# A NEW TYPE OF SIMPLIFIED INVERSE LAX-WENDROFF BOUNDARY TREATMENT FOR HYPERBOLIC CONSERVATION LAWS 

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#### Abstract

In this paper, we design a new type of simplified high order inverse Lax-Wendroff (ILW) boundary treatment for solving hyperbolic conservation laws with finite difference method on a Cartesian mesh, in which both scalar equations and systems are considered. This new ILW method decomposes the construction of ghost points into two steps: interpolation and extrapolation. At first, we approximate some special point values through an interpolation polynomial based on interior points near the boundary. Then, we construct a Hermite extrapolation polynomial based on those special point values and spatial derivatives at the boundary obtained through the ILW process. This extrapolation polynomial will give us the approximation of the ghost point values. Through an appropriate selection of the interpolation points, high-order accuracy and stable results can be achieved. We use the eigenvalue analysis method to analyze the stability of our new boundary treatment. The analysis results show that the new method can improve the computational efficiency while maintaining accuracy and stability. Numerical tests for one- and two-dimensional problems indicate that our method has high order accuracy for smooth solutions and non-oscillatory property for shock solution near the boundary.


Key Words: Inverse Lax-Wendroff method; numerical boundary treatment; high order accuracy; finite difference method; hyperbolic conservation laws; eigenvalue analysis.

1. Introduction. In this paper, we propose a new high order accuracy boundary treatment based on finite difference methods with fixed Cartesian mesh for hyperbolic conservation law equations. For problems on complex domain under such mesh, there are often two main difficulties. First, the computational stencil of a high order finite difference scheme is often relatively wide, thus we need to evaluate the values at several ghost points near the boundary. Secondly, the physical boundary often does not happen to be on the grid points, so we need to design an algorithm to introduce the boundary conditions into our boundary scheme. If the boundary scheme is not well designed, it may require an extremely small time step to ensure stability, resulting in low computational efficiency.

A common treatment is to use body-fitted grid. That is to establish appropriate body-fitted coordinates so that the grid points coincide with the physical boundary, and then to solve the partial differential equation in the new coordinate system. In this approach, the boundary conditions can be given directly on the grid points. The advantage of this method is that it can accurately meet the given boundary conditions.

[^0]Its disadvantage is that the generation of body-fitted grid could be difficult. The quality of the grid directly determines the computational efficiency and accuracy. Especially for problems with moving boundary, the management of the moving grid is generally complex, which will increase the computational cost greatly. In addition, the governing equation needs to be changed during computation. The transformed PDE is often more complex than the original equation, which will also increase the computational cost.

For non body-fitted mesh methods, many scholars have also proposed some methods, such as the embedded boundary method $[1,9-11,21,27]$, the immersed boundary method [4, 20, 22-24, 33], the ILW (inverse Lax-Wendroff) method [28-31] and so on. In this paper, the method we will introduce is a type of the ILW method.

The prototype of the earliest ILW method comes from the simulation of pedestrian flow $[7,34]$. The pedestrian walking direction can be determined by solving an Eikonal equation. They deal with the boundary conditions by transforming the normal derivative into the tangential derivative. Later, this method was extended to hyperbolic conservation law equations by Tan and Shu [28]. They transformed the normal derivative into time derivative and tangential derivative to deal with the corresponding inflow boundary conditions (different from the original Lax-Wendroff scheme, which transformed the time derivative into spatial derivative, hence the meaning of "inverse"), and applied this method to inviscid compressible fluids.

After the ILW method was proposed, many scholars have done a series of work, which have greatly developed this method. To deal with the heavy algebra of the original ILW method for nonlinear systems (especially in the high-dimensional case), the simplified ILW (SILW) method was proposed in [31], which greatly reduced the computational cost of the ILW method for solving systems. Lu et al. [19] proposed an ILW method to deal with sonic points by evaluating the flux values at ghost points, so it can deal with problems with changing wind direction. Ding et al. [5] redefined the concept of "conservation" for finite difference schemes, and gave an ILW method satisfying conservation in the new sense. In addition to hyperbolic conservation law equations, the ILW method was also applied to other types of equations, such as convection diffusion equation $[13,15,18]$ and Boltzmann equation [6]. For the moving boundary problem, Tan and Shu extended the ILW method to deal with the compressible inviscid fluid containing moving (translational) boundary in [29]. By redefining the material derivative on the boundary, in [3], Cheng et al. extended the method to deal with the arbitrary motion of the boundary, and used it to simulate the interaction between shock wave and rigid body. Liu et al. [17] extended this ILW method to convection-diffusion equations on moving domain, in which a unified algorithm was design for five cases: pure convection, convection-dominated, convectiondiffusion, diffusion-dominated and pure diffusion cases. For the three-dimensional moving boundary problem, Liu et al. [16] extended the moving boundary treatment to the three-dimensional case, and simulated the interaction between inviscid / viscous fluid and three-dimensional rigid body. References [13-15, 32] have analyzed the linear stability of ILW and SILW methods, which provide guidelines for us to design stable ILW boundary treatments.

In this paper, we will design a new type of SILW method for conservation law equations. The new ILW method decomposes the construction of the ghost points into two steps: interpolation and extrapolation. At first, we approximate some special point values through an interpolation polynomial based on interior points near the boundary. Then, we will construct a Hermite extrapolation polynomial based on those special point values and spatial derivatives at the boundary obtained through
the ILW process. This extrapolation polynomial will give us the approximation of the the ghost point values. Through an appropriate selection of the interpolation points, high-order accuracy and stable results can be achieved. The eigenvalue analysis method is used to help us select these interpolation points to ensure the stability of the numerical scheme. The analysis results show that the new method can improve the computational efficiency while maintaining accuracy and stability. Finally, we apply our method to the simulation of inviscid compressible fluid.

The organization of this paper is as follows. In Section 2, we will give the description of the new ILW method for one-dimensional scalar conservation law equations, and use the eigenvalue analysis method to perform the linear stability analysis. In Section 3, we will extend this algorithm to system and high-dimensional cases. The high order accuracy and robustness of our algorithm will be shown through numerical tests in Section 4. Conclusion remarks will be given in Section 5 .
2. The one-dimensional scalar conservation law case. Consider the scalar hyperbolic conservation law in the following form:

$$
\begin{cases}u_{t}+f(u)_{x}=0, & x \in(-1,1), t>0,  \tag{2.1}\\ u(-1, t)=g_{l}(t), & t>0, \\ u(x, 0)=u_{0}(t), & x \in[-1,1] .\end{cases}
$$

We assume that $f^{\prime}(u(-1, t))>0$, such that the left boundary $x=-1$ is an inflow boundary, where a boundary condition needs to be given. We also assume that $f^{\prime}(u(1, t))>0$. Hence the right boundary $x=1$ is an outflow boundary, where no boundary condition is required.

Suppose the domain is divided by the uniform mesh:

$$
\begin{equation*}
-1+C_{a} \Delta x=x_{0}<\cdots<x_{N}=1-C_{b} \Delta x \tag{2.2}
\end{equation*}
$$

with uniform mesh size $\Delta x=2 /\left(C_{a}+C_{b}+N\right)$ and $C_{a}, C_{b} \in[0,1)$. Note that we have deliberately allowed the physical boundary $x= \pm 1$ not coinciding with grid points.

We use the framework of method of lines (MOL) to construct a semi-discrete scheme on the interior point $x_{j}, j=0,1,2, \cdots, N$ :

$$
\begin{equation*}
\frac{d}{d t} u_{j}=L_{h}(u)_{j}, \tag{2.3}
\end{equation*}
$$

where,

$$
L_{h}=-\frac{1}{\Delta x}\left(\hat{f}_{j+1 / 2}-\hat{f}_{j-1 / 2}\right) \approx-\left.f(u)_{x}\right|_{x_{j}}
$$

is the spatial discrete operator. Here, $u_{j}(t)$ is the numerical approximation to the exact solution $u\left(x_{j}, t\right)$, and $\hat{f}_{j+1 / 2}$ is the numerical flux. In this paper, we will use an upwind-biased finite difference conservative scheme to construct $\hat{f}_{j+1 / 2}$, such as the WENO scheme [8].

After the spatial discretization, the semi-discrete scheme (2.3) is a system of ordinary differential equations. For time discretization, we use the total variation diminishing (TVD) Runge-Kutta (RK) scheme [26]. From time level $t^{n}$ to $t^{n+1}$, the third order TVD RK scheme is given as

$$
\begin{align*}
& u_{j}^{(1)}=u_{j}^{n}+\Delta t L_{h}\left(u^{n}\right)_{j}, \\
& u_{j}^{(2)}=\frac{3}{4} u_{j}^{n}+\frac{1}{4} u_{j}^{(1)}+\frac{1}{4} \Delta t L_{h}\left(u^{(1)}\right)_{j},  \tag{2.4}\\
& u_{j}^{n+1}=\frac{1}{3} u_{j}^{n}+\frac{2}{3} u_{j}^{(2)}+\frac{2}{3} \Delta t L_{h}\left(u^{(2)}\right)_{j} .
\end{align*}
$$

In particular, [2] pointed out that the boundary conditions in the intermediate stages of the above RK scheme should be modified as follows to avoid order reduction:

$$
\begin{align*}
& u^{n} \sim g_{l}\left(t_{n}\right), \\
& u^{(1)} \sim g_{l}\left(t_{n}\right)+\Delta t g_{l}^{\prime}\left(t_{n}\right),  \tag{2.5}\\
& u^{(2)} \sim g_{l}\left(t_{n}\right)+\frac{1}{2} \Delta t g_{l}^{\prime}\left(t_{n}\right)+\frac{1}{4} \Delta t^{2} g_{l}^{\prime \prime}\left(t_{n}\right) .
\end{align*}
$$

Note that for a high order finite difference scheme, a wide computational stencil is generally required. Hence, it is inevitable that some points in the computational stencil are not in our computational domain,

$$
x_{-p}=x_{0}-p \Delta x, \quad x_{N+p}=x_{N}+p \Delta x, \quad p=1,2, \cdots .
$$

Therefore, we can regard the boundary treatment problem as construction of the ghost point values. In the following, we will first review the original (S)ILW method proposed by Tan et al. $[28,31]$. And then, a new SILW method will be proposed to improve the computational efficiency while maintaining accuracy and stability. Linear stability analysis will be given to demonstrate the advantage of the new proposed method.
2.1. Review of the original SILW method. The main idea of the original inverse Lax-Wendroff method for hyperbolic conservation law equations [28] is to convert the spatial derivatives into the time derivatives through the PDE and boundary conditions at the inflow boundary. At the outflow boundary, the spatial derivatives of each order are approximated by extrapolation. After that, the values of the ghost points outside the computational domain are obtained by Taylor expansion at the boundary. More specifically, the ghost points near outflow boundaries, such as the right boundary $x=1$ in our example problem (2.1), can be obtained by extrapolation directly. We can choose the traditional Lagrange extrapolation with appropriate accuracy when the solution is smooth near the boundary, or least square extrapolation / WENO type extrapolation $[19,28,31]$ when the solution contains discontinuities near the boundary.

For the inflow boundary, such as the left boundary $x=-1$ in our example problem (2.1), to ensure our boundary treatment has $d$-th order accuracy, the value of the ghost points near $x=-1$ will be obtained by Taylor expansion:

$$
\begin{equation*}
u_{j}=\sum_{k=0}^{d-1} \frac{\left(x_{j}+1\right)^{k}}{k!} u^{*(k)}, \quad j=-1,-2, \cdots \tag{2.6}
\end{equation*}
$$

where, $u^{*(k)}$ is the approximation of $\left.\partial_{x}^{(k)} u\right|_{x=-1}$ with at least $(d-k)$-th order accuracy. Using PDE and boundary condition repeatedly, we have that

$$
\begin{aligned}
\left.\partial_{x}^{(0)} u\right|_{x=-1} & =g_{l}(t) \\
\left.\partial_{x}^{(1)} u\right|_{x=-1} & =\frac{g_{l}^{\prime}(t)}{-f\left(g_{l}(t)\right)}, \\
\left.\partial_{x}^{(2)} u\right|_{x=-1} & =\frac{f^{\prime}\left(g_{l}(t)\right) g_{l}^{\prime \prime}(t)-2 f^{\prime \prime}\left(g_{l}(t)\right) g_{l}^{\prime}(t)^{2}}{f^{\prime}\left(g_{l}(t)\right)^{3}},
\end{aligned}
$$

...

Table 1. The table of $\left(k_{d}\right)_{\min }$ for original SILW method.

| $d$ | 3 | 5 | 7 | 9 | 11 | 13 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\left(k_{d}\right)_{\min }$ | 2 | 3 | 4 | 6 | 8 | 10 |

Thus, we can set

$$
u^{*(k)}=\left.\partial_{x}^{(k)} u\right|_{x=-1} .
$$

To avoid the very heavy algebra of above original ILW method when calculating the high order space derivatives, the simplified ILW (SILW) method was proposed in [31]. Specifically, $u^{*(0)}, u^{*(1)}$ are constructed by the original ILW procedure, i.e., converting the spatial derivatives into the time derivatives through the PDE and boundary condition. The higher order spatial derivatives $u^{*(k)}, 2 \leq k \leq d-1$, are extrapolated from the interior points directly. This method can greatly improve the computational efficiency, especially for high-dimensional systems. However, [14] analyzed the linear stability of the SILW method through the eigenvalue method, showing that the SILW method [31] is stable for any $C_{a} \in[0,1)$ only when $d=3$, but unstable for $d>3$. In order to guarantee the stability, more ILW procedure need to be used to construct higher order spatial derivatives at the boundary.

Suppose that for a $d$-th order scheme, $u^{*(k)}$ is obtained through ILW procedure if $k \leq k_{d}-1$, or by extrapolation if $k_{d} \leq k \leq d-1$. For different high order schemes, [14] used the eigenvalue analysis to find out the minimum $k_{d}$, denoted by $\left(k_{d}\right)_{\text {min }}$, to make sure the scheme is stable for all $C_{a} \in[0,1)$. The values of $\left(k_{d}\right)_{\min }$ for a variety of $d$ are shown in Table 1. It can be seen that the $\left(k_{d}\right)_{\text {min }}$ is still large for high order scheme. This results in difficulty in writing codes and affects computational efficiency for high dimensional systems.

In summary, the above SILW method can be divided into the following two steps, i.e., "interpolation" and "extrapolation":

Step 1. Construct an interpolation polynomial $p(x)$ of degree $d-1$ with interior points $\left\{x_{0}, \cdots, x_{d-1}\right\}$, and obtain the approximation of spatial derivatives of each order on the boundary

$$
\left.u^{*(k)} \approx \partial_{x}^{k} p\right|_{x=-1}, \quad k=k_{d}, \cdots, d-1
$$

Step 2. Construct the extrapolation polynomial $q(x)$ of degree $d-1$ satisfying

$$
q^{(k)}(-1)=u^{*(k)}, \quad k=0, \cdots, d-1,
$$

where, $u^{*(k)}$ for $k<k_{d}$ are obtained by the ILW procedure (2.7), and the else are obtained by Step 1. Actually, in this case, $q(x)$ is the Taylor expansion polynomial. Then, we can get the ghost points values

$$
u_{j}=q\left(x_{j}\right), \quad j=-1,-2, \ldots .
$$

The above (S)ILW method is based on Taylor expansion. When assembling the extrapolation polynomial $q(x)$ in the second step, all information used is on the boundary. In fact, the information we can use to construct the extrapolation polynomial $q(x)$ mainly consists of two parts. One is the spatial derivatives on the boundary obtained by the ILW procedure, and the other part is the polynomial $p(x)$ constructed by interior points. In order to make the algorithm more efficient, we hope to use the first part information as less as possible under the premise of ensuring the stability of the scheme.
2.2. A new SILW method. In the following, we will describe our new SILW method for the scalar conservation law equation in (2.1), hoping $\left(k_{d}\right)_{\min }$ would be smaller for the same $d$. The key difference between the new method and the original one is that the extrapolation polynomial $q(x)$ will employ the point values on some special points in computational domain,

$$
u(-1+k \alpha \Delta x, t) \approx p(-1+k \alpha \Delta x), \quad k=1,2 \ldots
$$

instead of using the high order derivatives of the interpolation polynomial $p(x)$ at the boundary. Here, $\alpha \geq 0$ is a parameter to be determined such that the $\left(k_{d}\right)_{\min }$ would be smaller.

Specifically, we summarize the procedure of our new SILW method with $d$-th order accuracy in the following:
Step 1. Obtain the interpolation polynomial $p(x)$ of degree $d-1$ based on the points $\left\{x_{0}, \cdots, x_{d-1}\right\}$. Let

$$
u_{k *}=p(-1+k \alpha \Delta x), \quad 1 \leq k \leq d-k_{d} .
$$

Step 2. Construct the extrapolation polynomial $q(x)$ of degree $d-1$ to satisfy the following conditions:

$$
\begin{aligned}
& q^{(k)}(-1)=\left.\partial_{x}^{(k)} u\right|_{x=-1}, \quad 0 \leq k \leq k_{d}-1, \\
& q(-1+k \alpha \Delta x)=u_{k *}, \quad 1 \leq k \leq d-k_{d}
\end{aligned}
$$

where, $\left.\partial_{x}^{(k)} u\right|_{x=-1}$ is obtained by the ILW procedure. Let the ghost point values be the values of the extrapolation polynomial $q(x)$ at the corresponding points:

$$
u_{j}=q\left(x_{j}\right) \quad j=-1,-2, \ldots
$$

In the next subsection, we will show that through adjusting the value of $\alpha$, our new SILW method is better than the original SILW method in computational efficiency and stability.
2.3. Linear stability analysis. Here, we will give the stability analysis of the fully discrete schemes using the eigenvalue spectrum visualization. We consider the case of $d=2 k-1(k=2,3,4,5,6,7)$ and assume that $f^{\prime}(u)>0$. The conservative linear upwind scheme is used for spatial discretization. That is, $L_{h}$ in the scheme (2.3) is in the following form: $d=3$ :

$$
L_{h}(u)_{j}=-\frac{1}{\Delta x}\left(\frac{1}{6} f_{j-2}-f_{j-1}+\frac{1}{2} f_{j}+\frac{1}{3} f_{j+1}\right),
$$

$d=5:$

$$
L_{h}(u)_{j}=-\frac{1}{\Delta x}\left(-\frac{1}{30} f_{j-3}+\frac{1}{4} f_{j-2}-f_{j-1}+\frac{1}{3} f_{j}+\frac{1}{2} f_{j+1}-\frac{1}{20} f_{j+2}\right)
$$

$d=7:$

$$
\begin{aligned}
L_{h}(u)_{j}=-\frac{1}{\Delta x} & \left(\frac{1}{140} f_{j-4}-\frac{7}{105} f_{j-3}+\frac{3}{10} f_{j-2}-f_{j-1}+\frac{1}{4} f_{j}+\frac{3}{5} f_{j+1}\right. \\
& \left.-\frac{1}{10} f_{j+2}+\frac{1}{105} f_{j+3}\right)
\end{aligned}
$$

$$
d=9:
$$

$$
\begin{aligned}
L_{h}(u)_{j}=-\frac{1}{\Delta x} & \left(-\frac{1}{630} f_{j-5}+\frac{1}{56} f_{j-4}-\frac{2}{21} f_{j-3}+\frac{1}{3} f_{j-2}-f_{j-1}+\frac{1}{5} f_{j}\right. \\
& \left.+\frac{2}{3} f_{j+1}-\frac{1}{7} f_{j+2}+\frac{1}{42} f_{j+3}-\frac{1}{504} f_{j+4}\right)
\end{aligned}
$$

$d=11:$

$$
\begin{aligned}
L_{h}(u)_{j}=-\frac{1}{\Delta x} & \left(\frac{1}{2772} f_{j-6}-\frac{1}{210} f_{j-5}+\frac{5}{168} f_{j-4}-\frac{5}{42} f_{j-3}+\frac{5}{14} f_{j-2}\right. \\
& -f_{j-1}+\frac{1}{6} f_{j}+\frac{5}{7} f_{j+1}-\frac{5}{28} f_{j+2}+\frac{5}{126} f_{j+3}-\frac{1}{168} f_{j+4} \\
& \left.+\frac{1}{2310} f_{j+5}\right)
\end{aligned}
$$

$d=13:$

$$
\begin{aligned}
L_{h}(u)_{j}=-\frac{1}{\Delta x} & \left(-\frac{1}{12012} f_{j-7}+\frac{1}{792} f_{j-6}-\frac{1}{110} f_{j-5}+\frac{1}{24} f_{j-4}-\frac{5}{36} f_{j-3}\right. \\
& +\frac{3}{8} f_{j-2}-f_{j-1}+\frac{1}{7} f_{j}+\frac{3}{4} f_{j+1}-\frac{5}{24} f_{j+2}+\frac{1}{18} f_{j+3} \\
& \left.-\frac{1}{188} f_{j+4}+\frac{1}{660} f_{j+5}-\frac{1}{10296} f_{j+6}\right) .
\end{aligned}
$$

In particular, for the linear case $f(u)=u$, the semi-discrete scheme (2.3) can be written in the matrix-vector form,

$$
\frac{d \boldsymbol{U}}{d t}=\frac{1}{\Delta x} \boldsymbol{Q} \boldsymbol{U}
$$

where, $\boldsymbol{U}=\left(u_{0}, u_{2}, \ldots, u_{N}\right)^{T}$ and $\boldsymbol{Q}$ is the coefficient matrix of the spatial discretization.

References $[14,32]$ pointed out that we only need to care about the fixed eigenvalues of the matrix $\boldsymbol{Q}$ with the increase of grid points for stability analysis. If we use the third-order TVD RK time discretization (2.4), the stability region can be expressed as

$$
\begin{equation*}
|z(\mu)| \leq 1, \quad z(\mu)=1+\mu+\frac{\mu^{2}}{2}+\frac{\mu^{2}}{6} \tag{2.8}
\end{equation*}
$$

where, $\mu=s \frac{\Delta t}{\Delta x}$, and $s$ is the fixed eigenvalue of $\boldsymbol{Q}$ we are concerned about. Notice that $z$ may not exist or there may be more than one. If there is more than one $z$, we consider the largest $|z(\mu)|$. Let the CFL number be $\lambda_{c f l}=\frac{\Delta t}{\Delta x}$. We hope that the CFL number is independent of $C_{a}$. Hence, we discuss stability on the premise of maximum CFL number $\left(\lambda_{c f l}\right)_{\max }$, where, $\left(\lambda_{c f l}\right)_{\max }$ is the maximum CFL number for the corresponding Cauchy problem, and their specific values are shown in the Table 2.

We select several groups of different $\alpha$ and $k_{d}$ for linear stability analysis for different schemes, and the range of $\alpha$ is given as $[0,10]$. We compute the largest $|z(\mu)|$ for all the eigenvalues $s$. By using the software Matlab, we show the max $|z(\mu)|$ for $C_{a} \in[0,1)$ with different $\alpha$. For instance, we get the results of the third order

Table 2. The maximum CFL number for Cauchy problem.

| $d$ | 3 | 5 | 7 | 9 | 11 | 13 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\left(\lambda_{c f l}\right)_{\max }$ | 1.62 | 1.43 | 1.24 | 1.12 | 1.04 | 0.99 |

scheme with the new SILW procedure with $k_{d}=2$ as in Figure 1. When $\alpha=0.60$, $\max |z(\mu)|>1$ when $C_{a}$ approaches 0 . However, $\max |z(\mu)| \leq 1$ for all $C_{a} \in[0,1)$ if $\alpha=0.61$. This indicates that we should take $\alpha \geq 0.61$ to guarantee the scheme is stable with $k_{d}=2$. More cases are placed in Appendix A. A numerical test is also given to verify the stability analysis results. Finally, the minimum $k_{d}$ and the corresponding appropriate $\alpha$ range are shown in Table 3 .


Fig. 1. The third order scheme with the new SILW procedure with $k_{d}=2$. The horizontal axis represents $C_{a}$ and the vertical axis represents the largest $|z(\mu)|$.

As can be seen from the Table 3, compared with the original SILW method, we can construct the stable boundary treatments with smaller $k_{d}$ by adjusting $\alpha$. Here, we plot $C_{a}$ versus max $|z(\mu)|$ for both boundary treatments with the same order of accuracy and the same CFL number to compare the original and the new SILW methods. The fifth order schemes with $k_{d}=2$ are taken as an example, see Figure 2. It is observed that with the original SILW method, the fifth order upwind scheme is unstable when the first grid point is either close to or far from the left boundary, while the new SILW method is stable for all the $C_{a} \in[0,1]$.

Table 3. Linear stability analysis results of the new SILW method

| $d$ | 3 | 5 | 7 | 9 | 11 | 13 |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\left(k_{d}\right)_{\min }$ | 2 | 2 | 2 | 3 | 3 | 4 |
| $\alpha$ | $[0.61,10]$ | $[0.92,5.11]$ | $[1.34,1.99]$ | $[1.29,2.43]$ | $[1.42,1.70]$ | $[1.49,2.08]$ |

## 3. The new SILW method for one- and two-dimensional systems.

3.1. The new SILW method for one-dimensional Euler equation. We consider the following one-dimensional compressible Euler equation:

$$
\begin{equation*}
\boldsymbol{U}_{t}+\boldsymbol{F}(\boldsymbol{U})_{x}=\mathbf{0}, \quad x \in(-1,1), t>0 \tag{3.1}
\end{equation*}
$$



Fig. 2. The stability analysis results of the fifth order scheme with $k_{d}=2$, the horizontal axis represents $C_{a}$ and the vertical axis represents the largest $|z(\mu)|$.
where

$$
\boldsymbol{U}=\left(\begin{array}{c}
U_{1} \\
U_{2} \\
U_{3}
\end{array}\right)=\left(\begin{array}{c}
\rho \\
\rho u \\
E
\end{array}\right), \quad \boldsymbol{F}(\boldsymbol{U})=\left(\begin{array}{c}
\rho u \\
\rho u^{2}+p \\
u(E+p)
\end{array}\right)
$$

Here, $\rho, u, p$ and $E$ represent the density, velocity, pressure and total energy per volume, respectively. In order to close the system, we give the following equation of state of ideal gas:

$$
E=\frac{p}{\gamma-1}+\frac{1}{2} \rho u^{2}
$$

Here, $\gamma$ is the adiabatic constant, which equals to 1.4 for an ideal polytropic gas.
We consider the boundary treatment of left boundary $x=-1$ as an example. The original Euler equation (3.1) can be rewritten into the following nonconservative form:

$$
\boldsymbol{U}_{t}+\boldsymbol{A}(\boldsymbol{U}) \boldsymbol{U}_{x}=\mathbf{0}
$$

where, $\boldsymbol{A}(\boldsymbol{U})=\boldsymbol{F}^{\prime}(\boldsymbol{U})$ is the Jacobi matrix,

$$
\boldsymbol{A}(\boldsymbol{U})=\left(\begin{array}{ccc}
0 & 1 & 0  \tag{3.2}\\
\frac{1}{2}(\gamma-3) u^{2} & (3-\gamma) u & \gamma-1 \\
\frac{1}{2}(\gamma-1) u^{3}-u H & H-(\gamma-1) u^{2} & \gamma u
\end{array}\right)=\left(\begin{array}{l}
\boldsymbol{a}_{1}(\boldsymbol{U}) \\
\boldsymbol{a}_{2}(\boldsymbol{U}) \\
\boldsymbol{a}_{3}(\boldsymbol{U})
\end{array}\right)
$$

with $H=(E+p) / \rho$. And the Jacobi matrix is diagonalizable:

$$
\boldsymbol{A}(\boldsymbol{U})=\boldsymbol{F}^{\prime}(\boldsymbol{U})=\boldsymbol{R} \boldsymbol{\Lambda} \boldsymbol{L}
$$

Here, $\boldsymbol{\Lambda}=\operatorname{diag}(u-c, u, u+c), c=\sqrt{\gamma p / \rho}$ is the speed of sound, $\boldsymbol{R}$ and $\boldsymbol{L}$ are matrices as follows:

$$
\boldsymbol{R}(\boldsymbol{U})=\left(\begin{array}{ccc}
1 & 1 & 1 \\
u-c & u & u+c \\
H-u c & \frac{1}{2} u^{2} & H+u c
\end{array}\right)
$$

$\boldsymbol{L}(\boldsymbol{U})=\frac{1}{c^{2}}\left(\begin{array}{ccc}\frac{1}{2} u c+\frac{1}{4}(\gamma-1) u^{2} & -\frac{1}{2}(\gamma-1) u-\frac{1}{2} c & \frac{1}{2}(\gamma-1) \\ c^{2}-\frac{1}{2}(\gamma-1) u^{2} & (\gamma-1) u & 1-\gamma \\ -\frac{1}{2} u c+\frac{1}{4}(\gamma-1) u^{2} & -\frac{1}{2}(\gamma-1) u+\frac{1}{2} c & \frac{1}{2}(\gamma-1)\end{array}\right)=\left(\begin{array}{l}\boldsymbol{l}_{1}(\boldsymbol{U}) \\ \boldsymbol{l}_{2}(\boldsymbol{U}) \\ \boldsymbol{l}_{3}(\boldsymbol{U})\end{array}\right)$.
The number of boundary conditions we need to give is determined by the sign of the eigenvalues of Jacobi matrix $\boldsymbol{A}(\boldsymbol{U})$ at the boundary. Specifically, on the left boundary $x=-1$, it can be divided into the following cases:
Case 1. $u-c>0$ : three boundary conditions need to be given;
Case 2. $u-c \leq 0, u>0$ : two boundary conditions need to be given;
Case 3. $u \leq 0, u+c>0$ : only one boundary condition need to be given;
Case 4. $u+c \leq 0$ : no boundary conditions are required.
In particular, for case 3 , when the given boundary condition is $U_{2}(-1, t)=0$, we get the free slip solid wall boundary condition.

Here, we take the case 2 as an example to describe our algorithm. Suppose two boundary conditions are given at the left boundary,

$$
U_{1}(-1, t)=g_{1}(t), \quad \text { and } \quad U_{2}(-1, t)=g_{2}(t)
$$

Again, we perform uniform mesh generation on the computational domain with mesh size $\Delta x$,

$$
-1+C_{a} \Delta x=x_{0}<\cdots<x_{N}=1-C_{b} \Delta x .
$$

Employ the finite difference methods to get the semi-discrete scheme:

$$
\begin{equation*}
\frac{d}{d t} \boldsymbol{U}_{j}=-\frac{1}{\Delta x}\left(\hat{\boldsymbol{F}}_{j+1 / 2}-\hat{\boldsymbol{F}}_{j-1 / 2}\right), \quad j=0, \cdots, N \tag{3.3}
\end{equation*}
$$

Here, $\boldsymbol{U}_{j}(t)$ is the approximation of $\boldsymbol{U}\left(x_{j}, t\right)$, and numerical flux $\hat{\boldsymbol{F}}_{j+1 / 2}$ can be obtained by the WENO reconstruction. We take the fifth order scheme as an example to describe our boundary algorithm, and other high order schemes can be obtained similarly. For the fifth order WENO scheme, we need the values of at three ghost points near the boundary $x=-1$, which are $\boldsymbol{U}_{-1}, \boldsymbol{U}_{-2}$ and $\boldsymbol{U}_{-3}$.

Similar to the case of scalar equation, in order to obtain the ghost point values, we will construct the extrapolation polynomials $\boldsymbol{q}(x)$ near the boundary. It can be seen from Table 3 that, for the fifth order scheme, we have to use the point values and first order spatial derivatives on the boundary through the ILW procedure. To ensure the order of accuracy, the value $\boldsymbol{U}^{*(0)}$ and $\boldsymbol{U}^{*(1)}$ should be 5 th and 4 th order approximations of $\left.\boldsymbol{U}\right|_{x=-1}$ and $\left.\boldsymbol{U}_{x}\right|_{x=-1}$, respectively.

Specifically, we use the left characteristic matrix $\boldsymbol{L}=\boldsymbol{L}\left(\boldsymbol{U}_{0}\right)$ to do the characteristic projection $\boldsymbol{V}=\boldsymbol{L} \boldsymbol{U}$. Here,

$$
\boldsymbol{V}=\left(V_{1}, V_{2}, V_{3}\right)^{T}
$$

is the characteristic variable. In case $2, V_{1}$ is the outflow variable, $V_{2}$ and $V_{3}$ are the inflow variables. Combined with the boundary conditions, we can obtain the following linear system:

$$
\begin{align*}
U_{1}^{*(0)} & =g_{1}(t), \\
U_{2}^{*(0)} & =g_{2}(t),  \tag{3.4}\\
\boldsymbol{l}_{1} \cdot \boldsymbol{U}^{*(0)} & =V_{1}^{*(0)},
\end{align*}
$$

where, $V_{1}^{*(0)}$ is the 5 th order approximation of $\left.V_{1}\right|_{x=-1}$ and can be extrapolated from the interior grid points. By solving the above system, we can get the value of $\boldsymbol{U}^{*(0)}$.

For $\boldsymbol{U}^{*(1)}$, applying the ILW procedure, we have

$$
\begin{align*}
& \boldsymbol{a}_{1}\left(\boldsymbol{U}^{*(0)}\right) \cdot \boldsymbol{U}^{*(1)}=-g_{1}^{\prime}(t), \\
& \boldsymbol{a}_{2}\left(\boldsymbol{U}^{*(0)}\right) \cdot \boldsymbol{U}^{*(1)}=-g_{2}^{\prime}(t), \tag{3.5a}
\end{align*}
$$

Combine with outflow conditions,

$$
\begin{equation*}
\boldsymbol{l}_{1} \cdot \boldsymbol{U}^{*(1)}=V_{1}^{*(1)} \tag{3.5b}
\end{equation*}
$$

where, $V_{1}^{*(1)}$ is the 4 th order approximation of $\left.\left(V_{1}\right)_{x}\right|_{x=-1}$ and can be extrapolated from the interior grid points. By combining and solving the above equations, we can get the value of $\boldsymbol{U}^{*(1)}$.

At this time, we can construct the point values of the ghost points through the new SILW method. The specific construction method is as follows:
Step 1. Construct the vector of interpolation polynomial $\boldsymbol{p}(x)$ of degree four with the interior grid points values $\left\{\boldsymbol{U}_{0}, \cdots, \boldsymbol{U}_{4}\right\}$. Let

$$
\boldsymbol{U}_{k *}=\boldsymbol{p}(-1+k \alpha \Delta x), \quad 1 \leq k \leq 3
$$

Here, $\alpha$ can be selected as any number in $[0.92,5.11]$ according to the stability analysis results in Table 3.
Step 2. Construct the extrapolation polynomial $\boldsymbol{q}(x)$ of degree four to satisfy:

$$
\begin{aligned}
& \boldsymbol{q}(-1)=\boldsymbol{U}^{*(0)} \\
& \boldsymbol{q}^{\prime}(-1)=\boldsymbol{U}^{*(1)} \\
& \boldsymbol{q}(-1+k \alpha \Delta x)=\boldsymbol{U}_{k *}, \quad 1 \leq k \leq 3
\end{aligned}
$$

Step 3. Let the ghost points values be the values of the extrapolation polynomial $\boldsymbol{q}(x)$ at the corresponding points:

$$
\boldsymbol{U}_{j}=\boldsymbol{q}\left(x_{j}\right) \quad j=-1,-2,-3 .
$$

In particular, if the solution of the equation has discontinuities near the boundary, we can apply the WENO extrapolation technique when constructing the extrapolation polynomial $\boldsymbol{q}(x)$ in step 3 . More details about WENO extrapolation can be found in $[19,28,31]$. In addition, in many cases, boundary conditions are not directly given to the conserved variables. For example, the temperature or pressure is given on the boundary. For these cases, when dealing with the boundary conditions, we need to convert the conservation equation into equations in terms of primitive variables. This process is used in $[3,16,17,30]$, which is very similar to the above process of conserved variables. We will not expand it here.
3.2. The new SILW method for two-dimensional Euler equation. Consider the two-dimensional Euler equation as follows:

$$
\begin{equation*}
\frac{\partial \boldsymbol{U}}{\partial t}+\frac{\partial \boldsymbol{F}(\boldsymbol{U})}{\partial x}+\frac{\partial \boldsymbol{G}(\boldsymbol{U})}{\partial y}=\mathbf{0}, \quad(x, y)^{T} \in \Omega \tag{3.6}
\end{equation*}
$$

where,

$$
\boldsymbol{U}=\left(\begin{array}{c}
\rho  \tag{3.7}\\
\rho u \\
\rho v \\
E
\end{array}\right), \quad \boldsymbol{F}(\boldsymbol{U})=\left(\begin{array}{c}
\rho u \\
\rho u^{2}+p \\
\rho u v \\
u(E+p)
\end{array}\right), \quad \boldsymbol{G}(\boldsymbol{U})=\left(\begin{array}{c}
\rho v \\
\rho u v \\
\rho v^{2}+p \\
v(E+p)
\end{array}\right)
$$

Here, $\rho, \boldsymbol{u}=(u, v)^{T}, p$ and $E$ represent the density, velocity, pressure and total energy per volume, respectively. In order to close the system, we give the following equation of state of ideal gas:

$$
E=\frac{p}{\gamma-1}+\frac{1}{2} \rho\left(u^{2}+v^{2}\right)
$$

Here, $\gamma$ is the adiabatic constant, which equals to 1.4 for an ideal polytropic gas.
We use a uniform non body-fitted Cartesian mesh to divide the domain

$$
x_{i+1}=x_{i}+\Delta x, \quad y_{j+1}=y_{j}+\Delta y
$$

with mesh size $\Delta x$ and $\Delta y$ in $x$ - and $y$-direction, respectively. Discretize the equation into the following conservative semi-discrete scheme:

$$
\frac{d \boldsymbol{U}_{i, j}}{d t}+\frac{\hat{\boldsymbol{F}}_{i+\frac{1}{2}, j}-\hat{\boldsymbol{F}}_{i-\frac{1}{2}, j}}{\Delta x}+\frac{\hat{\boldsymbol{G}}_{i, j+\frac{1}{2}}-\hat{\boldsymbol{G}}_{i, j-\frac{1}{2}}}{\Delta y}=\mathbf{0}
$$

where, $\boldsymbol{U}_{i, j}(t)$ is approximation to the exact solution $\boldsymbol{U}\left(x_{i}, y_{j}, t\right), \hat{\boldsymbol{F}}_{i+\frac{1}{2}, j}$ and $\hat{\boldsymbol{G}}_{i, j+\frac{1}{2}}$ are numerical fluxes, which can be obtained by the 5 th order WENO reconstruction.


Fig. 3. The local coordinate rotation diagram.

Suppose $P_{i j}=\left(x_{i}, y_{j}\right)$ is a ghost point near the boundary. At first, we find its foot point $P_{a} \in \partial \Omega\left(t_{n}\right)$, so that the normal $\mathbf{n}$ at $P_{a}$ goes through $P_{i j}$, as shown in Figure 3. Assume the normal vector from $P_{a}$ to $P_{i, j}$ is $\mathbf{n}=(\cos \theta, \sin \theta)^{T}$. In order to simplify the algorithm, we perform a local coordinate rotation transformation at $P_{a}$,

$$
\binom{\hat{x}}{\hat{y}}=\left(\begin{array}{cc}
\cos \theta & \sin \theta \\
-\sin \theta & \cos \theta
\end{array}\right)\binom{x}{y} .
$$

In the new coordinate system, the equation (3.6) can be rewritten as

$$
\begin{equation*}
\frac{\partial \hat{\boldsymbol{U}}}{\partial t}+\frac{\partial \boldsymbol{F}(\hat{\boldsymbol{U}})}{\partial \hat{x}}+\frac{\partial \boldsymbol{G}(\hat{\boldsymbol{U}})}{\partial \hat{y}}=\mathbf{0} \tag{3.8}
\end{equation*}
$$

where,

$$
\hat{\boldsymbol{U}}=\left(\begin{array}{c}
\rho \\
\rho \hat{u} \\
\rho \hat{v} \\
E
\end{array}\right)=\left(\begin{array}{c}
\hat{U}_{1} \\
\hat{U}_{2} \\
\hat{U}_{3} \\
\hat{U}_{4}
\end{array}\right), \quad\binom{\hat{u}}{\hat{v}}=\left(\begin{array}{cc}
\cos \theta & \sin \theta \\
-\sin \theta & \cos \theta
\end{array}\right)\binom{u}{v} .
$$

$$
\boldsymbol{A}(\hat{\boldsymbol{U}})=\left(\begin{array}{l}
\boldsymbol{a}_{1}(\hat{\boldsymbol{U}}) \\
\boldsymbol{a}_{2}(\hat{\boldsymbol{U}}) \\
\boldsymbol{a}_{3}(\hat{\boldsymbol{U}}) \\
\boldsymbol{a}_{4}(\hat{\boldsymbol{U}})
\end{array}\right)=\boldsymbol{F}^{\prime}(\hat{\boldsymbol{U}}), \quad \text { Res }=-\frac{\partial \boldsymbol{G}(\hat{\boldsymbol{U}})}{\partial \hat{y}}=\left(\begin{array}{l}
R_{e s_{1}} \\
R_{2} \\
R e s_{2} \\
R_{2}
\end{array}\right) .
$$

Then, the equations can be written in the following non conservative form

$$
\begin{equation*}
\hat{\boldsymbol{U}}_{t}+\boldsymbol{A}(\hat{\boldsymbol{U}}) \hat{\boldsymbol{U}}_{x}=\boldsymbol{R e s} \tag{3.9}
\end{equation*}
$$

The original equation is hyperbolic, so $\boldsymbol{A}(\hat{\boldsymbol{U}})$ is diagonalizable:

$$
\boldsymbol{A}(\hat{\boldsymbol{U}})=\boldsymbol{R}(\hat{\boldsymbol{U}}) \boldsymbol{\Lambda}(\hat{\boldsymbol{U}}) \boldsymbol{L}(\hat{\boldsymbol{U}})
$$

Here,

$$
\boldsymbol{\Lambda}(\hat{\boldsymbol{U}})=\operatorname{diag}(\hat{u}-c, \hat{u}, \hat{u}, \hat{u}+c)
$$

$$
\boldsymbol{L}(\hat{\boldsymbol{U}})=\left(\begin{array}{c}
\boldsymbol{l}_{1}(\hat{\boldsymbol{U}}) \\
\boldsymbol{l}_{2}(\hat{\boldsymbol{U}}) \\
\boldsymbol{l}_{3}(\hat{\boldsymbol{U}}) \\
\boldsymbol{l}_{4}(\hat{\boldsymbol{U}})
\end{array}\right) .
$$

The number of boundary conditions that should be given at the boundary point $P_{a}$ is related to the eigenvalues $\hat{u}-c, \hat{u}, \hat{u}, \hat{u}+c$ at this point. Specifically, it can be divided into the following situations:
Case 1: $\hat{u}-c>0$, no boundary conditions are required;
Case 2: $\hat{u}-c \leq 0, \quad \hat{u}>0$, only one boundary condition needs to be given;
Case 3: $\hat{u} \leq 0, \quad \hat{u}+c>0$, three boundary conditions need to be given;
Case 4: $\hat{u}+c \leq 0$, four boundary conditions need to be given.
We take case 2 as an example to describe our algorithm. Suppose the boundary condition given at the boundary point $P_{a}$ is

$$
\hat{U}_{2}=g(t)
$$

As before, we can transform the construction problem of ghost point values into the construction problem of extrapolation polynomial $\boldsymbol{q}(s)$. It also can be seen from Table 3 that, for the fifth order scheme, we need to use the 0th and 1st order normal direction derivatives on the boundary obtained from the ILW procedure when constructing $\boldsymbol{q}(s)$. That is, we need to get the value of $\hat{\boldsymbol{U}}^{*(0)}$ and $\hat{\boldsymbol{U}}^{*(1)}$, which are the 5 th and 4th order approximations of $\left.\hat{\boldsymbol{U}}\right|_{P_{a}}$ and $\left.\hat{\boldsymbol{U}}_{\hat{x}}\right|_{P_{a}}$ respectively, through the ILW procedure.

Specifically, we use the left characteristic matrix $\boldsymbol{L}=\boldsymbol{L}\left(\hat{\boldsymbol{U}}_{\text {ext }}\right)$ to do the characteristic projection $\boldsymbol{V}=\boldsymbol{L} \hat{\boldsymbol{U}}$. Here, $\hat{\boldsymbol{U}}_{\text {ext }}$ is the extrapolation value at $P_{a}, \boldsymbol{V}=$ $\left(V_{1}, V_{2}, V_{3}, V_{4}\right)^{T}$ is the characteristic variable. For case $2, V_{2}, V_{3}, V_{4}$ are the outflow variables, $V_{1}$ is the inflow variable.

Combined with the boundary conditions, we can obtain the following linear system:

$$
\begin{align*}
\hat{U}_{2}^{*(0)} & =g(t), \\
\boldsymbol{l}_{2} \cdot \hat{\boldsymbol{U}}^{*(0)} & =V_{2}^{*(0)},  \tag{3.10}\\
\boldsymbol{l}_{3} \cdot \hat{\boldsymbol{U}}^{*(0)} & =V_{3}^{*(0)}, \\
\boldsymbol{l}_{4} \cdot \hat{\boldsymbol{U}}^{*(0)} & =V_{4}^{*(0)},
\end{align*}
$$

where, $V_{2}^{*(0)}, V_{3}^{*(0)}$ and $V_{4}^{*(0)}$ can be extrapolated from the interior grid points. By solving the above system, we can get the value of $\hat{\boldsymbol{U}}^{*(0)}$.

For $\hat{\boldsymbol{U}}^{*(1)}$, apply the ILW procedure and we have

$$
\begin{equation*}
\boldsymbol{a}_{2}\left(\hat{\boldsymbol{U}}^{*(0)}\right) \cdot \hat{\boldsymbol{U}}^{*(1)}=-g^{\prime}(t)+\operatorname{Res}_{2} . \tag{3.11a}
\end{equation*}
$$

Combine with the outflow conditions,

$$
\begin{align*}
& \boldsymbol{l}_{2} \cdot \hat{\boldsymbol{U}}^{*(1)}=V_{1}^{*(2)}, \\
& \boldsymbol{l}_{3} \cdot \hat{\boldsymbol{U}}^{*(1)}=V_{1}^{*(3)},  \tag{3.11b}\\
& \boldsymbol{l}_{4} \cdot \hat{\boldsymbol{U}}^{*(1)}=V_{1}^{*(4)},
\end{align*}
$$

where, $V_{2}^{*(1)}, V_{3}^{*(1)}$ and $V_{4}^{*(1)}$ can be extrapolated from the interior grid points. By solving the above equations, we can get the value of $\hat{\boldsymbol{U}}^{*(1)}$.


Fig. 4. Two dimensional new SILW method diagram

Next, we use the new SILW method to construct the value of ghost point $P_{i, j}$. It is divided into the following steps:
Step 1. Obtain the interpolation polynomial $\boldsymbol{p}(x, y)$ of degree 4 with the values internal grid points near $P_{a}$.
Step 2. Let

$$
\boldsymbol{U}_{k *}=\boldsymbol{p}\left(P_{k}\right), \quad 1 \leq k \leq 3,
$$

where,

$$
P_{k}=P_{a}-k \alpha h \boldsymbol{n}, \quad 1 \leq k \leq 3
$$

As show in Figure 4, $\left.\left\{P_{k}\right\}\right|_{k=1} ^{3}$ are some non grid points in the interior area on the normal line. Here, $h=\sqrt{\Delta x^{2}+\Delta y^{2}}, \alpha$ can be chosen as any number in $\left[0.92 \frac{\max (\Delta x, \Delta y)}{h}, 5.11 \frac{\min (\Delta x, \Delta y)}{h}\right]$.
Step 3 . Let $\boldsymbol{q}(s)$ be the unique polynomial of degree 4 satisfy

$$
\begin{aligned}
& \boldsymbol{q}(0)=\boldsymbol{U}^{*(0)} \\
& \boldsymbol{q}^{\prime}(0)=\boldsymbol{U}^{*(1)} \\
& \boldsymbol{q}(-k \alpha h)=\boldsymbol{U}_{k *}, \quad 1 \leq k \leq 3 .
\end{aligned}
$$

Step 4. Take the function value at $P_{i j}$

$$
\boldsymbol{U}_{i j}=\boldsymbol{q}\left(\left|P_{i j}-P_{a}\right|\right)
$$

It should be noted that when $g(t)=0$, we actually get the non penetrating free slip boundary condition:

$$
\boldsymbol{u} \cdot \boldsymbol{n}=0
$$

As in the one-dimensional case, if the solution is discontinuous near the boundary, we can use a one-dimensional WENO extrapolation technique in step 3.

Notice that, for the problems with changing wind direction, the above inverse LaxWendroff procedure may involve solving an ill-conditioned linear algebraic system, which may ruin the accuracy or even lead to blowing up. There are two ways to deal with this problem. One is mentioned in [28], which adds additional extrapolation equations and solves a least squares problem whenever one of the eigenvalues very closed to 0 . The other method is proposed in [19], which evaluates the solution values and the flux values at ghost points separately. In the test examples of this paper, we use the first method. The performance of the second method applied to our new SILW boundary treatment is unclear.
4. Numerical tests. We take some numerical tests to show the efficiency and stability of our new proposed SILW method. We use the third and fifth order FDWENO scheme for spatial discretization. Correspondingly, the new SILW boundary treatment with third order and fifth order accuracy will be used, respectively. For all the one-dimensional numerical tests we take the parameter $\alpha=1.0$, for all the two-dimensional numerical tests we take $\alpha=1.25$. The third order TVD RK scheme (2.4) is employed for time discretization, with the time step

$$
\Delta t=\mathrm{CFL} \frac{\Delta x^{k / 3}}{c}
$$

for one-dimensional problems, and

$$
\Delta t=\frac{\mathrm{CFL}}{c_{x} / \Delta x^{k / 3}+c_{y} / \Delta y^{k / 3}} .
$$

for two-dimensional problems. Here, the index $k / 3$ help us to guarantee $k$-th order in time. $c=c_{x}=\max _{\boldsymbol{U}}\left|\lambda\left(\boldsymbol{F}^{\prime}(\boldsymbol{U})\right)\right|, c_{y}=\max _{\boldsymbol{U}}\left|\lambda\left(\boldsymbol{g}^{\prime}(\boldsymbol{u})\right)\right|$, and $\lambda$ is the eigenvalue of the Jacobian matrix. Throughout our numerical tests, the CFL number is taken as 0.6 .

Example 1. At first, we consider the accuracy test of the new SILW on the onedimensional Euler equation on the computational domain as $[-\pi, \pi]$. We choose suitable boundary conditions such that the exact solution is:

$$
\left\{\begin{array}{l}
\rho(x, t)=1-0.2 \sin (2 t-x)  \tag{4.1}\\
u(x, t)=2 \\
p(x, t)=2
\end{array}\right.
$$

In order to verify the applicability of our algorithm to the "cut cell" problem, we test with difference choices of $C_{a}$ and $C_{b}$. The computational errors about density $\rho$ at final time $t_{\text {end }}=1$ are shown in Table $4-5$. We can see that for all cases, the schemes can achieve the designed order accuracy with mesh refinements.

Table 4. Example 1: errors and orders of accuracy of $\rho$ with third order scheme.

| $N$ | $C_{a}=0.0001, C_{b}=0.7$ |  |  |  | $C_{a}=0.9999, C_{b}=0.7$ |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| 20 | $1.67 \mathrm{E}-004$ | - | $5.73 \mathrm{E}-004$ | - | $1.45 \mathrm{E}-004$ | - | $4.80 \mathrm{E}-004$ | - |
| 40 | $1.58 \mathrm{E}-005$ | 3.40 | $4.78 \mathrm{E}-005$ | 3.58 | $1.00 \mathrm{E}-005$ | 3.85 | $4.39 \mathrm{E}-005$ | 3.45 |
| 80 | $2.07 \mathrm{E}-006$ | 2.92 | $5.30 \mathrm{E}-006$ | 3.17 | $8.52 \mathrm{E}-007$ | 3.56 | $4.57 \mathrm{E}-006$ | 3.26 |
| 160 | $2.69 \mathrm{E}-007$ | 2.94 | $7.30 \mathrm{E}-007$ | 2.86 | $8.38 \mathrm{E}-008$ | 3.34 | $5.24 \mathrm{E}-007$ | 3.12 |
| 320 | $3.43 \mathrm{E}-008$ | 2.97 | $9.50 \mathrm{E}-008$ | 2.94 | $9.15 \mathrm{E}-009$ | 3.19 | $6.27 \mathrm{E}-008$ | 3.06 |
| 640 | $4.32 \mathrm{E}-009$ | 2.98 | $1.21 \mathrm{E}-008$ | 2.96 | $1.06 \mathrm{E}-009$ | 3.10 | $7.66 \mathrm{E}-009$ | 3.03 |

Table 5. Example 1: errors and orders of accuracy of $\rho$ with fifth order scheme.

| $N$ | $C_{a}=0.0001, C_{b}=0.7$ |  |  |  | $C_{a}=0.9999, C_{b}=0.7$ |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| 20 | $9.33 \mathrm{E}-005$ | - | $1.76 \mathrm{E}-004$ | - | $7.41 \mathrm{E}-005$ | - | $1.36 \mathrm{E}-004$ | - |
| 40 | $2.99 \mathrm{E}-006$ | 4.96 | $6.09 \mathrm{E}-006$ | 4.84 | $2.62 \mathrm{E}-006$ | 4.81 | $5.47 \mathrm{E}-006$ | 4.63 |
| 80 | $9.28 \mathrm{E}-008$ | 5.01 | $1.99 \mathrm{E}-007$ | 4.93 | $8.65 \mathrm{E}-008$ | 4.92 | $1.81 \mathrm{E}-007$ | 4.91 |
| 160 | $2.88 \mathrm{E}-009$ | 5.01 | $6.06 \mathrm{E}-009$ | 5.04 | $2.77 \mathrm{E}-009$ | 4.96 | $5.78 \mathrm{E}-009$ | 4.97 |
| 320 | $8.89 \mathrm{E}-011$ | 5.01 | $1.77 \mathrm{E}-010$ | 5.09 | $8.72 \mathrm{E}-011$ | 4.99 | $1.73 \mathrm{E}-010$ | 5.05 |

Example 2. Next, we consider the example given in [12] to test the accuracy of our method. The governing equations is still the one-dimensional compressible Euler equation, with following initial condition:

$$
\left\{\begin{array}{l}
\rho(x, 0)=\frac{1+0.2 \sin (x)}{2 \sqrt{3}}  \tag{4.2}\\
u(x, 0)=\sqrt{\gamma} \rho(x, 0) \\
p(x, 0)=\rho(x, 0)^{\gamma}
\end{array}\right.
$$

The computational domain is taken as $[0,2 \pi]$. We choose the parameter $\gamma=3$. Consequently, the exact solution is

$$
\rho(x, t)=\frac{\mu(x, t)}{2 \sqrt{3}}, \quad u(x, t)=\sqrt{\gamma} \rho(x, t), \quad p(x, t)=\rho(x, t)^{\gamma}
$$

where $\mu(x, t)$ is the solution of the following Burgers' equation:

$$
\left\{\begin{array}{l}
\mu_{t}+\left(\frac{\mu^{2}}{2}\right)_{x}=0, \quad 0<x<2 \pi, \quad t>0  \tag{4.3}\\
\mu(x, 0)=1+0.2 \sin (x), \quad 0 \leq x \leq 2 \pi
\end{array}\right.
$$

We take boundary conditions from the exact solution whenever needed.
We consider the extrema situation and set $C_{a}=0.0001, C_{b}=0.9999$. The computational errors about density $\rho$ and orders of accuracy at time $t_{\text {end }}=3.0$ are shown in Table 6 , indicating that our methods can achieve the designed third order or fifth order accuracy.

Example 3. Now we consider the interaction of two blast waves [28]. In this problem, multiple reflections occur between shock and rarefaction off the walls. The initial

Table 6. Example 2: errors and orders of accuracy of $\rho$.

| $N$ | third order scheme |  |  |  | fifth order scheme |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| 40 | $1.30 \mathrm{E}-003$ | - | $2.49 \mathrm{E}-003$ | - | $4.01 \mathrm{E}-004$ | - | $1.04 \mathrm{E}-003$ | - |
| 80 | $2.04 \mathrm{E}-004$ | 2.67 | $6.20 \mathrm{E}-004$ | 2.00 | $2.24 \mathrm{E}-005$ | 4.15 | $8.03 \mathrm{E}-005$ | 3.69 |
| 160 | $2.84 \mathrm{E}-005$ | 2.84 | $8.50 \mathrm{E}-005$ | 2.86 | $7.16 \mathrm{E}-007$ | 4.97 | $2.98 \mathrm{E}-006$ | 4.75 |
| 320 | $3.60 \mathrm{E}-006$ | 2.97 | $9.93 \mathrm{E}-006$ | 3.09 | $2.38 \mathrm{E}-008$ | 4.90 | $1.00 \mathrm{E}-007$ | 4.89 |
| 640 | $4.52 \mathrm{E}-007$ | 2.99 | $1.18 \mathrm{E}-006$ | 3.06 | $8.04 \mathrm{E}-010$ | 4.89 | $2.90 \mathrm{E}-009$ | 5.10 |



Fig. 5. Example 3: Density profiles. $h=1 / 640$. The solid line represents the reference solution and the circle represents the numerical solution.
condition is

$$
\boldsymbol{U}(x, 0)= \begin{cases}\boldsymbol{U}_{L}, & x<0.1  \tag{4.4}\\ \boldsymbol{U}_{M}, & 0.1<x<0.9 \\ \boldsymbol{U}_{R}, & x>0.9\end{cases}
$$

Here, $\rho_{L}=\rho_{M}=\rho_{R}=1, u_{L}=u_{M}=u_{R}=0, p_{L}=10^{3}, p_{M}=10^{-2}$, and $p_{R}=10^{2}$. We take $t_{\text {end }}=0.038$ and $C_{a}=0.0001, C_{b}=0.7$. At the same time, we use a very dense grid with $\Delta x=1 / 2560$ and the original ILW method to obtain the reference solution. The numerical results are shown in Figure 5. We can see that the new ILW method can distinguish the structure of the solution well, and higher order scheme has a better approximation to the complex structure.

Example 4. We consider two-dimensional linear scalar equations on a disk:

$$
\begin{equation*}
u_{t}+u_{x}+u_{y}=0, \quad(x, y)^{T} \in \Omega=\left\{(x, y): x^{2}+y^{2}<0.5\right\} . \tag{4.5}
\end{equation*}
$$

The initial condition is given as

$$
u(x, y, 0)=0.25+0.5 \sin [\pi(x+y)]
$$

and the boundary is given whenever needed such that the exact solution is

$$
u(x, y, t)=0.25+0.5 \sin [\pi(x+y-2 t)] .
$$

The domain is discretized by embedding the domain in a regular Cartesian mesh with $x_{i}=\left(i-\frac{1}{2}\right) \Delta x, y_{j}=\left(j-\frac{1}{2}\right) \Delta y$, and the non body-fitted Cartesian mesh $h=\Delta x=\Delta y$. We show Figure 6 as an example. The final time is taken as $t_{\text {end }}=1.0$. The numerical results are given in Table 7, indicating that our schemes are stable and can achieve the design order of accuracy.


Fig. 6. Example 4: Non body-fitted Cartesian mesh. The red points are the interior points.

Table 7. Example 4: errors and orders of accuracy .

|  | third order scheme |  |  |  | fifth order scheme |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $h$ | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| $1 / 10$ | $1.28 \mathrm{E}-004$ | - | $4.83 \mathrm{E}-004$ | - | $4.16 \mathrm{E}-004$ | - | $1.42 \mathrm{E}-003$ | - |
| $1 / 20$ | $1.33 \mathrm{E}-005$ | 3.26 | $4.54 \mathrm{E}-005$ | 3.41 | $1.51 \mathrm{E}-005$ | 4.77 | $1.38 \mathrm{E}-004$ | 3.36 |
| $1 / 40$ | $1.43 \mathrm{E}-006$ | 3.21 | $5.84 \mathrm{E}-006$ | 2.95 | $3.47 \mathrm{E}-007$ | 5.44 | $5.54 \mathrm{E}-006$ | 4.64 |
| $1 / 80$ | $1.46 \mathrm{E}-007$ | 3.28 | $6.68 \mathrm{E}-007$ | 3.12 | $1.17 \mathrm{E}-008$ | 4.89 | $2.71 \mathrm{E}-007$ | 4.35 |
| $1 / 160$ | $1.19 \mathrm{E}-008$ | 3.61 | $8.82 \mathrm{E}-008$ | 2.92 | $4.47 \mathrm{E}-010$ | 4.70 | $9.72 \mathrm{E}-009$ | 4.80 |

Example 5. We test the vortex evolution problem for the 2D Euler equation. The mean flow is $\rho=u=v=p=1$ with following isentropic vortex perturbation centered at $\left(x_{0}, y_{0}\right)=(0,0)$ (perturbation in $(u, v)$ and temperature $T=p / \rho$, no perturbation in the entropy $\left.S=p / \rho^{\gamma}\right)$ :

$$
\begin{align*}
(\delta u, \delta v) & =\frac{\epsilon}{2 \pi} e^{0.5\left(1-r^{2}\right)}(-\bar{y}, \bar{x}) \\
\delta T & =-\frac{(\gamma-1) \epsilon^{2}}{8 \gamma \pi^{2}} e^{\left(1-r^{2}\right)}  \tag{4.6}\\
\delta S & =0
\end{align*}
$$

where $(\bar{x}, \bar{y})=\left(x-x_{0}, y-y_{0}\right), r^{2}=\bar{x}^{2}+\bar{y}^{2}$, and the vortex strength $\epsilon=5$. It is clear that the exact solution is just the passive convection of the vortex with the mean velocity. The computational domain is taken as $(-0.5,1) \times(-0.5,1)$ and the final time is taken as $t_{\text {end }}=1.0$. The boundary conditions are taken from the exact
solution whenever needed. We divide the domain with the uniform Cartesian mesh $x_{i}=\left(i-\frac{1}{2}\right) h$ and $y_{j}=\left(j-\frac{1}{2}\right) h$, with mesh size $h=1.5 / N$. The numerical results in Table 8 show that the schemes are stable and can reach the designed high order.

Table 8. Example 5: errors and orders of accuracy of $\rho$.

|  | third order scheme |  |  |  | fifth order scheme |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $h$ | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| $3 / 40$ | $1.73 \mathrm{E}-004$ | - | $3.05 \mathrm{E}-004$ | - | $3.39 \mathrm{E}-005$ | - | $7.12 \mathrm{E}-005$ | - |
| $3 / 80$ | $2.17 \mathrm{E}-005$ | 2.99 | $4.10 \mathrm{E}-005$ | 2.89 | $1.09 \mathrm{E}-006$ | 4.95 | $2.33 \mathrm{E}-006$ | 4.93 |
| $3 / 160$ | $2.51 \mathrm{E}-006$ | 3.10 | $4.93 \mathrm{E}-006$ | 3.05 | $3.46 \mathrm{E}-008$ | 4.98 | $1.08 \mathrm{E}-007$ | 4.43 |
| $3 / 320$ | $2.99 \mathrm{E}-007$ | 3.07 | $6.32 \mathrm{E}-007$ | 2.96 | $1.12 \mathrm{E}-009$ | 4.94 | $4.71 \mathrm{E}-009$ | 4.51 |
| $3 / 640$ | $3.63 \mathrm{E}-008$ | 3.04 | $8.34 \mathrm{E}-008$ | 2.92 | $3.77 \mathrm{E}-011$ | 4.89 | $1.90 \mathrm{E}-010$ | 4.63 |

Example 6. Next, we consider 2D version of Example 2 [12]. The governing equation is the two-dimensional compressible Euler equations with following initial condition:

$$
\left\{\begin{array}{l}
\rho(x, y, 0)=\frac{1+0.2 \sin \left(\frac{x+y}{2}\right)}{\sqrt{6}}  \tag{4.7}\\
u(x, y, 0)=v(x, y, 0)=\sqrt{\frac{\gamma}{2}} \rho(x, y, 0) \\
p(x, y, 0)=\rho(x, y, 0)^{\gamma}
\end{array}\right.
$$

We choose the parameter $\gamma=3$, such that the exact solution is
$\rho(x, y, t)=\frac{\mu(x, y, t)}{\sqrt{6}}, \quad u(x, y, t)=v(x, y, t)=\sqrt{\frac{\gamma}{2}} \rho(x, y, t), \quad p(x, y, t)=\rho(x, y, t)^{\gamma}$,
where $\mu(x, y, t)$ is the solution of the following 2 D Burgers' equation:

$$
\left\{\begin{array}{l}
\mu_{t}+\left(\frac{\mu^{2}}{2}\right)_{x}+\left(\frac{\mu^{2}}{2}\right)_{y}=0, \quad(x, y) \in \Omega  \tag{4.8}\\
\mu(x, y, 0)=1+0.2 \sin \left(\frac{x+y}{2}\right)
\end{array}\right.
$$

We consider following two computational domains:

$$
\begin{align*}
& \Omega=[0,4 \pi] \times[0,4 \pi]  \tag{4.9a}\\
& \Omega=\left\{(x, y) \mid x^{2}+y^{2}<(1.5 \pi)^{2}\right\} \tag{4.9b}
\end{align*}
$$

and take boundary conditions from the exact solution whenever needed. For the square domain (4.9a), we use a grid similar to example 5. And for the circular domain (4.9b), we use a non body-fitted grid similar to example 4. The numerical results at final time $t_{\text {end }}=1$ are shown in Table 9-10. We can see that our schemes are stable and high order accuracy for all cases.

Example 7. We consider a flow around a cylinder. The center of the bottom surface of the cylinder is at the origin with a radius of 1 . At the initial moment, a fluid with Mach 3 moves towards the cylinder. In consideration of the symmetry, we only consider the problem of an upper half plane. For the lower boundary of the computation area at $y=0$, we use the reflection technique; for the left boundary

Table 9. Example 6: The errors and the orders of accuracy of $\rho$ on the square domain $\Omega=[0,4 \pi] \times[0,4 \pi]$.

| third order scheme | fifth order scheme |  |  |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
|  | $4.09 \mathrm{E}-004$ | - | $1.13 \mathrm{E}-005$ | - | $1.42 \mathrm{E}-005$ | - | $2.08 \mathrm{E}-006$ | - |
| $4 \pi / 150$ | $1.22 \mathrm{E}-004$ | 2.98 | $3.51 \mathrm{E}-006$ | 2.89 | $1.93 \mathrm{E}-006$ | 4.92 | $2.89 \mathrm{E}-007$ | 4.86 |
| $4 \pi / 200$ | $5.18 \mathrm{E}-005$ | 2.98 | $1.66 \mathrm{E}-006$ | 2.60 | $4.59 \mathrm{E}-007$ | 4.99 | $7.55 \mathrm{E}-008$ | 4.67 |
| $4 \pi / 250$ | $2.66 \mathrm{E}-005$ | 2.98 | $9.14 \mathrm{E}-007$ | 2.67 | $1.49 \mathrm{E}-007$ | 5.01 | $2.77 \mathrm{E}-008$ | 4.48 |
| $4 \pi / 300$ | $1.54 \mathrm{E}-005$ | 2.99 | $5.54 \mathrm{E}-007$ | 2.74 | $6.01 \mathrm{E}-008$ | 5.01 | $1.20 \mathrm{E}-008$ | 4.56 |

Table 10. Example 6: The errors and the orders of accuracy of $\rho$ on the circular domain $\Omega=\left\{(x, y) \mid x^{2}+y^{2}<(1.5 \pi)^{2}\right\}$.

|  | third order scheme |  |  |  | fifth order scheme |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $h$ | $L^{1}$ error | order | $L^{\infty}$ error | order | $L^{1}$ error | order | $L^{\infty}$ error | order |
| $4 \pi / 100$ | $1.46 \mathrm{E}-004$ | - | $1.14 \mathrm{E}-005$ | - | $1.48 \mathrm{E}-005$ | - | $8.59 \mathrm{E}-006$ | - |
| $4 \pi / 150$ | $4.20 \mathrm{E}-005$ | 3.08 | $4.28 \mathrm{E}-006$ | 2.42 | $2.13 \mathrm{E}-006$ | 4.78 | $1.28 \mathrm{E}-006$ | 4.69 |
| $4 \pi / 200$ | $1.81 \mathrm{E}-005$ | 2.91 | $1.79 \mathrm{E}-006$ | 3.03 | $6.65 \mathrm{E}-007$ | 4.05 | $3.51 \mathrm{E}-007$ | 4.49 |
| $4 \pi / 250$ | $9.20 \mathrm{E}-006$ | 3.05 | $8.90 \mathrm{E}-007$ | 3.12 | $2.21 \mathrm{E}-007$ | 4.91 | $1.16 \mathrm{E}-007$ | 4.95 |
| $4 \pi / 300$ | $5.44 \mathrm{E}-006$ | 2.87 | $5.23 \mathrm{E}-007$ | 2.91 | $8.42 \mathrm{E}-008$ | 5.31 | $4.72 \mathrm{E}-008$ | 4.93 |

of the computation region at $x=-4$, we give the inflow boundary condition; for the right boundary $x=0$ and the upper boundary $y=6$ of the computation area, the outflow boundary conditions are given. On the surface of a cylinder, our new ILW method is used to deal with a no-penetration boundary condition. As before, a uniform non body-fitted Cartesian mesh is used, which is shown in Figures 7. Figure 8 show the numerical results. We can see that the results are comparable with those in $[5,19,28,31]$.


Fig. 7. Example 7: The non body-fitted Cartesian mesh with near the cylinder boundary. The red points are the interior points.

Example 8. We consider the double Mach reflection problem. At the initial moment,


Fig. 8. Example 7: Pressure contour of flow past a cylinder, 20 contours from 2 to 15. $\Delta x=\Delta y=1 / 40$. left: third order scheme. right: fifth order scheme.
a horizontally moving Mach 10 shock wave passes through a wedge with an inclination angle of $30^{\circ}$. In common practice, the wedge is placed horizontally to apply reflective boundary conditions. At this time, the shock wave forms an angle of $60^{\circ}$ with the wall. In $[8,25]$, the original double Mach number reflection problem is computed respectively. With the ILW method, people can also do numerical simulation on the original region $[28,31]$. Here, we use the new ILW method to simulate this problem. In detail, at the top of the calculation area, we give the exact flow value according to the shock Mach number. at the left and right boundary, we give the supersonic inlet and outlet boundary conditions respectively. On the lower right boundary, a new ILW method is adopted. The discretization of space and time is consistent with the previous example. Figure 9 show the computational region and density contour respectively. The zoomed in region near the double Mach stem is presented in Figure 10. We rotate and translate the region for ease of comparison. It is observed that the new ILW method captures the shock wave well, and it is comparable with the previous results.
5. Concluding remarks. In this paper, we propose a new SILW method for conservation laws, which decomposes the procedure of construction ghost points into two steps: interpolation and extrapolation. At first, we approximate some special point values through an interpolation polynomial based on interior points near the boundary. Then, we construct a Hermite extrapolation polynomial based on the special point values and spatial derivatives at the boundary obtained through the ILW process. After that, we can get the approximation of the the ghost point values. Through the linear stability analysis with the eigenvalue method, we can conclude that our new SILW method is more efficient than the original SILW method while ensuring the stability. Then we extend our new SILW method to systems and high-dimensional


Fig. 9. Example 8: Left:The computational region of the double mach reflection problem. The dashed line indicates the computational domain used in [8, 25]. Right: The density contour. 30 contours from 1.731 to $20.92 . \Delta x=\Delta y=1 / 320$.


Fig. 10. Example 8: Density contour on the local area. 30 contours from 1.731 to 20.92 .
cases, and carry out a series of numerical experiments. The numerical results show that our new SILW method is stable and can achieve the expected accuracy. In the future, we are going to extend this new SILW method to deal with the initial-boundary value problems of diffusion equations and convection-diffusion equations.

## Appendix A. More results about linear stability analysis.

The linear stability analysis results of the new SILW method with different internal schemes and different $k_{d}$ are shown in Figure 11-13. The parameter $\alpha$ is taken as the critical value between stable and unstable. These figures verify the correctness and the optimality of the $\alpha$ range given in Table 3.

Next, we want to verify the results of the above stability analysis numerically.


Fig. 11. The result of linear stability analysis with $k_{d}=2$. The horizontal axis represents $C_{a}$ and the vertical axis represents the largest $|z(\mu)|$.

Consider the following problem:
(A.1)

$$
\left\{\begin{aligned}
u_{t}+u_{x} & =0, \quad-1<x<1, t>0 \\
u(x, 0) & =0.25+0.5 \sin (\pi x), \quad-1 \leq x \leq 1 \\
u(-1, t) & =0.25+0.5 \sin (\pi t), \quad t>0
\end{aligned}\right.
$$

The exact solution is

$$
u(x, t)=0.25+0.5 \sin (\pi(x-t))
$$

We use the $d$-th order upwind scheme for spatial discretization and the third-order


Fig. 12. The result of linear stability analysis with $k_{d}=3$. The horizontal axis represents $C_{a}$ and the vertical axis represents the largest $|z(\mu)|$.

TVD RK scheme for time discretization. Let $t_{\text {end }}=30, N=200$. Take time step

$$
\Delta t=\left(\lambda_{c f l}\right)_{\max } \Delta x
$$

We test the problem with $\alpha$ in or out the range given in Table 3. The numerical results are shown in Table 11. It can be observed that when $\alpha$ falls in the range, the scheme will be stable for all tested $C_{a}$. Otherwise, if $\alpha$ is out of the range, we can always find one $C_{a}$ such that the scheme is unstable.

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Fig. 13. The result of linear stability analysis with $k_{d}=4$. The horizontal axis represents $C_{a}$ and the vertical axis represents the largest $|z(\mu)|$.

Table 11. Numerical verification results of linear stability analysis

| d | Stable $\alpha$ in Table 3 | $\alpha$ | Result |
| :---: | :---: | :---: | :---: |
| 3 | [0.61,10] | $\begin{aligned} & \hline 0.60 \\ & 1.00 \end{aligned}$ | Unstable for $C_{a}=10^{-6}$ Stable for all tested $C_{a}$ |
| 5 | [0.92,5.11] | $\begin{aligned} & 0.91 \\ & 1.00 \\ & 5.12 \end{aligned}$ | Unstable for $C_{a}=0.38$ Stable for all tested $C_{a}$ Unstable for $C_{a}=0.70$ |
| 7 | [1.34,1.99] | $\begin{aligned} & \hline 1.33 \\ & 1.50 \\ & 2.00 \\ & \hline \end{aligned}$ | Unstable for $C_{a}=0.40$ Stable for all tested $C_{a}$ Unstable for $C_{a}=0.40$ |
| 9 | [1.29,2.43] | $\begin{aligned} & \hline 1.28 \\ & 1.50 \\ & 2.44 \\ & \hline \end{aligned}$ | Unstable for $C_{a}=0.85$ Stable for all tested $C_{a}$ Unstable for $C_{a}=0.03$ |
| 11 | [1.42,1.70] | $\begin{aligned} & \hline 1.41 \\ & 1.50 \\ & 1.71 \\ & \hline \end{aligned}$ | Unstable for $C_{a}=0.93$ Stable for all tested $C_{a}$ Unstable for $C_{a}=0.01$ |
| 13 | [1.49,2.08] | $\begin{aligned} & \hline 1.48 \\ & 1.75 \\ & 2.09 \\ & \hline \end{aligned}$ | Unstable for $C_{a}=1-10^{-6}$ <br> Stable for all tested $C_{a}$ <br> Unstable for $C_{a}=1-10^{-6}$ |

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