma t})},$$
(1.16)

where

$$\coth(x) = \frac{\cosh(x)}{\sinh(x)} = \frac{e^x + e^{-x}}{e^x - e^{-x}}$$

is the hyperbolic cotangent function. The solution can be seen in Figure 2.7 on page 28.

Trigonometric solution. Our last example is the trigonometric solution, which can be obtained transforming a separable solution to the linear heat equation (1.11)

$$v(t, x) = a + b e^{-\sigma \lambda t} \cos(\sqrt{\lambda} x),$$

where a > b must hold to ensure that it is a positive everywhere and $\lambda > 0$ is a positive constant. As discussed previously, any positive solution to the linear heat equation can be transformed to a solution of the viscous Burgers' equation (1.9) using

the relationship (1.13). After performing the necessary calculations we obtain

$$u(t,x) = \frac{2\sigma b\sqrt{\lambda}\sin\sqrt{\lambda}x}{a e^{\sigma \lambda t} + b\cos\sqrt{\lambda}x}.$$
 (1.17)

The solution can be seen in Figure 2.8 on page 28.

1.2 Traffic flow

Similar equation describes the change of the density of cars in continuum trafficflow models. Let us give a brief derivation of one of the simplest such model, known as the LWR model after Lighthill, Whitham [20], and Richards [28].

We start with the conservation of cars which states that on a given section of a one-lane highway the number of cars changes only due to cars leaving or coming at the ends of the interval.

In the derivation we assume that the number of cars is large enough on the interval of our interest so that it a reasonable assumption that the cars are smoothly distributed and define the density of cars $\rho(t,x)$ in units cars per car length. For simplicity we assume that every car has the same length. Then we have $\rho = 0$ and $\rho = 1$ for empty road and bumper to bumper traffic, respectively. Doing this way the number of cars on a given interval $[x_1, x_2]$ of the road in a given time t can be expressed as

$$\int_{x_1}^{x_2} \rho(t, x) \, \mathrm{d}x.$$

Let us denote the velocity of cars as v(t, x) in the Eulerian description of the traffic. We can express the cars that flow through a given point x over a time interval $[t_1, t_2]$ as

$$\int_{t}^{t_2} \rho(t, x) v(t, x) \, \mathrm{d}t.$$

Assuming that the velocity is positive in the direction of the x axis we can express the conservation of cars as

$$\int\limits_{\underbrace{x_1}}^{x_2} \rho(t_2,x) \,\mathrm{d}x - \int\limits_{x_1}^{x_2} \rho(t_1,x) \,\mathrm{d}x = \int\limits_{\underbrace{t_1}}^{t_2} \rho(t,x_1) v(t,x_1) \,\mathrm{d}t - \int\limits_{\underbrace{t_1}}^{t_2} \rho(t,x_2) v(t,x_2) \,\mathrm{d}t \,.$$
 change of the number of cars from t_1 to t_2 inflow at x_1 outflow at x_2

Assuming that the functions are smooth, using the Newton-Leibniz theorem we can rewrite the law as

$$\int_{x_1}^{x_2} \int_{t_1}^{t_2} \frac{\partial}{\partial t} \rho(t, x) dt dx = -\int_{t_1}^{t_2} \int_{x_1}^{x_2} \frac{\partial}{\partial x} \left(\rho(t, x_1) v(t, x_1) \right) dx dt.$$

If the intervals can be arbitrarily chosen, we can write the conservation of cars in differential form:

$$\rho_t + (\rho \, v)_x = 0,\tag{1.18}$$

knowing that ρ and v are functions of time and space, the subscripts denote partial derivatives.

It is reasonable to expect that a car is moving slower in a denser traffic. In other words, the speed of a car decreases as the density increases. This suggests to express the velocity as a function of the density $v = v(\rho)$. Then we can define the flux of cars as $f(\rho) = \rho v(\rho)$ and write the conservation law (1.18) as

$$\rho_t + f(\rho)_x = 0. \tag{1.19}$$

It is not a trivial task to define the relationship between the velocity and density. There are several to choose from [17, 30]. Greenshields [14] proposed a linear relationship between the density and velocity, where he relied on experiments performed in equilibrium traffic, in which case the density and velocity are constants. In the Greenshields model the velocity is defined as

$$v(\rho) = v_{max}(1 - \rho), \tag{1.20}$$

where the constant v_{max} denotes the maximal velocity.

This assumption leads to a concave quadratic flux function

$$f(\rho) = \rho \, v_{max} \, (1 - \rho)$$

A way to extend this model [33], is to assume that the flux also depends on the density gradient. If the traffic is getting denser more rapidly the flux of cars reduces more. Applying this yields a modified flux function

$$f(\rho, \rho_x) = \rho \, v_{max} (1 - \rho) - D\rho_x, \tag{1.21}$$

where D is a diffusion coefficient [33, 17, 30]. Substituting (1.21) to (1.19) and putting the diffusion term to the right-hand side we get

$$\rho_t + u(\rho)\rho_x = D\rho_{xx},\tag{1.22}$$

where we denoted

$$u(\rho) = v_{max}(1 - 2\rho). \tag{1.23}$$

Multiplying both sides by $u'(\rho)$ we obtain

$$u'(\rho)\rho_t + u(\rho)u'(\rho)\rho_x = Du'(\rho)\rho_{xx},$$

which can be rewritten as

$$u_t + uu_x = Du_{xx} - Du''(\rho)\rho_x^2.$$

Since $u''(\rho) = 0$ we end up with the viscous Burgers' equation (1.9), where σ corresponds to the diffusion coefficient D. By solving (1.9) and using the relationship (1.23) we easily obtain a relation for the density

$$\rho(u) = \frac{1}{2} \left(1 - \frac{u}{v_{max}} \right). \tag{1.24}$$

The relationship (1.24) is used to obtain exact solutions for (1.22), which is then compared to our numerical solution, where we solve (1.22) directly.

The ideas discussed above were further extended to so called higher order models. First such model is known as the Payne-Whitham(PW) model [17, 30, 33]. In addition to the conservation of cars they added a second equation that mimics the momentum equation (1.8) expressing the acceleration of cars in the Eulerian description. There exist other models presented, e.g. in [30, 1, 35].

Chapter 2

Inflow-implicit/outflow-explicit method

2.1 Introduction

There is only a relatively small number of cases, when we can find an analytical solution for a differential equation. Therefore the development of numerical methods for finding approximate solutions with the desired accuracy is very important. For solving partial differential equations, several numerical methods have been developed. Among the most popular numerical techniques are the finite difference method, the finite volume method and the finite element method. These methods reduce the problem of solving a differential equation to solving a system of equations. Below we give a brief description of these methods for the 1D case, which is relevant to this work.

The finite difference method [19] is the oldest and in some ways the simplest approach. The idea is to replace derivatives with differences. The domain of interest is divided into a given number of elements connected by nodes. The differences are calculated from the approximations at the nodes. The way, in which we calculate the differences come from accuracy and stability considerations. The method is well suited for regular rectangular meshes, but becomes impractical for complicated domains, which are often the subject of interest in applications.

The finite volume method uses a different approach [18, 26]. First we subdivide our spatial domain of interest into "finite volumes" or grid cells. In 1D these are intervals

on the real line. We denote the *i*-th grid cell by

$$p_i = (x_{i-1/2}, x_{i+1/2}).$$

The method is based on the conservation of a quantity u(t,x). Instead of looking for approximations of the values of u at the nodes, as we do in the finite difference method, we are interested in the approximations of the integral of u over each interval. We assume that the change of the integral over a time interval $[t^{n-1}, t^n]$ for n = 1, 2, ... is due to the fluxes f(t,x) at the endpoints of the cells:

$$\int_{n_i} u(t^n, x) dx - \int_{n_i} u(t^{n-1}, x) dx = \int_{t^{n-1}}^{t^n} f(t, x_{i-1/2}) dt - \int_{t^{n-1}}^{t^n} f(t, x_{i+1/2}) dt.$$
 (2.1)

Our goal is to calculate how the average of u changes over time, i.e. if we know the integral at time t^{n-1} , what is its value at a later time t^n . For this purpose we approximate the average value of u by a constant representative value denoted as \bar{u}_i^{n-1} inside the i-th cell at time t^{n-1} . The approximation of the integral becomes

$$\int_{n_i} u(t^{n-1}, x) \, \mathrm{d}x \approx \bar{u}_i^{n-1} \, h,$$

where $h = x_{i+1/2} - x_{i-1/2}$ is the length of a cell. We can also express

$$\int_{t^{n-1}}^{t^n} f(t, x_{i-1/2}) \, \mathrm{d}t \approx F_{i-1/2} \, \tau,$$

where $F_{i-1/2}$ is an approximation of the average flux at $x_{i-1/2}$ and over time interval $[t^{n-1}, t^n]$ with $\tau = t^n - t^{n-1}$. After substitution the conservation law takes the form

$$(\bar{u}_i^n - \bar{u}_i^{n-1})h = (F_{i-1/2} - F_{i+1/2})\tau$$

Dividing by the length of the cell h, rearranging terms we get an expression for the approximate average value at time t^n as

$$\bar{u}_i^n = \bar{u}_i^{n-1} - \frac{\tau}{h} (F_{i+1/2} - F_{i-1/2}). \tag{2.2}$$

If we assume dependence of the flux at a given interface on the representative values in the neighboring cells in space and time we get a fully discrete method. The scheme (2.2) has a discrete conservation property, which we can see by summing $\bar{u}_i^n h$, the approximation of the integral of a quantity u, over any set of cells

$$h\sum_{i=I}^{J} \bar{u}_{i}^{n} = h\sum_{i=I}^{J} \bar{u}_{i}^{n-1} - \tau \sum_{i=I}^{J} (F_{i+1/2} - F_{i-1/2}).$$

It can be seen, that the terms in the sum of fluxes on the right-hand side cancel out, except the fluxes at the boundaries I and J, i.e.

$$h\sum_{i=I}^{J} \bar{u}_{i}^{n} = h\sum_{i=I}^{J} \bar{u}_{i}^{n-1} - \tau \left(F_{J+1/2} - F_{I-1/2} \right).$$

The scheme is said to be in conservation form, because it mimics the property of the conservation law (2.1), i.e. the change of a quantity inside an interval is due to fluxes at the endpoints. This property makes the method well suited for solving conservation laws. Unlike the finite difference method, it can be used even for complicated domains.

A very successful method for solving differential equations is the finite element method [27, 25]. Similarly as in the case of the previous methods, we subdivide the domain of interest into smaller pieces, finite elements. It is based either on a minimization principle, if possible, or on the concept of a weak solution to a differential equation. It is well suited for complicated geometries similarly to the finite volume method.

In this work we apply the inflow-implicit/outflow-explicit (IIOE) scheme to advection terms of the nonlinear equations discussed in the previous chapter. The method is based on finite volume space discretization and a semi-implicit discretization in time. Inflows to the cells are treated implicitly and outflows are treated explicitly. We could explain this idea intuitively that we know what is flowing out from a cell at a given time but leave the method to solve a system of equations determined by the inflows to obtain the solution values at the new time step. In earlier works, [23, 24] it was shown that the IIOE scheme is formally second order accurate in space and time for 1D linear advection problems with variable velocity. In this work, we show that the IIOE method is formally second order consistent also in case of nonlinear scalar conservation laws for smooth solutions. Combining with the Crank-Nicolson scheme for the diffusion term, we get a new numerical scheme for the nonlinear advection-diffusion problem (1.9). We also present a limited IIOE scheme for inviscid conservation laws, which is important for problems, where solutions tend to develop shocks.

2.2 Advection with variable velocity

For reader convenience we present here the derivation of the IIOE scheme as in [21, 22, 23, 24] for the equation

$$u_t + vu_x = 0, (2.3)$$

where v = v(x). We rewrite (2.3) in the equivalent form

$$u_t + (vu)_x - uv_x = 0, (2.4)$$

and integrating over a grid cell p_i with cell center x_i , length h, left border $x_{i-\frac{1}{2}}$, right border $x_{i+\frac{1}{2}}$ yields

$$\int\limits_{p_i} u_t \, dx + \int\limits_{p_i} (vu)_x \, dx - \int\limits_{p_i} uv_x \, dx = 0$$

Let us denote $v_i = v(x_i)$, $v_{i-\frac{1}{2}} = v(x_{i-\frac{1}{2}})$, $v_{i+\frac{1}{2}} = v(x_{i+\frac{1}{2}})$. Let us denote by u_i^n a value of the solution inside the *i*-th finite volume cell at time step n computed by the numerical scheme. We use a constant representation of the solution inside a cell p_i denoted by \bar{u}_i and constant representative values at the cell interfaces denoted by $\bar{u}_{i-\frac{1}{2}}$, $\bar{u}_{i+\frac{1}{2}}$ respectively. Using the Newton-Leibniz formula we obtain

$$\int_{v_i} u_t \, dx + v_{i+\frac{1}{2}} \bar{u}_{i+\frac{1}{2}} - v_{i-\frac{1}{2}} \bar{u}_{i-\frac{1}{2}} - \bar{u}_i (v_{i+\frac{1}{2}} - v_{i-\frac{1}{2}}) = 0.$$

By rearranging terms we get

$$\int_{p_i} u_t \, dx + v_{i-\frac{1}{2}} (\bar{u}_i - \bar{u}_{i-\frac{1}{2}}) + (-v_{i+\frac{1}{2}}) (\bar{u}_i - \bar{u}_{i+\frac{1}{2}}) = 0.$$

 $v_{i-\frac{1}{2}} > 0$ represents inflow from the left cell interface, while $(-v_{i+\frac{1}{2}}) > 0$ represents inflow from the right cell interface. Otherwise they represent outflows. Thus we define

$$a_{i-\frac{1}{2}}^{in} = \max(v_{i-\frac{1}{2}}, 0), \quad a_{i-\frac{1}{2}}^{out} = \min(v_{i-\frac{1}{2}}, 0),$$
 (2.5)

$$a_{i+\frac{1}{2}}^{in} = \max(-v_{i+\frac{1}{2}}, 0), \ a_{i+\frac{1}{2}}^{out} = \min(-v_{i+\frac{1}{2}}, 0).$$
 (2.6)

We use a simple finite difference approximation for the time derivative $\frac{u_i^n - u_i^{n-1}}{\tau}$, where τ is a uniform time step, take inflow implicitly, outflow explicitly and use the straightforward reconstructions $\bar{u}_i^n = u_i^n$, $\bar{u}_{i-\frac{1}{2}}^n = \frac{1}{2}(u_i^n + u_{i-1}^n)$, $\bar{u}_{i+\frac{1}{2}}^n = \frac{1}{2}(u_i^n + u_{i+1}^n)$, we obtain

the basic one-dimensional IIOE scheme for variable velocity:

$$u_{i}^{n} + \frac{\tau}{2h} a_{i-\frac{1}{2}}^{in} \left(u_{i}^{n} - u_{i-1}^{n} \right) + \frac{\tau}{2h} a_{i+\frac{1}{2}}^{in} \left(u_{i}^{n} - u_{i+1}^{n} \right) =$$

$$u_{i}^{n-1} - \frac{\tau}{2h} \left(a_{i-\frac{1}{2}}^{out} \left(u_{i}^{n-1} - u_{i-1}^{n-1} \right) + a_{i+\frac{1}{2}}^{out} \left(u_{i}^{n-1} - u_{i+1}^{n-1} \right) \right).$$

$$(2.7)$$

2.3 Nonlinear advection with diffusion

We can extend these ideas for velocities depending on the solution. For the advective part of the viscous Burgers' equation (1.9) we calculate the velocities from the solution at the cell interfaces. This leads us to a nonlinear system of equations, which we solve iteratively. The time dependent velocities in the k-th iteration become

$$v_{i-\frac{1}{2}}^{n,k}=(u_i^{n,k-1}+u_{i-1}^{n,k-1})/2,\quad v_{i+\frac{1}{2}}^{n,k}=(u_i^{n,k-1}+u_{i+1}^{n,k-1})/2,\quad k=1,2,3,....$$
 and $u_i^{n,0}=u_i^{n-1}.$

When solving the traffic flow problem (1.22), we replace u_i^n by ρ_i^n . Considering that characteristic speed in the equations of traffic flow (1.22) equals $1 - 2\rho$, using the same reconstructions as for u_i^n and cancelling common factors, for $v_{max} = 1$ the time dependent velocities according to (1.23) in the k-th iteration become

$$v_{i-\frac{1}{2}}^{n,k} = 1 - (\rho_i^{n,k-1} + \rho_{i-1}^{n,k-1}), \quad v_{i+\frac{1}{2}}^{n,k} = 1 - (\rho_i^{n,k-1} + \rho_{i+1}^{n,k-1}).$$

In both cases we consider the time dependent splitting to inflows and outflows

$$a_{i-\frac{1}{2}}^{in,n-1} = \max(\operatorname{sgn}(v_{i-\frac{1}{2}}^{n-1}),0), \quad a_{i-\frac{1}{2}}^{out,n-1} = \max(\operatorname{sgn}(-v_{i-\frac{1}{2}}^{n-1}),0), \tag{2.8}$$

$$a_{i+\frac{1}{2}}^{in,n-1} = \min(\operatorname{sgn}(v_{i+\frac{1}{2}}^{n-1}),0), \quad a_{i+\frac{1}{2}}^{out,n-1} = \min(\operatorname{sgn}(-v_{i+\frac{1}{2}}^{n-1}),0). \tag{2.9}$$

In order to keep second order accuracy of the IIOE scheme, the diffusion part is treated by the Crank-Nicolson approach. Finally we end up with the following system

$$\begin{aligned} u_i^{n,k} + \frac{\tau}{2h} \left(a_{i-\frac{1}{2}}^{in,n-1} v_{i-\frac{1}{2}}^{n,k} + \frac{\sigma}{h} \right) \left(u_i^{n,k} - u_{i-1}^{n,k} \right) + \frac{\tau}{2h} \left(a_{i+\frac{1}{2}}^{in,n-1} v_{i+\frac{1}{2}}^{n,k} + \frac{\sigma}{h} \right) \left(u_i^{n,k} - u_{i+1}^{n,k} \right) \neq &2.10) \\ u_i^{n-1} - \frac{\tau}{2h} \left(\left(a_{i-\frac{1}{2}}^{out,n-1} v_{i-\frac{1}{2}}^{n-1} + \frac{\sigma}{h} \right) \left(u_i^{n-1} - u_{i-1}^{n-1} \right) + \left(a_{i+\frac{1}{2}}^{out,n-1} v_{i+\frac{1}{2}}^{n-1} + \frac{\sigma}{h} \right) \left(u_i^{n-1} - u_{i+1}^{n-1} \right) \right), \end{aligned}$$

which we solve iteratively. In every iteration, we calculate the residuum defined as

$$\frac{1}{N}||A(u^{n,k})u^{n,k} - Bu^{n-1}||,$$

where $A(u^{n,k})$ and B are coefficient matrices obtained by writing the system (2.10) using matrix notation and N is the number of unknowns. Our goal was to reach

second order accuracy. For this purpose, while solving the Burgers' equation (1.9), it is sufficient to stop doing iterations when the residuum drops below 10^{-6} . To achieve the desired accuracy for the traffic flow problem (1.22), we stopped the calculation when the residuum drops below 10^{-7} . It means that we have to solve few times (usually from 3 to 6) a tridiagonal system in every time step of the IIOE scheme in case of nonlinear advection-diffusion problems. The tridiagonal system is solved using the Thomas' algorithm. All the numerical experiments are performed on a bounded interval, at the endpoints we have Dirichlet boundary conditions, calculated from the exact solutions at every time step.

2.3.1 Numerical experiments - viscous Burgers' equation

In order to test the numerical scheme (2.10), we chose 4 representative solutions of the viscous Burgers' equation: the traveling-wave solution, the rarefaction-wave solution, the triangular-wave solution and a trigonometric solution. In an earlier work [21] the authors have used the traveling-wave solution to test the numerical scheme similar to (2.10). These experiments were already presented in [16].

Traveling wave. The traveling-wave solution can be obtained by looking for special solutions of the Burgers' equation (1.9) that are only shifted in time. This way we can obtain, see, e.g. [25]

$$u(t,x) = u_r + \frac{1}{2}(u_l - u_r) \left(1 - \tanh\left(\frac{(u_l - u_r)(x - st)}{4\sigma}\right) \right), \qquad (2.11)$$

where $s = (u_l + u_r)/2$ is the propagation speed. Here we chose $u_l = 1$, $u_r = 0$. We will use it as our first example to test the numerical scheme (2.10) on space interval (-0.5, 0.5) and time interval (0, 0.48). First we solve the problem with $\sigma = 0.01$. In this test example we chose a time step $\tau = 4h$. It means that for h = 0.01(n = 100) we use a time step $\tau = 0.04$. Then one can refine the time step and grid size in order to check that the scheme is second order accurate, cf. Table 2.1. The visual comparisons of the numerical and exact solutions for n = 100 are presented in Figure 2.1.

It is interesting to consider what happens if we do not perform any nonlinear iteration. Which means that we calculate the inflow coefficients using the values of the solution from the previous time step. This way we get EOC = 1, as it is documented in Table 2.2. In Figure 2.2 we can observe that the propagation speed of the numerical

solution differs from the exact one. By refining the grid, the speed gets closer to the exact one.

In the second case we decreased the viscosity to $\sigma = 0.001$. In Table 2.3 we show errors and EOC for refined and coarsened grids and time step. One can again see that EOC = 2 also in this example when refining the grid. The lower convergence rate in the beginning is caused by oscillations when the grid size is not sufficiently fine, but as we can see from Figure 2.3 these oscillations are "stable", they do not increase in time and by refining the spatial resolution they are removed completely as documented in Figure 2.4.

Table 2.1: Report on the L_2 errors of IIOE method for the traveling-wave solution (2.11) of the viscous Burgers' equation (1.9) with $\sigma = 0.01$.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.01	0.04	12	$5.0 \ 10^{-3}$	
200	0.005	0.02	24	$1.22 \ 10^{-3}$	2.03
400	0.0025	0.01	48	$3.03 \ 10^{-4}$	2.01
800	0.00125	0.005	96	$7.74 \ 10^{-5}$	1.97
1600	0.000625	0.025	192	$1.90 \ 10^{-5}$	2.02

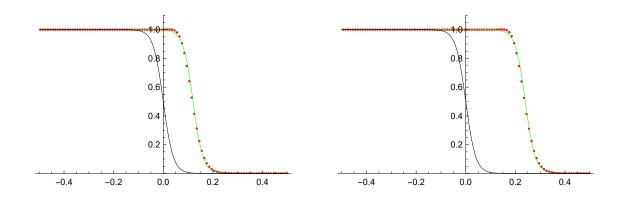


Figure 2.1: Comparing the IIOE scheme with the exact traveling-wave solution (2.11) in time t = 0.24 (left) and t = 0.48 (right), with $\sigma = 0.01$, n = 100, $\tau = 4h$

Table 2.2: Report on the L_2 errors of IIOE method for the traveling-wave solution (2.11) of the viscous Burgers' equation (1.9) with $\sigma = 0.01$, number of nonlinear iterations = 1.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.01	0.04	12	$4.41 \ 10^{-2}$	
200	0.005	0.02	24	$2.32 \ 10^{-2}$	0.93
400	0.0025	0.01	48	$1.18 \ 10^{-2}$	0.97
800	0.00125	0.005	96	$5.96 \ 10^{-3}$	0.99

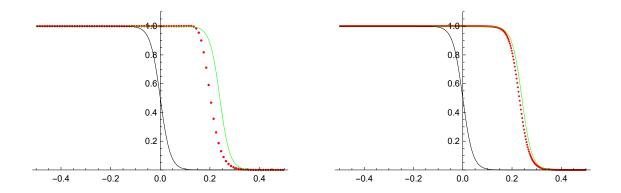


Figure 2.2: Comparing the IIOE scheme with the exact traveling-wave solution (2.11) in time t=0.48, with $\sigma=0.01$, n=100(left), n=400(right), $\tau=4h$, number of nonlinear iterations = 1. We can see that after refining the grid, the propagation speed of the numerical solution is getting closer to the exact speed.

Table 2.3: Report on the L_2 errors of IIOE method for the traveling-wave solution (2.11) of the viscous Burgers' equation (1.9) with $\sigma = 0.001$.

n	h	au	NTS	$L_2(I, L_2)$	EOC
250	0.004	0.016	30	$2.01 \ 10^{-2}$	
500	0.002	0.08	60	$6.84 \ 10^{-3}$	1.55
1000	0.001	0.04	120	$1.79 \ 10^{-3}$	1.94
2000	0.0005	0.02	240	$4.55 \ 10^{-4}$	1.97

Rarefaction wave. The next example is the rarefaction-wave solution (1.15). We chose $u_l = 0$, $u_r = 1$ and $\sigma = 0.01$. First, equation (1.9) is solved by the scheme (2.10) on space interval (-0.5, 0.5) and time interval [0.01, 0.41). Since the exact solution

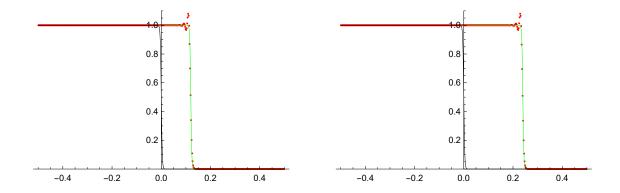


Figure 2.3: Comparing the IIOE scheme with the exact traveling-wave solution (2.11) in time t = 0.24 (left) and t = 0.48 (right), with $\sigma = 0.001$, n = 500, $\tau = 4h$. We can observe small nonincreasingly propagating oscillations.

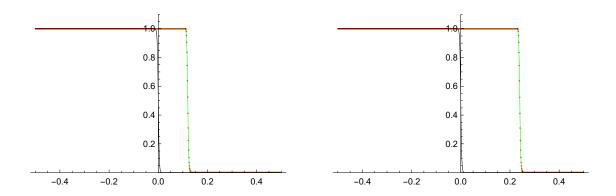


Figure 2.4: Comparing the IIOE scheme with the exact traveling-wave solution (2.11) in time t = 0.24 (left) and t = 0.48 (right), with $\sigma = 0.001$, n = 1000, $\tau = 4h$. On the refined grid the oscillations are gone.

(1.15) is defined for t > 0, we decided to initialize the calculation at time 0.01. The time step τ was chosen to be equal to 4h again. The numerical solution is visually compared to the exact solution in Figure 2.5. The errors are presented in Table 2.4.

Table 2.4: Report on the L_2 errors of IIOE method for the rarefaction-wave solution (1.15) of the viscous Burgers' equation (1.9) with $\sigma = 0.01$.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.01	0.04	10	$5.48 \ 10^{-3}$	
200	0.005	0.02	20	$1.56 \ 10^{-3}$	1.82
400	0.0025	0.01	40	$4.01\ 10^{-4}$	1.96
800	0.00125	0.005	80	$1.00 \ 10^{-4}$	2.00

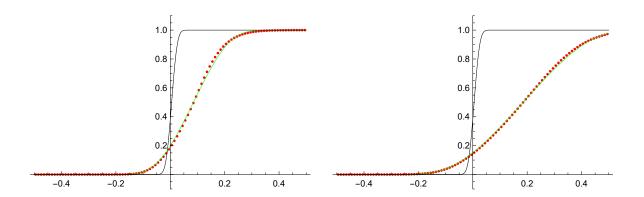


Figure 2.5: Comparing the IIOE scheme with the exact rarefaction-wave solution (1.15) in time t = 0.17 (left) and t = 0.41 (right), with $\sigma = 0.01$, n = 100, $\tau = 4h$

Triangular wave. Our third example is the triangular-wave solution (1.16). In this case the problem (1.9) is solved by the scheme (2.10) on space interval (-0.5, 1.5) and time interval (0.01, 0.51) with $\sigma = 0.02$. Again, the exact solution (1.16) is defined for t > 0 so the numerical calculation was initialized at time 0.01. In Table 2.5 we show the errors for refined grids and time step. When the grid size is not sufficiently fine, we can observe oscillations at the peak of the wave in Figure 2.6. These oscillations do not grow unbounded and can be removed by refining the grid as it is shown in Figure 2.7.

Trigonometric solution. Our last example is the trigonometric solution (1.17) with a = 1.0025, b = 1, $\lambda = \pi^2$ and $\sigma = 0.01$. The errors are reported in Table 2.6. A visual comparison of the numerical results with the exact solution is presented in Figure 2.8.

Table 2.5: Report on the L_2 errors of IIOE method for the triangular-wave solution (1.16) of the viscous Burgers' equation (1.9) with $\sigma = 0.02$.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.02	0.08	5	$3.07 \ 10^{-1}$	
200	0.01	0.04	10	$1.30 \ 10^{-1}$	1.24
400	0.005	0.02	20	$3.72 \ 10^{-2}$	1.81
800	0.0025	0.01	40	$9.02 \ 10^{-3}$	2.04

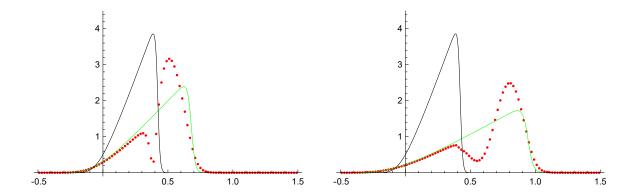


Figure 2.6: Comparing the IIOE scheme with the exact triangular wave solution (1.16) in time t = 0.26 (left), t = 0.50 (right) for n = 100, with $\sigma = 0.02$, $\tau = 4h$.

Table 2.6: Report on the L_2 errors of IIOE method for the trigonometric solution (1.17) of the viscous Burgers' equation (1.9) with $\sigma = 0.01$.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.02	0.08	15	$1.66 \ 10^{-2}$	
200	0.01	0.04	30	$3.18 \ 10^{-3}$	2.38
400	0.005	0.02	60	$5.30 \ 10^{-4}$	2.58
800	0.0025	0.01	120	$1.00 \ 10^{-4}$	2.41

2.3.2 Numerical experiments - traffic flow

In this section we solve the traffic flow equation (1.22) with $v_{max} = 1$. As discussed earlier, by making certain assumptions about the flux function, we can transform the solutions of viscous Burgers' equation (1.9) to obtain a solution for the density of the cars in the traffic flow problem (1.22). Below we give an interpretation and comparison

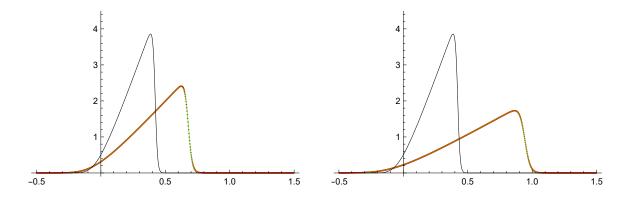


Figure 2.7: Comparing the IIOE scheme with the exact triangular wave solution (1.16) in time t = 0.26 (left), t = 0.50 (right) for n = 800 with $\sigma = 0.02$, $\tau = 4h$.

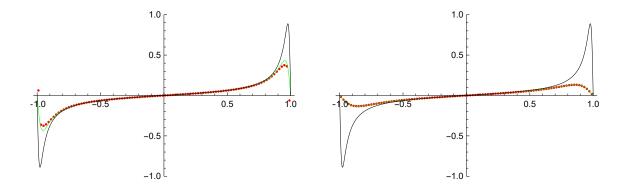


Figure 2.8: Comparing the IIOE scheme with the exact trigonometric solution (1.17) in time t=0.08 (left) and t=0.96 (right) for n=100, with $\sigma=0.01$, b=1, a=1.0025, $\tau=4h$.

of the numerical and exact solutions of the traveling wave and the rarefaction wave solution in a traffic flow context. The experiments were presented in [16].

Traveling wave. A traveling wave can form in a traffic, when incoming cars are approaching a congestion or a line of cars staying behind a red light waiting for turning to green. In the congestion, the density $\rho_r = 1$, which means bumper to bumper traffic. The density of the incoming traffic was chosen to be $\rho_l = 0.1$, which means that there is one car/10 car lengths.

The exact solution of the density ρ was obtained as follows: considering the densities $\rho_l = 0.1$, $\rho_r = 1$ and using the relationship (1.23) we can calculate $u_l = 0.8$ and $u_r = -1$ for the traveling wave solution (2.11) of the Burgers' equation. Then using the relationship (1.24) we obtain the solution for the density ρ . This solution for the traffic density ρ was compared to the numerical solution using the IIOE scheme (2.10).

The errors are documented in Table 2.7. The solution for D=0.01 is presented in Figure 2.9.

Table 2.7: Report on the L_2 errors of IIOE method for the traffic flow problem (1.22) with D = 0.01.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.01	0.04	12	$9.82 \ 10^{-4}$	
200	0.005	0.02	24	$2.36 \ 10^{-4}$	2.05
400	0.0025	0.01	48	$5.94 \ 10^{-5}$	1.99
800	0.00125	0.005	96	$1.53 \ 10^{-5}$	1.96

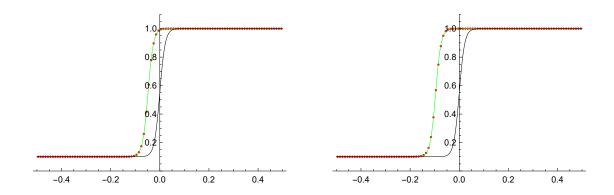


Figure 2.9: Cars from left with traffic density $\rho_l = 0.1$ are approaching the congested region with $\rho_r = 1$. This would happen, for example, when cars are staying behind a red light. The black line is the initial condition for the numerical computation, the red dots are the values computed by the IIOE scheme. The green line is the exact solution. The results of the numerical solution are shown at time t = 0.96, for n = 100, $\tau = 4h$, D = 0.01.

Rarefaction wave. We also interpret the rarefaction wave solution in a traffic flow context. Imagine that the road is divided into two parts by the traffic light positioned at the origin. Cars to the left are staying behind the red light for t < 0, so the density $\rho_l = 1$ there. On the right there is an empty road, $\rho_r = 0$. At time t = 0 the light turns green. This corresponds to the initial condition for a traffic light positioned at x = 0

$$\rho(0,x) = \begin{cases} 1, & x \le 0, \\ 0, & x > 0, \end{cases}$$

Right after the light turned green, the cars start to leave, the traffic rarefies. The exact solution for the density of equation 1.22) was obtained in the same way as in our previous example. First, considering the values $\rho_l = 1$, $\rho_r = 0$ we calculate the corresponding values $u_l = -1$, $u_r = 1$ according to (1.23) for the rarefaction wave solution of the Burgers' equation. Then substituting the exact solution (1.15) to (1.24) we get the density function ρ . Since the initial condition (2.12) is discontinuous we started the computation at time $t^0 = 0.01$. The numerical solution for the density obtained by the IIOE scheme (2.10) is presented for D = 0.01 in Figure 2.10. The errors are presented in Table 2.8.

Table 2.8: Report on the L_2 errors of IIOE method for the traffic flow problem (1.22) with D = 0.01.

n	h	au	NTS	$L_2(I, L_2)$	EOC
100	0.01	0.04	9	$2.49 \ 10^{-3}$	
200	0.005	0.02	18	$9.02 \ 10^{-4}$	1.46
400	0.0025	0.01	36	$2.72 \ 10^{-4}$	1.73
800	0.00125	0.005	72	$7.62 \ 10^{-5}$	1.84

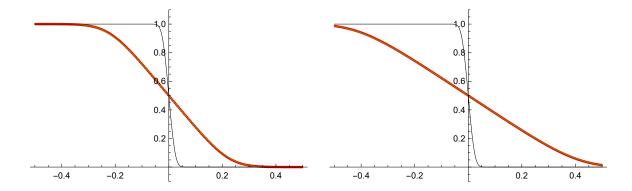


Figure 2.10: At t < 0 cars to the left of the origin are staying behind a red light, the density $\rho_l = 1$. To the right there is an empty road, $\rho_r = 0$. At time t = 0 the traffic light at the origin turns green and cars start to leave. The exact solution(black) is compared with the numerical solution obtained by the IIOE scheme (2.10)(red) at time t = 0.18(left) and t = 0.36(right) for n = 200, $\tau = 4h$, D = 0.01.The initial condition at time $t^0 = 0.01$ is given in black.

2.4 Scalar conservation laws

Let us consider the scalar conservation law of the form

$$u_t + f(u)_x = 0, (2.12)$$

where u = u(t, x) is the unknown scalar quantity, f(u) is a flux function. We can equivalently write the equation in quasi-linear or advective form [18, 33] as

$$u_t + f'(u)u_x = 0 (2.13)$$

where f'(u) is the characteristic speed.

The equation

$$u_t + vu_x = 0, (2.14)$$

is the simplest scalar conservation law with the flux function f(u) = vu with constant characteristic speed f'(u) = v. Considering positive characteristic speed v > 0, the IIOE scheme (2.7) can be written as

$$u_i^n = u_i^{n-1} - \frac{\tau}{h} \left(\left(v u_i^n + v u_{i+1}^{n-1} \right) - \left(v u_{i-1}^n + v u_i^{n-1} \right) \right). \tag{2.15}$$

Similarly, for v < 0 the scheme (2.7) reads

$$u_i^n = u_i^{n-1} - \frac{\tau}{h} \left[\left(v u_i^n + v u_{i-1}^{n-1} \right) - \left(v u_{i+1}^n + v u_i^{n-1} \right) \right]. \tag{2.16}$$

Notice that both (2.15) and (2.16) have the form of a conservative finite volume scheme (2.2), if we define the numerical flux as

$$F_{i-1/2} = \begin{cases} \frac{1}{2} \left(v u_{i-1}^n + v u_i^{n-1} \right) & \text{if } v > 0 \\ \\ \frac{1}{2} \left(v u_i^n + v u_{i-1}^{n-1} \right) & \text{if } v < 0. \end{cases}$$

$$(2.17)$$

Based on this observation, we suggest that for scalar conservation laws (2.12) we should write the IIOE scheme as a conservative finite volume method (2.2) with numerical flux depending on the sign characteristic speed f'(u). We denote $f_i^n := f(u_i^n)$, then the average flux is defined by

$$F_{i-1/2} = \begin{cases} \frac{1}{2} \left(f_{i-1}^n + f_i^{n-1} \right) & \text{if } f'(u) \big|_{x_{i-1/2}} > 0 \\ \\ \frac{1}{2} \left(f_{i-1}^{n-1} + f_i^n \right) & \text{if } f'(u) \big|_{x_{i-1/2}} < 0, \end{cases}$$

$$(2.18)$$

where we calculate the characteristic speed at a given interface from the previous time step n-1, i.e.

$$f'(u)|_{x_{i-1/2}} = \frac{f'(u_{i-1}^{n-1}) + f'(u_i^{n-1})}{2}.$$
 (2.19)

It should be noted that (2.15), (2.16) was already recognized in [23, 24].

If we consider the nonlinear transport equation

$$u_t + \left(\frac{u^2}{2}\right)_x = 0,$$
 (2.20)

with flux given by $f(u) = u^2/2$ and characteristic speed f'(u) = u, which is known as the inviscid Burgers' equation, since it is (1.9) with zero viscosity $\sigma = 0$, the numerical flux at cell interfaces is defined by

$$F_{i-1/2} = \begin{cases} \frac{1}{2} \left[\frac{(u_{i-1}^n)^2}{2} + \frac{(u_i^{n-1})^2}{2} \right] & \text{if } u_{i-1/2} > 0\\ \frac{1}{2} \left[\frac{(u_i^n)^2}{2} + \frac{(u_{i-1}^{n-1})^2}{2} \right] & \text{if } u_{i-1/2} < 0 \end{cases}$$

$$(2.21)$$

where $u_{i-1/2} = (u_{i-1}^{n-1} + u_i^{n-1})/2$. Numerical experiments for the inviscid Burgers' equation will be presented in sections 2.4.2 and 2.5.2.

Remark. For nonlinear conservation laws, where the solutions can develop shocks, it is important that the numerical scheme is conservative, i.e. it can be written in the

form of (2.2). The Lax-Wendroff theorem tells us, that if this holds, the numerical solution of a stable method converges to a weak solution of the differential equation (2.12), as the grid is refined. To ensure that the weak solution is a physically relevant one, the numerical scheme must also satisfy an entropy condition [18].

2.4.1 Formal second order consistency

Theorem 1. For scalar conservation laws (2.12) the IIOE scheme with numerical fluxes at the interfaces defined by (2.18) is formally second order for smooth solutions on a uniform grid and the consistency error is of order $\mathcal{O}(h^2) + \mathcal{O}(\tau h) + \mathcal{O}(\tau^2)$.

Proof. Let us consider, that $u_i^n := u(t^n, x_i)$ is a solution of (2.12) and $f_i^n := f(u(t^n, x_i))$ in the scheme (2.18). The characteristic speed at the interfaces is calculated according to (2.19) from the previous time step. As the numerical fluxes, thus the scheme, depend on the sign of the characteristic speed at the interfaces, we have to distinguish four cases:

Case 1: $f'(u)|_{x_{i-1/2}}$, $f'(u)|_{x_{i+1/2}} > 0$. In that case the IIOE scheme reads

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^n + f_{i+1}^{n-1}) - (f_{i-1}^n + f_i^{n-1}) \right].$$

The Taylor expansion of u in time yields

$$u_i^n = u_i^{n-1} + \tau \partial_t u_i^{n-1} + \frac{\tau^2}{2} \partial_t^2 u_i^{n-1} + \mathcal{O}(\tau^3), \quad u_i^{n-1} = u_i^n - \tau \partial_t u_i^n + \frac{\tau^2}{2} \partial_t^2 u_i^n + \mathcal{O}(\tau^3).$$

Subtracting these two expressions we get

$$u_i^n - u_i^{n-1} = \frac{\tau}{2} \left(\partial_t u_i^{n-1} + \partial_t u_i^n \right) + \underbrace{\frac{\tau^2}{4} \left(\partial_t^2 u_i^{n-1} - \partial_t^2 u_i^n \right)}_{\mathcal{O}(\tau^3)} + \mathcal{O}(\tau^3). \tag{2.22}$$

The Taylor expansion of f in space yields

$$f_{i-1}^{n} = f_{i}^{n} - h\partial_{x}f_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}) = f_{i}^{n} + h\partial_{t}u_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}), \qquad (2.23)$$

$$f_{i+1}^{n-1} = f_i^{n-1} + h\partial_x f_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3) = f_i^{n-1} - h\partial_t u_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3).$$
(2.24)

In both cases we used the relationship between partial derivatives $\partial_t u = \partial_x f$, assuming that the differential equation holds. Now we subtract (2.24) from (2.23) to get

$$f_{i+1}^{n-1} - f_{i-1}^n = f_i^{n-1} - f_i^n - h\left(\partial_t u_i^{n-1} + \partial_t u_i^n\right) + \underbrace{\frac{h^2}{2}\left(\partial_x^2 f_i^{n-1} - \partial_x^2 f_i^n\right)}_{\mathcal{O}(h^2\tau)} + \mathcal{O}(h^3).$$

Multiplying both sides by $\tau/(2h)$ and by algebraic manipulation we get

$$\frac{\tau}{2} \left(\partial_t u_i^{n-1} + \partial_t u_i^n \right) = -\frac{\tau}{2h} \left[\left(f_i^n + f_{i+1}^{n-1} \right) - \left(f_{i-1}^n + f_i^{n-1} \right) \right] + \mathcal{O}(h\tau^2) + \mathcal{O}(h^2\tau). \quad (2.25)$$

Substituting (2.25) to (2.22) we get

$$u_i^n - u_i^{n-1} = -\frac{\tau}{2h} \left[\left(f_i^n + f_{i+1}^{n-1} \right) - \left(f_{i-1}^n + f_i^{n-1} \right) \right] + \mathcal{O}(h\tau^2) + \mathcal{O}(h^2\tau) + \mathcal{O}(\tau^3),$$

where we can recognize the IIOE scheme. Dividing by τ we get the consistency error

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^n + f_{i+1}^{n-1}) - (f_{i-1}^n + f_i^{n-1}) \right] + \mathcal{O}(h\tau) + \mathcal{O}(h^2) + \mathcal{O}(\tau^2). \tag{2.26}$$

Case 2: $f'(u)|_{x_{i-1/2}}$, $f'(u)|_{x_{i+1/2}} \leq 0$, the IIOE scheme then reads

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^{n-1} + f_{i+1}^n) - (f_{i-1}^{n-1} + f_i^n) \right].$$

The Taylor expansion of f in space yields

$$f_{i-1}^{n-1} = f_i^{n-1} - h\partial_x f_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3) = f_i^{n-1} + h\partial_t u_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3),$$
(2.27)

$$f_{i+1}^{n} = f_{i}^{n} + h\partial_{x}f_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}) = f_{i}^{n} - h\partial_{t}u_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}).$$
 (2.28)

Now we subtract (2.27) from (2.28) to get

$$f_{i+1}^{n} - f_{i-1}^{n-1} = f_{i}^{n} - f_{i}^{n-1} - h(\partial_{t}u_{i}^{n} + \partial_{t}u_{i}^{n-1}) + \underbrace{\frac{h^{2}}{2}\left(\partial_{x}^{2}f_{i}^{n} - \partial_{x}^{2}f_{i}^{n-1}\right)}_{\mathcal{O}(h^{2}\tau)} + \mathcal{O}(h^{3}).$$

Multiplying both sides by $\tau/(2h)$ and by algebraic manipulation we get

$$\frac{\tau}{2} \left(\partial_t u_i^n + \partial_t u_i^{n-1} \right) = -\frac{\tau}{2h} \left[\left(f_i^{n-1} + f_{i+1}^n \right) - \left(f_{i-1}^{n-1} + f_i^n \right) \right] + \mathcal{O}(h\tau^2) + \mathcal{O}(h^2\tau). \quad (2.29)$$

Substituting (2.29) to (2.22) we get

$$u_i^n - u_i^{n-1} = -\frac{\tau}{2h} \left[\left(f_i^{n-1} + f_{i+1}^n \right) - \left(f_{i-1}^{n-1} + f_i^n \right) \right] + \mathcal{O}(h\tau^2) + \mathcal{O}(h^2\tau) + \mathcal{O}(\tau^3),$$

where we can recognize the IIOE scheme. Dividing by τ we get the consistency error

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^{n-1} + f_{i+1}^n) - (f_{i-1}^{n-1} + f_i^n) \right] + \mathcal{O}(h\tau) + \mathcal{O}(h^2) + \mathcal{O}(\tau^2). \quad (2.30)$$

Two other cases remain, when the characteristic speed has different sign at the interfaces.

Case 3: $f'(u)|_{x_{i-1/2}} < 0$, $f'(u)|_{x_{i+1/2}} > 0$, for which the scheme reads

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^n + f_{i+1}^{n-1}) - (f_{i-1}^{n-1} + f_i^n) \right],$$

which reduces to

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{f_{i+1}^{n-1} - f_{i-1}^{n-1}}{2h},\tag{2.31}$$

where we can recognize a centered difference for the spatial derivative of the flux function. By Taylor expansion in space of f at time step n-1 we get

$$f_{i+1}^{n-1} = f_i^{n-1} + h\partial_x f_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3) = f_i^{n-1} - h\partial_t u_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3),$$

$$f_{i-1}^{n-1} = f_i^{n-1} - h\partial_x f_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3) = f_i^{n-1} + h\partial_t u_i^{n-1} + \frac{h^2}{2}\partial_x^2 f_i^{n-1} + \mathcal{O}(h^3).$$

After subtraction and dividing by (-2h) we get

$$-\frac{f_{i+1}^{n-1} - f_{i-1}^{n-1}}{2h} = \partial_t u_i^{n-1} + \mathcal{O}(h^2). \tag{2.32}$$

By Taylor expansion of u in time we get

$$u_i^n = u_i^{n-1} + \tau \partial_t u_i^{n-1} + \frac{\tau^2}{2} \partial_t^2 u_i^{n-1} + \mathcal{O}(\tau^3).$$
 (2.33)

By rearranging terms and dividing by τ we get

$$\frac{u_i^n - u_i^{n-1}}{\tau} = \partial_t u_i^{n-1} + \frac{\tau}{2} \partial_t^2 u_i^{n-1} + \mathcal{O}(\tau^2). \tag{2.34}$$

Substituting (2.32) we end up with

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{f_{i+1}^{n-1} - f_{i-1}^{n-1}}{2h} + \frac{\tau}{2} \partial_t^2 u_i^{n-1} + \mathcal{O}(\tau^2) + \mathcal{O}(h^2). \tag{2.35}$$

We want to show that the term $\frac{\tau}{2}\partial_t^2 u_i^{n-1}$ is $\mathcal{O}(\tau h)$. From the conservation law (2.12) we know that $\partial_t u_i^{n-1} = -\partial_x f(u_i^{n-1}) = -f'(u_i^{n-1})\partial_x u_i^{n-1}$. Then

$$\partial_t^2 u_i^{n-1} = -\partial_t \left(f'(u_i^{n-1}) \partial_x u_i^{n-1} \right) = -f''(u_i^{n-1}) \underbrace{\partial_t u_i^{n-1}}_{=-f'(u_i^{n-1}) \partial_x u_i^{n-1}} \partial_x u_i^{n-1} - f'(u_i^{n-1}) \partial_t \partial_x u_i^{n-1} = -f'(u_i^{n-1}) \partial_x u_i^{n-1}$$

$$= f'(u_i^{n-1}) \left(f''(u_i^{n-1}) (\partial_x u_i^{n-1})^2 - \partial_t \partial_x u_i^{n-1} \right), \qquad (2.36)$$

where we took out the common factor $f'(u_i^{n-1})$. Now, similarly as in [23, 24], we want to use the fact that the characteristic speed changes sign inside the cell. If f'(u) is smooth and $f'(u)|_{x_{i-1/2}} < 0$, $f'(u)|_{x_{i+1/2}} > 0$ then there exists some $x_* \in (x_{i-1/2}, x_{i+1/2})$ for which $f'(u)|_{x_*} = 0$. In that case both fluxes represent outflows so we calculate them in the previous time step n-1. From the Taylor expansion of the speed around x_* we get

$$f'(u_i^{n-1}) = \underbrace{f'(u_*^{n-1})}_{=0} + \underbrace{(x_i - x_*)}_{\mathcal{O}(h)} \partial_x f'(u_*^{n-1}) + \mathcal{O}(h^2),$$

from which we see that $\partial_t^2 u_i^{n-1}$ is $\mathcal{O}(h)$ so the term $\frac{\tau}{2} \partial_t^2 u_i^{n-1}$ in (2.35) is $\mathcal{O}(\tau h)$.

Case 4: The last possible case is when $f'(u)|_{x_{i-1/2}} > 0$, $f'(u)|_{x_{i+1/2}} < 0$, for which we have

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{1}{2h} \left[(f_i^{n-1} + f_{i+1}^n) - (f_{i-1}^n + f_i^{n-1}) \right],$$

which simplifies to

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{f_{i+1}^n - f_{i-1}^n}{2h},$$

where similarly as in the previous case, we can recognize a centered difference for the flux function on the right hand side calculated at time step n. The derivation is almost the same as for **Case 3**. We need the following Taylor expansions of f

$$f_{i+1}^{n} = f_{i}^{n} + h\partial_{x}f_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}) = f_{i}^{n} - h\partial_{t}u_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}), \qquad (2.37)$$

$$f_{i-1}^{n} = f_{i}^{n} - h\partial_{x}f_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}) = f_{i}^{n} + h\partial_{t}u_{i}^{n} + \frac{h^{2}}{2}\partial_{x}^{2}f_{i}^{n} + \mathcal{O}(h^{3}).$$
 (2.38)

Subtracting (2.38) from (2.37) and dividing by (-2h) we get

$$-\frac{f_{i+1}^n - f_{i-1}^n}{2h} = \partial_t u_i^n + \mathcal{O}(h^2). \tag{2.39}$$

We need the Taylor expansion of u

$$u_i^{n-1} = u_i^n - \tau \partial_t u_i^n + \frac{\tau^2}{2} \partial_t^2 u_i^n + \mathcal{O}(\tau^3).$$
 (2.40)

By rearranging terms and dividing by τ we get

$$\frac{u_i^n - u_i^{n-1}}{\tau} = \partial_t u_i^n + \frac{\tau}{2} \partial_t^2 u_i^n + \mathcal{O}(\tau^2). \tag{2.41}$$

Substituting (2.39) we end up with

$$\frac{u_i^n - u_i^{n-1}}{\tau} = -\frac{f_{i+1}^n - f_{i-1}^n}{2h} + \frac{\tau}{2}\partial_t^2 u_i^n + \mathcal{O}(\tau^2) + \mathcal{O}(h^2). \tag{2.42}$$

It remains to show that the term $\frac{\tau}{2}\partial_t^2 u_i^n$ is $\mathcal{O}(\tau h)$. We use the notation $f'(u(t^n, x_i)) = f'(u_i^n)$. From the conservation law (2.12) we know that $\partial_t u_i^n = -\partial_x f(u_i^n) = -f'(u_i^n)\partial_x u_i^n$. Then

$$\partial_t^2 u_i^n = -\partial_t \left(f'(u_i^n) \partial_x u_i^n \right) = -f''(u_i^n) \underbrace{\partial_t u_i^n}_{=-f'(u_i^n) \partial_x u_i^n} \partial_x u_i^n - f'(u_i^n) \partial_t \partial_x u_i^n =$$

$$= -f'(u_i^n) \left(f''(u_i^n) (\partial_x u_i^n)^2 - \partial_t \partial_x u_i^n \right), \tag{2.43}$$

where we took out the common factor $f'(u_i^n)$. As before, we use the fact that the characteristic speed changes sign inside the cell. We assume, that f'(u) is smooth in space and time, then $f'(u_i^n) = f'(u_i^{n-1}) + \mathcal{O}(\tau)$. We have already shown, that $f'(u_i^{n-1})$

is of order $\mathcal{O}(h)$, from which we see that $\partial_t^2 u_i^n$ is $\mathcal{O}(h) + \mathcal{O}(\tau)$ so the term $\frac{\tau}{2} \partial_t^2 u_i^n$ in (2.42) is $\mathcal{O}(\tau h) + \mathcal{O}(\tau^2)$.

2.4.2 Numerical experiments

In order to have some evidence that the IIOE scheme with numerical fluxes defined by (2.18) is second order accurate for smooth solutions of conservation laws of the form (2.12), we performed numerical experiments for solutions of the advection equation with constant speed (2.14) and nonlinear transport equation (2.20). In both cases, an exact Dirichlet boundary condition is given at the inflow endpoint.

Advection with constant speed. First we chose a solution of the advection equation (2.14) with v = 1 and initial condition given by

$$u^{0}(x) = 1 + \frac{1}{2} \left(1 - \tanh\left(\frac{x + 0.5}{0.2}\right) \right).$$
 (2.44)

The exact solution is given by

$$u(t,x) = u^0(x-t).$$

The numerical solution was obtained by the IIOE scheme with numerical flux defined by (2.15). The computation was performed on a space interval (-1,1) and time interval (0, 1). The results are reported in Table 2.9. A visual comparison of the numerical solution with the exact one can be seen in Figure 2.11.

Table 2.9: Report on the $L_1(I, L_1)$ errors of the IIOE scheme for an exact smooth solution of the linear advection equation (2.14), $\tau = 4h$.

n	h	au	NTS	$L_1(I,L_1)$	EOC
80	0.025	0.1	10	$2.66 \ 10^{-2}$	
160	0.0125	0.05	20	$7.40 \ 10^{-3}$	1.85
320	0.00625	0.025	40	$1.89 \ 10^{-3}$	1.97
640	0.003125	0.0125	80	$4.71 \ 10^{-4}$	2.00

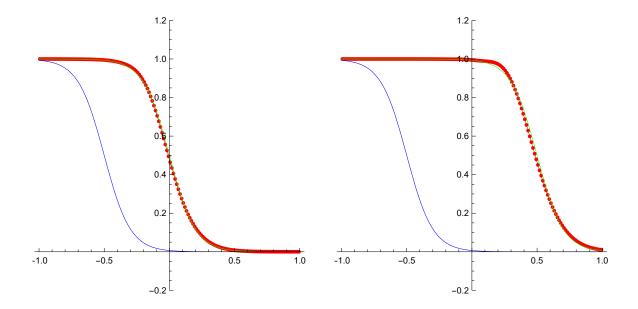


Figure 2.11: Comparing the IIOE scheme with an exact smooth solution for the linear advection equation (2.14) with constant speed v=1 with numerical flux defined by (2.17) in time t=0.5(left) and t=1(right), n=160, $\tau=4h$. The initial profile (2.44) is given in blue, the exact solution in green and the numerical solution in red.

Inviscid Burgers' equation. For the nonlinear equation (2.20) we chose a smooth rarefaction wave solution with given initial condition

$$u^{0}(x) = \frac{\arctan(10 x)}{\pi} + 1/2. \tag{2.45}$$

The exact solution can be obtained by the method of characteristics, see, e.g. [25, 18, 33]. It can be shown, that if the initial profile is non-decreasing everywhere, then the solution does not develop discontinuities for t > 0 [25, 33]. The numerical solution was computed on space interval (-2, 2) and time (0, 1). The $L_1(I, L_1)$ errors and EOC are reported in Table 2.10. The numerical solution is visually compared to the exact one in Figure 2.12.

Table 2.10: Report on the $L_1(I, L_1)$ errors of the IIOE scheme for a smooth rarefaction wave solution of the inviscid Burgers' equation (2.20) with numerical flux defined by (2.21), $\tau = 4h$.

n	h	au	NTS	$L_1(I,L_1)$	EOC
80	0.05	0.2	5	$2.32 \ 10^{-2}$	
160	0.025	0.1	10	$6.93 \ 10^{-3}$	1.74
320	0.0125	0.05	20	$1.86 \ 10^{-3}$	1.90
640	0.00625	0.025	40	$4.70 \ 10^{-4}$	1.98

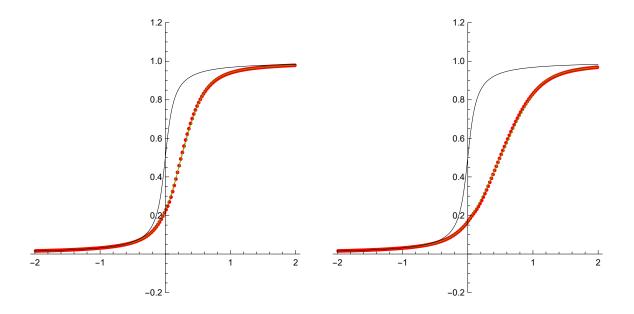


Figure 2.12: Comparing the IIOE scheme with the exact smooth rarefaction wave solution of the inviscid Burgers' equation (2.20) with numerical flux defined by (2.21) in time t = 0.5(left) and t = 1(right), n = 160, $\tau = 4h$. The initial profile (2.45) is given in black, the exact solution in green and the numerical solution in red.

2.5 Flux Limited IIOE method

Higher order schemes suffer from numerical dispersion. Nonphysical oscillations can occur near sharp changes of the solution. A remedy to this is to use limiters [18].

While calculating the flux at a given interface, previously we took both terms in (2.18) with equal weight 1/2. This suggests that we should, as a possible improvement, calculate a weighted average instead:

$$F_{i-1/2}^{\theta} = \begin{cases} (1 - \theta_{i-1/2}) f_{i-1}^{n} + \theta_{i-1/2} f_{i}^{n-1} & \text{if } f'(u) \big|_{x_{i-1/2}} > 0 \\ \theta_{i-1/2} f_{i-1}^{n-1} + (1 - \theta_{i-1/2}) f_{i}^{n} & \text{if } f'(u) \big|_{x_{i-1/2}} < 0 \end{cases}$$

$$(2.46)$$

for some weighting coefficient $\theta_{i-1/2}$. One should notice that for $\theta_{i-1/2} = \frac{1}{2}$ we get the basic IIOE scheme and for $\theta_{i-1/2} = 0$ we get the implicit upwind scheme.

It is important that we assign a weighting coefficient to a given interface, so the method stays conservative. This is not true for the limiter used in previous works [23, 24].

Remark. The weighted averaging of implicit and explicit terms in (2.46) can be interpreted as a flux-limiter method [18], where we combine a high-order scheme with a low-order one. The goal is to use the higher-order scheme as much as possible without generating nonphysical oscillations. The flux is calculated as follows:

$$F_{i-1/2}^{\Phi} = F_{i-1/2}^{L} + \Phi_{i-1/2}(F_{i-1/2}^{H} - F_{i-1/2}^{L}), \tag{2.47}$$

where $F_{i-1/2}^L$ is a flux of a low-order scheme, which does not make new minima or maxima, neither accentuates already existing ones, $F_{i-1/2}^H$ is a high-order flux, which is more accurate but nonphysical oscillations might occur in the numerical solution, $\Phi_{i-1/2}$ is a flux limiter $0 \leq \Phi_{i-1/2} \leq 1$. Notice that for $\Phi_{i-1/2} = 0$ the flux consists only of the low-order flux, for $\Phi_{i-1/2} = 1$ we have the higher-order flux. The weighted averaging in (2.46) described previously is equivalent to flux-limiter method, if the flux of the higher-order scheme is the basic IIOE (2.18) and the low-order flux comes from the implicit upwind scheme. Then the limited flux (2.47) for f'(u) > 0 can be written as

$$F_{i-1/2}^{\Phi} = f(u_{i-1}^n) + \Phi_{i-1/2} \left(\frac{1}{2} f(u_{i-1}^n) + \frac{1}{2} f(u_i^{n-1}) - f(u_{i-1}^n) \right)$$

$$= \left(1 - \frac{\Phi_{i-1/2}}{2} \right) f(u_{i-1}^n) + \frac{\Phi_{i-1/2}}{2} f(u_i^{n-1}).$$
(2.48)

We can see that the two approaches are equivalent by choosing $\theta_{i-1/2} = \frac{\Phi_{i-1/2}}{2}$. This implies that we should expect that our weighting coefficient is bounded $0 \le \theta_{i-1/2} \le 1/2$. We calculate $F_{i+1/2}^{\Phi}$ analogously.

An interesting question is how to calculate the weighting coefficient. We can use a similar approach as in [4, 34, 23, 24], which means that we calculate the weighting coefficients in a way to ensure that the numerical solution does not accentuates existing minima and maxima.

2.5.1 Advection with constant speed

Here we present a possible way for the limiting process in case of linear advection (2.14), considering the characteristic speed v > 0 is positive everywhere. In this case the scheme reads

$$u_i^n = u_i^{n-1} - v \frac{\tau}{h} \left[\left((1 - \theta_{i+1/2}) u_i^n + \theta_{i+1/2} u_{i+1}^{n-1} \right) - \left((1 - \theta_{i-1/2}) u_{i-1}^n + \theta_{i-1/2} u_i^{n-1} \right) \right].$$
(2.49)

In this case the only unknowns in the equation are the weighting coefficients and the value u_i^n . We calculate the weighting coefficients to ensure that the value in the new time step will not be greater as an appropriately defined maximum value $u_{i,max}^n$ neither smaller then a minimum value $u_{i,min}^n$, i.e.

$$u_{i,min}^n \le u_i^n \le u_{i,max}^n$$
.

A simple choice for $u_{i,min}^n$, $u_{i,max}^n$ is for example

$$u^n_{i,min} = \min(u^{n-1}_{i-|c|}, \ u^{n-1}_{i-|c|-1}), \quad u^n_{i,max} = \max(u^{n-1}_{i-|c|}, \ u^{n-1}_{i-|c|-1})$$

where c is the Courant number $c = v\tau/h$. We apply the floor function $\lfloor c \rfloor$, which returns the largest integer less then or equal to c. A similar approach was used in [23, 24]. We start the calculation at the left end of the interval, which is the inflow boundary, where we know the value u_0^n from the given Dirichlet boundary condition. We set $\theta_{1/2} = 1/2$, $\theta_{3/2} = 1/2$, which means that we calculate u_1^n with the basic IIOE scheme. If $u_1^n < u_{i,min}^n$ we adjust $\theta_{3/2}$ in order to get $u_1^n \ge u_{i,min}^n$. The case if $u_1^n > u_{i,max}^n$ is treated analogously.

Below we describe the steps of our flux limiting approach, the FLIIOE method. In general, for the *i*-th cell we do the following:

- 1. u_{i-1}^n and $\theta_{i-1/2}$ are known.
- 2. Set $\theta_{i+1/2} = 1/2$.
- 3. Calculate u_i^n from (2.49)

$$u_i^n = \frac{u_i^{n-1} - c \left[\theta_{i+1/2} u_{i+1}^{n-1} - (1 - \theta_{i-1/2}) u_{i-1}^n - \theta_{i-1/2} u_i^{n-1} \right]}{1 + c(1 - \theta_{i+1/2})}.$$
 (2.50)

4. If $u_i^n < u_{i,min}^n$ it is desired that

$$u_{i,min}^{n} \le \frac{u_{i}^{n-1} - c \left[\theta_{i+1/2} u_{i+1}^{n-1} - (1 - \theta_{i-1/2}) u_{i-1}^{n} - \theta_{i-1/2} u_{i}^{n-1} \right]}{1 + c(1 - \theta_{i+1/2})},$$

from which we get

$$\theta_{i+1/2} \le \frac{u_i^{n-1} + c\left(\theta_{i-1/2}u_{i-1}^n + \theta_{i-1/2}u_i^{n-1}\right) - u_{i,min}^n(1+c)}{c(u_{i+1}^{n-1} - u_{i,min}^n)}.$$

Thus we redefine

$$\theta_{i+1/2} = \min\left(\max\left(\frac{u_i^{n-1} + c\left(\theta_{i-1/2}u_{i-1}^n + \theta_{i-1/2}u_i^{n-1}\right) - u_{i,min}^n(1+c)}{c(u_{i+1}^{n-1} - u_{i,min}^n)}, 0\right), 1/2\right). \quad (2.51)$$

5. If $u_i^n > u_{i,max}^n$ it is desired that

$$u_{i,max}^{n} \ge \frac{u_{i}^{n-1} - c \left[\theta_{i+1/2} u_{i+1}^{n-1} - (1 - \theta_{i-1/2}) u_{i-1}^{n} - \theta_{i-1/2} u_{i}^{n-1}\right]}{1 + c(1 - \theta_{i+1/2})},$$

from which we get

$$\theta_{i+1/2} \ge \frac{u_i^{n-1} + c\left(\theta_{i-1/2}u_{i-1}^n + \theta_{i-1/2}u_i^{n-1}\right) - u_{i,max}^n(1+c)}{c(u_{i+1}^{n-1} - u_{i,max}^n)}.$$

Thus we redefine

$$\theta_{i+1/2} = \min\left(\max\left(\frac{u_i^{n-1} + c\left(\theta_{i-1/2}u_{i-1}^n + \theta_{i-1/2}u_i^{n-1}\right) - u_{i,max}^n(1+c)}{c(u_{i+1}^{n-1} - u_{i,max}^n)}, 0\right), 1/2\right). \quad (2.52)$$

6. We calculate u_i^n by (2.50) as in step 3 with a possibly modified $\theta_{i+1/2}$.

We performed numerical experiments in section 2.5.3.

2.5.2 Inviscid Burgers' equation

We describe a possible way to limit the scheme also in the case of a nonlinear equation. Again, for simplicity we assume that the characteristic speed $f'(u) = u \ge 0$

is non-negative everywhere The scheme then reads

$$u_{i}^{n} = u_{i}^{n-1} - \frac{\tau}{h} \left[\left((1 - \theta_{i+1/2}) \frac{(u_{i}^{n})^{2}}{2} + \theta_{i+1/2} \frac{(u_{i+1}^{n-1})^{2}}{2} \right) - \left((1 - \theta_{i-1/2}) \frac{(u_{i-1}^{n})^{2}}{2} + \theta_{i-1/2} \frac{(u_{i}^{n-1})^{2}}{2} \right) \right]$$
(2.53)

We use similar ideas as in the case of linear transport. We want to ensure that our numerical solution does not violate the min-max principle. Since the speed is not constant, we calculate $u_{i,min}^n$ and $u_{i,max}^n$ differently. First we calculate the Courant number using the maximum value of the characteristic speed at the previous time step, which is $\max_i(u_i^{n-1})$ in the case of the inviscid Burgers' equation. We calculate the maximal Courant number

$$c = \max_{i} (u_i^{n-1})\tau/h. \tag{2.54}$$

Then the limiting values are calculated as

$$u_{i,min}^n = \min_{i-|c|-1}^i (u_i^{n-1}), \quad u_{i,max}^n = \max_{i-|c|-1}^i (u_i^{n-1}).$$

The procedure is then analogous to the linear case, except we use (2.53). Below we describe the steps of the flux limiting FLIIOE method for the inviscid Burgers' equation. In general, for the *i*-th cell we do the following:

- 1. u_{i-1}^n and $\theta_{i-1/2}$ are known.
- 2. We set $\theta_{i+1/2} = 1/2$.
- 3. Calculate u_i^n from (2.53)

$$u_i^n = \frac{-1 \pm \sqrt{1 + \frac{\tau}{h}(1 - \theta_{i+1/2})(\frac{\tau}{h}F_{i-1/2}^{\theta} + 2u_i^{n-1} - \frac{\tau}{h}\theta_{i+1/2}(u_{i+1}^{n-1})^2)}}{\frac{\tau}{h}(1 - \theta_{i+1/2})}, \quad (2.55)$$

where

$$F_{i-1/2}^{\theta} = (1 - \theta_{i-1/2}) \frac{(u_{i-1}^n)^2}{2} + \theta_{i-1/2} \frac{(u_i^{n-1})^2}{2}.$$

- We have to decide, which solution to choose. Since the numerical experiments we performed all have positive initial values, we chose the positive one from the two.
- If the expression under the square root is negative, we fail to calculate a real solution. In this case we set $\theta_{i+1/2} = 0$, i.e. at $x_{i-1/2}$ we use the implicit upwind flux.

4. If $u_i^n < u_{i,min}^n$ it is desired that

$$u_{i,min}^{n} \leq \frac{-1 + \sqrt{1 + \frac{\tau}{h}(1 - \theta_{i+1/2})(\frac{\tau}{h}F_{i-1/2}^{\theta} + 2u_{i}^{n-1} - \frac{\tau}{h}\theta_{i+1/2}(u_{i+1}^{n-1})^{2})}}{\frac{\tau}{h}(1 - \theta_{i+1/2})},$$

from which we get

$$\theta_{i+1/2} \le \frac{u_i^{n-1} - u_{i,min}^n + \frac{\tau}{h} (F_{i-1/2}^\theta - (u_{i,min}^n)^2 / 2)}{\frac{\tau}{2h} ((u_{i,min}^n)^2 - (u_{i+1}^{n-1})^2)}.$$
 (2.56)

Thus we redefine

$$\theta_{i+1/2} = \min\left(\max\left(\frac{u_i^{n-1} - u_{i,min}^n + \frac{\tau}{h}(F_{i-1/2}^\theta - (u_{i,min}^n)^2/2)}{\frac{\tau}{2h}((u_{i,min}^n)^2 - (u_{i+1}^{n-1})^2)}, 0\right), 1/2\right). \quad (2.57)$$

5. If $u_i^n > u_{i,max}^n$ we want

$$u_{i,max}^{n} \ge \frac{-1 + \sqrt{1 + \frac{\tau}{h}(1 - \theta_{i+1/2})(\frac{\tau}{h}F_{i-1/2}^{\theta} + 2u_{i}^{n-1} - \frac{\tau}{h}\theta_{i+1/2}(u_{i+1}^{n-1})^{2})}}{\frac{\tau}{h}(1 - \theta_{i+1/2})},$$

from which we get

$$\theta_{i+1/2} \ge \frac{u_i^{n-1} - u_{i,max}^n + \frac{\tau}{h} (F_{i-1/2}^{\theta} - (u_{i,max}^n)^2 / 2)}{\frac{\tau}{2h} ((u_{i,max}^n)^2 - (u_{i+1}^{n-1})^2)}.$$
 (2.58)

Thus we redefine

$$\theta_{i+1/2} = \min\left(\max\left(\frac{u_i^{n-1} - u_{i,max}^n + \frac{\tau}{h}(F_{i-1/2}^{\theta} - (u_{i,max}^n)^2/2)}{\frac{\tau}{2h}((u_{i,max}^n)^2 - (u_{i+1}^{n-1})^2)}, 0\right), 1/2\right). (2.59)$$

6. We calculate u_i^n by (2.55) with a possibly modified $\theta_{i+1/2}$.

We performed numerical experiments in section 2.5.4.

2.5.3 Numerical experiments - Advection with constant speed

In this section we present numerical experiments performed by the flux limited FLIIOE scheme in case of advection with constant speed. In this linear case (2.49) is compared with the stabilized schemes S¹IIOE (without quadratic reconstruction) and S²IIOE (with quadratic reconstruction) described in [23, 24]. For this purpose we use the same examples that can be found in these works: a **smooth profile** with initial condition

$$u^{0}(x) = \max(0, \cos^{5}(\pi(x+0.5))), \tag{2.60}$$

and a piecewise constant profile with initial condition

$$u^{0}(x) = \begin{cases} 1, & x \in [-0.75, -0.25], \\ 0, & \text{otherwise.} \end{cases}$$

In both cases we solve (2.14) numerically with v = 1 on space interval (-1, 1) and time interval (0, 1). The exact solution is calculated as $u(t, x) = u^0(x - t)$. At the inflow boundary exact Dirichlet boundary condition is given.

The values for S²IIOE (with quadratic reconstruction) are taken from previous works [22, 24]. For the smooth profile (2.60) we can observe that the flux limited scheme (2.49) performs better than the S¹IIOE and it is comparable with the S²IIOE in all cases. The results are documented in Table 2.11. In Figure 2.13 we compare S¹IIOE with the flux limited scheme (2.49) visually.

In the case of a discontinuous profile (2.61) the flux limited scheme (2.49) performs better than both stabilized schemes, c.f. Table 2.12. A visual comparison with S¹IIOE can be seen in Figure 2.14.

Table 2.11: Report on the $L_1(I, L_1)$ errors of the S¹HOE scheme without quadratic reconstruction, S²HOE scheme with quadratic reconstruction, the flux limited FLHOE scheme (2.49) for a smooth initial profile (2.60), $\tau = h$.

	S ¹ IIOE	S ¹ IIOE	S ² IIOE	S ² IIOE	FLIIOE	FLIIOE
n	$L_1(I, L_1)$	EOC	$L_1(I, L_1)$	EOC	$L_1(I, L_1)$	EOC
40	$9.82 \ 10^{-2}$		$9.19 \ 10^{-2}$		$7.27 \ 10^{-2}$	
80	$3.38 \ 10^{-2}$	1.54	$2.80 \ 10^{-2}$	1.71	$2.63 \ 10^{-2}$	1.47
160	$1.01 \ 10^{-2}$	1.74	$7.69 \ 10^{-3}$	1.86	$7.55 \ 10^{-3}$	1.80
320	$2.73 \ 10^{-3}$	1.89	$1.97 \ 10^{-3}$	1.96	$1.96 \ 10^{-3}$	1.95
640	$7.11 \ 10^{-4}$	1.94	$4.97 \ 10^{-4}$	1.99	$5.22 \ 10^{-4}$	1.91
1280	$1.83 \ 10^{-4}$	1.96	$1.24 \ 10^{-4}$	2.00	$1.41 \ 10^{-4}$	1.89

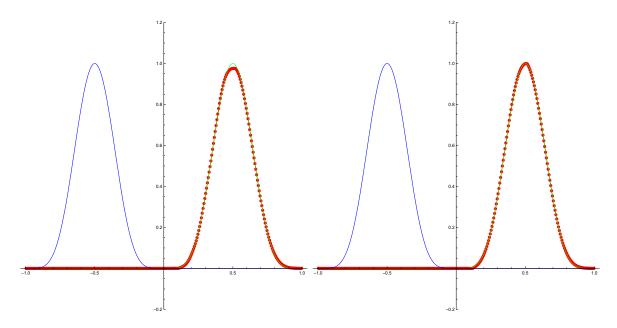


Figure 2.13: Comparing the S¹IIOE scheme without quadratic reconstruction (left) and the flux limited FLIIOE scheme (2.49) (right) with the exact solution of (2.14) for a smooth initial profile (2.60) in time t = 1, n = 320, $\tau = h$.

Table 2.12: Report on the $L_1(I, L_1)$ errors of the S¹IIOE scheme without quadratic reconstruction, S²IIOE scheme with quadratic reconstruction and the flux limited FLIIOE scheme (2.49) for a piecewise constant initial profile (2.61), $\tau = h$.

	S^1IIOE	S^1IIOE	S ² IIOE	S^2IIOE	FLIIOE	FLIIOE
n	L_1 error	EOC	L_1 error	EOC	L_1 error	EOC
40	$2.03 \ 10^{-1}$		$2.02 \ 10^{-1}$		$1.49 \ 10^{-1}$	
80	$1.31 \ 10^{-1}$	0.63	$1.30 \ 10^{-1}$	0.63	$9.40 \ 10^{-2}$	0.66
160	$8.38 \ 10^{-2}$	0.64	$8.38 \ 10^{-2}$	0.64	$5.91 \ 10^{-2}$	0.67
320	$5.35 \ 10^{-2}$	0.65	$5.34 \ 10^{-2}$	0.65	$3.72 \ 10^{-2}$	0.67
640	$3.41 \ 10^{-2}$	0.65	$3.41 \ 10^{-2}$	0.65	$2.35 \ 10^{-2}$	0.66
1280	$2.16 \ 10^{-2}$	0.66	$2.16 \ 10^{-2}$	0.66	$1.48 \ 10^{-2}$	0.67

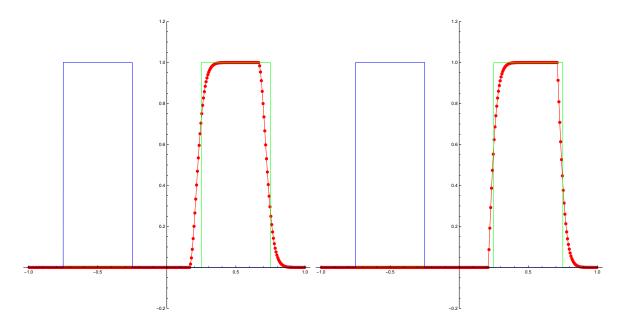


Figure 2.14: Comparing the S¹IIOE scheme without quadratic reconstruction (left) and the FLIIOE scheme (2.49) (right) with the exact solution of (2.14) with a piecewise constant initial profile (2.61) in time t = 1, n = 320, $\tau = h$.

2.5.4 Numerical experiments - Inviscid Burgers' equation

We performed numerical experiments also for the nonlinear transport equation (2.20). For this purpose the following examples where chosen: shock wave, rarefaction wave and triangular wave. As in the linear case, at the inflow boundary exact Dirichlet boundary condition is given.

Shock wave. In this case the scheme (2.53) is solved on the space interval (-0.5, 0.5) and time interval (0, 0.5) with a step function initial condition

$$u^{0}(x) = \begin{cases} 1, & x \le 0, \\ 0, & x > 0. \end{cases}$$

A physically relevant weak solution [25, 18, 33] is a traveling shock wave

$$u(t,x) = u^0(x - s t),$$

where s = 1/2 is the speed of the shock propagation. The $L_1(I, L_1)$ errors and EOC of the computation are documented in Table 2.13. A visual comparison of the numerical solution with the exact one can be seen in Figure 2.15.

Table 2.13: Report on the $L_1(I, L_1)$ errors of the flux limited FLIIOE scheme (2.53) for a shock wave solution (2.61) of the inviscid Burgers' equation (2.20). We used time steps $\tau = 4h$ in the first 4 rows and $\tau = 32h$ in the next 4 rows.

n	h	au	NTS	$L_1(I,L_1)$	EOC
80	0.0125	0.05	10	$5.03 \ 10^{-3}$	
160	0.0125	0.025	20	$2.52 \ 10^{-3}$	0.99
320	0.003125	0.0125	40	$1.26 \ 10^{-3}$	1.00
640	0.0015625	0.00625	80	$6.32 \ 10^{-3}$	1.00
320	0.03125	0.1	5	$6.78 \ 10^{-3}$	
640	0.0015625	0.05	10	$3.39 \ 10^{-3}$	1.00
1280	0.00078125	0.025	20	$1.69 \ 10^{-3}$	1.00
2560	0.000391	0.0125	40	$8.47 \ 10^{-4}$	1.00

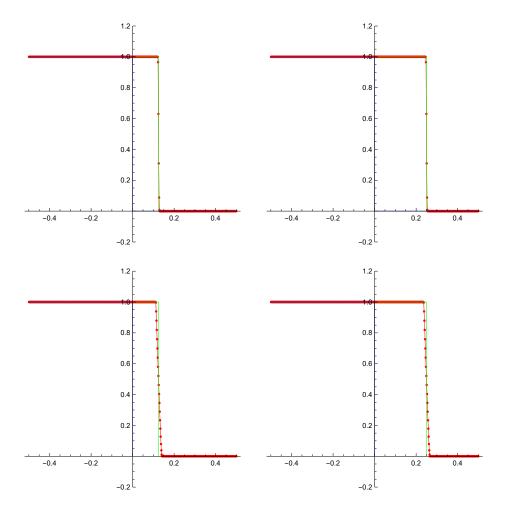


Figure 2.15: Results of the FLIIOE(flux limited) scheme (2.53) with the traveling shock wave solution (2.61) of the inviscid Burgers' equation (2.20) at time t = 0.25(left) and t = 0.5(right), n = 640 for relatively large time steps $\tau = 4h(\text{top})$, $\tau = 32h(\text{bottom})$. The initial profile is given in blue, the exact solution in green and the numerical solution in red.

Rarefaction wave. The next example is a solution of (2.20) with a discontinuous initial condition

$$u^{0}(x) = \begin{cases} 0, & x \le 0, \\ 1, & x > 0. \end{cases}$$

A physically relevant weak solution is a rarefaction wave [25, 18, 33] for t > 0

$$u(t,x) = \begin{cases} 1, & x > t, \\ x/t, & 0 \le x \le t, \\ 0, & x < 0. \end{cases}$$

We solved (2.20) numerically by (2.53) on space interval (-0.5, 1.5) and time interval (0, 1). The results are documented in Table 2.14. The numerical solution is compared visually with the exact one in Figure 2.16.

Table 2.14: Report on the $L_1(I, L_1)$ errors of the flux limited FLIIOE scheme (2.53) for a rarefaction wave solution (2.61) of the inviscid Burgers' equation (2.20). We used time steps $\tau = h(\text{top})$, $\tau = 2h(\text{bottom})$.

n	h	au	NTS	$L_1(I,L_1)$	EOC
80	0.025	0.025	40	$1.63 \ 10^{-2}$	
160	0.0125	0.0125	80	$8.90 \ 10^{-3}$	0.87
320	0.00625	0.00625	160	$4.70 \ 10^{-3}$	0.92
640	0.003125	0.003125	320	$2.42 \ 10^{-3}$	0.96
80	0.025	0.05	20	$2.21 \ 10^{-2}$	
160	0.0125	0.025	40	$1.22 \ 10^{-2}$	0.86
320	0.00625	0.0125	80	$6.58 \ 10^{-3}$	0.89
640	0.00625	0.003125	160	$3.45 \ 10^{-3}$	0.93

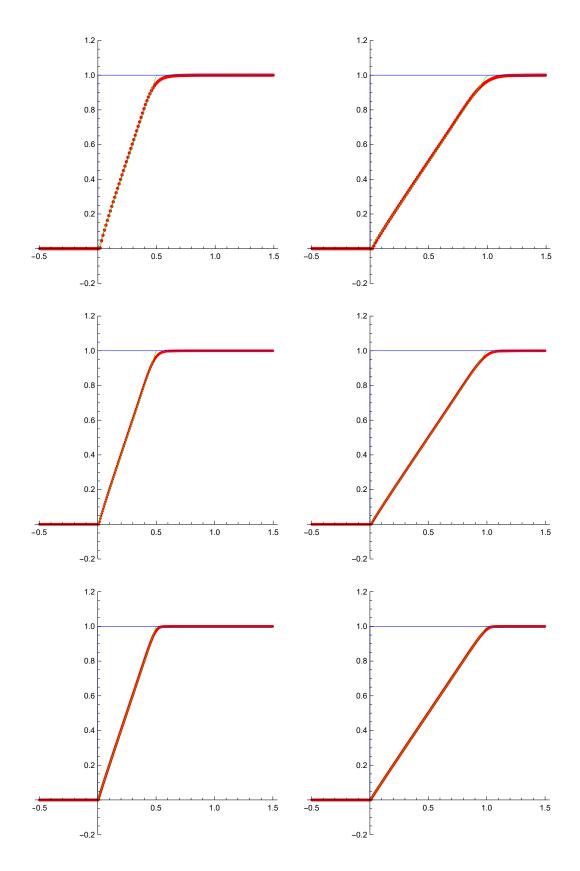


Figure 2.16: Results of the flux limited FLIIOE scheme (2.53) for a rarefaction wave solution (2.61) of the inviscid Burgers' equation (2.20) at time t=0.5(left) and t=1(right), n=160(top), n=320(center), n=640(bottom), $\tau=h$. The initial profile is given in blue, the exact solution in green and the numerical solution in red.

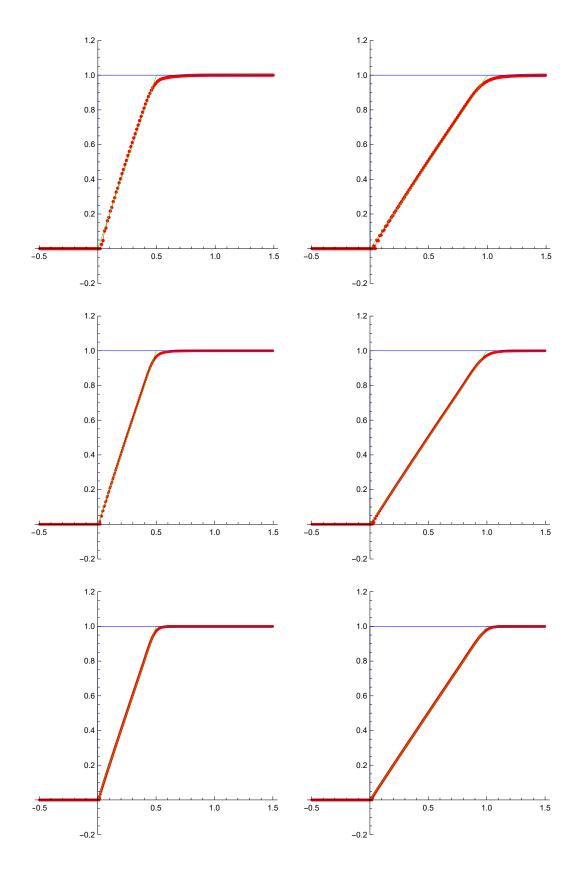


Figure 2.17: Results of the flux limited FLIIOE scheme (2.53) for a rarefaction wave solution (2.61) of the inviscid Burgers' equation (2.20) at time t=0.5(left) and t=1(right), n=160(top), n=320(center), n=640(bottom), $\tau=2h$. The initial profile is given in blue, the exact solution in green and the numerical solution in red.

Triangular wave The third example is another physically relevant weak solution of (2.20) [25, 18, 33] given for $t \ge 0$ as

$$u(t,x) = \begin{cases} x/(1+t), & 0 \le x \le \sqrt{1+t}, \\ 0, & \text{otherwise.} \end{cases}$$

We solved (2.20) numerically by (2.53) on space interval (-0.5, 1.5) and time interval (0, 1). The results are documented in 2.15. The numerical solution is compared visually with the exact one in 2.18.

Table 2.15: Report on the $L_1(I, L_1)$ errors of the flux limited FLIIOE scheme (2.53) for a triangular wave solution (2.61) of the inviscid Burgers' equation (2.20), with time steps $\tau = h$ in the first 4 rows and $\tau = 2h$ in the next 4 rows.

				FLIIOE	
n	h	au	NTS	$L_1(I,L_1)$	EOC
80	0.025	0.025	40	$1.62 \ 10^{-2}$	
160	0.0125	0.0125	80	$7.73 \ 10^{-3}$	1.07
320	0.00625	0.00625	160	$3.66 \ 10^{-3}$	1.08
640	0.003125	0.003125	320	$1.77 \ 10^{-3}$	1.05
80	0.025	0.05	20	$2.40 \ 10^{-2}$	
160	0.0125	0.025	40	$1.14 \ 10^{-2}$	1.07
320	0.00625	0.0125	80	$5.35 \ 10^{-3}$	1.09
640	0.00625	0.003125	160	$2.66 \ 10^{-3}$	1.01

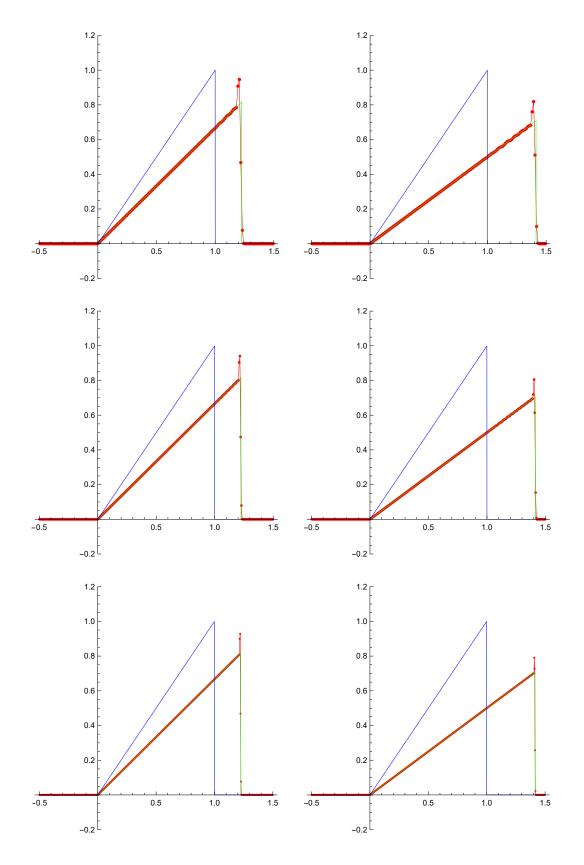


Figure 2.18: Results of the flux limited FLIIOE scheme (2.53) for a triangular wave solution (2.61) of the inviscid Burgers' equation (2.20) at time t = 0.5(left) and t = 1(right), n = 160(top), n = 320(center), n = 640(bottom), $\tau = h$. The initial profile is given in blue, the exact solution in green and the numerical solution in red.

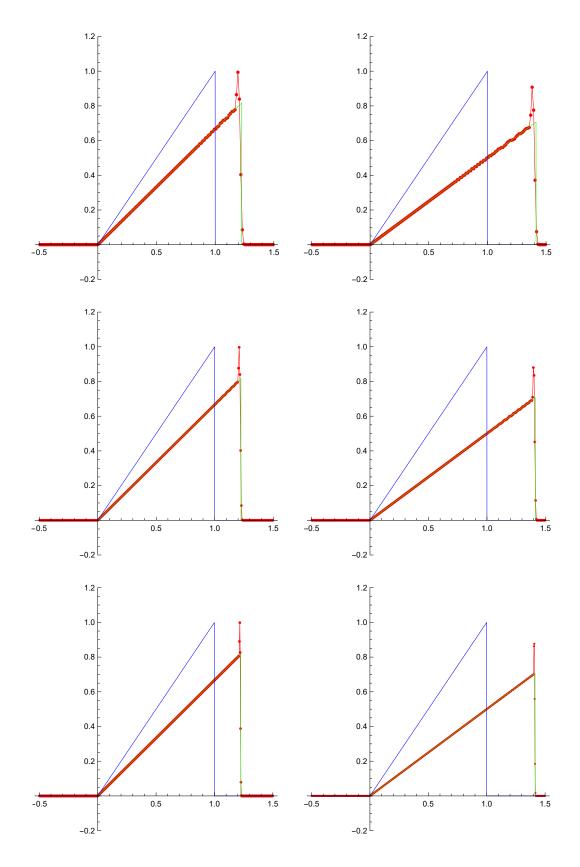


Figure 2.19: Results of the flux limited FLIIOE scheme (2.53) for a triangular wave solution (2.61) of the inviscid Burgers' equation (2.20) at time t = 0.5(left) and t = 1(right), n = 160(top), n = 320(center), n = 640(bottom), $\tau = 2h$. The initial profile is given in blue, the exact solution in green and the numerical solution in red.

Chapter 3

Conclusions

In this work we presented a scheme for solving nonlinear advection-diffusion equations similar to what can be found in [16], where for the advection term we applied the IIOE scheme and for the diffusion term we applied the Crank-Nicolson scheme. We tested the performance of the scheme on the viscous Burgers' equation (1.9) and the traffic flow equation (1.22) by comparing numerical solutions with exact ones. In numerical experiments we can observe second order convergence.

We have shown that the IIOE method can be written in the form of a conservative finite volume method (2.2), which is crucial for first order scalar conservation laws of the form (2.12). We have shown that on a uniform grid the scheme is formally second order accurate for smooth solutions of scalar conservation laws of the form (2.12). We performed numerical experiments for the inviscid Burgers' equation first in the case of a smooth solution, where the results are in accordance with our theoretical expectations.

For problems, where numerical solutions tend to develop spurious oscillations, a possible flux limiting process were presented, where we want to ensure that the numerical solutions is bounded by some minimum and maximum value.

Chapter 4

Thesis objectives and prospective contribution

In the future we plan to perform the following tasks:

- Generalizing the flux-limiter for arbitrary cases, applications for nonlinear inviscid equations and traffic flow problems
- In case of enough time, extension to higher dimensions for other fluid mechanics problems

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