Robust high order discontinuous Galerkin schemes for two-dimensional gaseous detonations

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Abstract

One of the main challenges in computational simulations of gas detonation propagation is that negative density or negative pressure may emerge during the time evolution, which will cause blow-ups. Therefore, schemes with provable positivity-preserving of certain quantities such as density and pressure are desired. First order and second order positivity-preserving schemes were well studied, e.g., \cite{10}. A simple solution for arbitrarily high order schemes was proposed recently in \cite{22}. For high order discontinuous Galerkin (DG) method, even though the characteristicwise TVB limiter in \cite{1, 2} can kill oscillations, it is not sufficient to maintain the positivity. In this paper, we first show an extension of the \cite{22, 23, 24} to design positivity-preserving arbitrarily high order DG schemes for reactive Euler equations. Then we show a new simpler and more robust implementation of the positivity-preserving limiter than the one in \cite{22}. Numerical tests show that the third order DG scheme with the new positivity-preserving limiter produces satisfying results even without the TVB limiter.

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1 Introduction

Gas detonation is supersonic flow phenomena that consist of a precursor shock that ignites a combustible mixture gas and of a thin reaction zone immediately behind the shock. Although detonation has been studied for many years, it remains an active area of research in both theoretical studies and in numerical simulations due to practical importance. To study the gaseous detonation numerically, the governing equations could be chosen as the Euler equations describing inviscid compressible flow with the chemical reaction added. There are many difficulties to design stable numerical schemes solving a general hyperbolic system with source terms accurately. For example, the width of reaction zone attached to the shock might be very small, see [3], and the source term might induce stiffness, see [9].

In this paper, we focus on how to render numerical schemes stable. In practice, there are a lot of cases in which the density or pressure of the numerical solutions may become negative easily. For instance, highly energetic flows often contain regions with the kinetic part dominant, thus the internal energy is relatively very small. Another example is computational simulation of gas detonation propagation through different geometries. The shock diffraction may result in very low density and pressure. Under such conditions, it has been observed that numerical schemes will produce negative density or pressure, even for non-reactive gas flows, which may lead to blow-ups. This phenomenon tends to be amplified by the chemical activity. Crude replacement of negative values by positive ones not only destroyed local and global conservation, but was also found to be unstable. Therefore, it is strongly motivated to design schemes with a provable positivity-preserving property. Moreover, a conservative positivity-preserving scheme can be easily prove to have $L^1$ stability.

First order and second order positivity-preserving schemes were well studied in the literature [10]. So we are mainly interested in high order positivity-preserving schemes. On the one hand, low order schemes have been used in the simulation of detonation waves [12, 13], but numerical results have some deviation with the experimental results. On the other hand, some high order schemes have been developed in recent years [4, 5, 16, 19, 20]. Successful high
order numerical schemes for hyperbolic conservation laws, for example, the Runge-Kutta discontinuous Galerkin (RKDG) method in [1, 2], the essentially non-oscillatory (ENO) finite volume and finite difference schemes in [7, 18], and the weighted ENO (WENO) finite volume and finite difference schemes in [11, 8], do not satisfy a strict positivity-preserving property. In fact, they all fail for certain low density or low pressure test cases. Special treatment for certain schemes may achieve the positivity and conservation, but it is very difficult to maintain high order accuracy for smooth solutions. Constructing high order schemes which automatically preserve the positivity of density and pressure is highly nontrivial. In [22, 23], two of the authors proposed an arbitrarily high order positivity-preserving Runge-Kutta discontinuous Galerkin method for compressible Euler equations. The main idea is to find some straightforward sufficient condition for DG method with first order Euler forward time discretization to keep positivity. A simple limiter which is easy and cheap to implement will enforce the sufficient condition without destroying conservation and accuracy. Strong stability preserving (SSP) high order Runge-Kutta or multi-step method will still keep the positivity since they are convex combinations of Euler forward. With the limiter, high order RKDG method will be positivity-preserving of density and pressure in the cell average sense during the time evolution.

We will show an extension of this method to Euler system with an Arrhenius form chemical reaction source term and an additional equation for the evolution of reaction rate, which are typical governing equations for modelling the gaseous detonation. The positivity of the reaction rate is also crucial to the stability of schemes in this model. We also propose a more robust new implementation of the positivity-preserving limiter. The DG scheme with only the new positivity-preserving limiter is stable even for very strong shocks. Numerical tests of the third order DG method will be reported.
2 Positivity-preserving high order discontinuous Galerkin method for two-dimensional reactive Euler equations

2.1 Preliminaries

We consider the dimensionless two-dimensional compressible Euler equations with a source term representing chemical reactions for the ideal gas,

$$ w_t + f(w)_x + g(w)_y = s(w), \quad t \geq 0, (x, y) \in \mathbb{R}^2, $$

$$ w = \begin{pmatrix} \rho \\ m \\ n \\ E \\ \rho Y \end{pmatrix}, \quad f(w) = \begin{pmatrix} m \\ \rho u^2 + p \\ \rho uv \\ (E + p)u \\ \rho uY \end{pmatrix}, \quad g(w) = \begin{pmatrix} n \\ \rho uv \\ \rho v^2 + p \\ (E + p)v \\ \rho vY \end{pmatrix}, \quad s(w) = \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ \omega \end{pmatrix} $$

with

$$ m = \rho u, \quad n = \rho v, \quad E = \frac{1}{2}\rho u^2 + \frac{1}{2}\rho v^2 + \frac{p}{\gamma - 1} + \rho qY, $$

where $q$ is the heat release of reaction, $\gamma$ is the specific heat ratio and $Y$ denotes the reactant mass fraction. The source term is assumed to be in an Arrhenius form

$$ \omega = -\tilde{K} \rho Y e^{-\tilde{T}/T}, $$

where $T = \frac{\rho}{p}$ is the temperature, $\tilde{T}$ is the activation temperature and $\tilde{K}$ is a constant. The eigenvalues of the Jacobian $f'(w)$ are $u - c, u, u, u, u + c$ and the eigenvalues of the Jacobian $g'(w)$ are $v - c, v, v, v, v + c$ where $c = \sqrt{\frac{\gamma E}{p}}$.

Define the set of admissible states by

$$ G = \left\{ w = \begin{pmatrix} \rho \\ m \\ n \\ E \\ \rho Y \end{pmatrix} \middle| \rho > 0 \quad \text{and} \quad p(w) \geq 0, \quad Y \geq 0 \right\}, $$

then $G$ is a convex set since $p$ is a concave function of $w$. We are interested in schemes for (2.1) producing the numerical solutions in the admissible set $G$. We start with the
one-dimensional non-reactive equation $w_t + f(w)_x = 0$ and the first order Lax-Friedrichs scheme

$$w_j^{n+1} = w_j^n - \lambda [\widehat{f}(w_j^n, w_{j+1}^n) - \widehat{f}(w_{j-1}^n, w_j^n)],$$

(2.3)

$$\widehat{f}(u, v) = \frac{1}{2} [f(u) + f(v) - a(v - u)], \quad a = ||(|u| + c)||_\infty,$$  

(2.4)

where $n$ refers to the time step and $j$ to the spatial cell, and $\lambda = \frac{At}{Ax}$ is the ratio of time and space mesh sizes. Following Remark 2.4 in [22], it is easy to check that, for scheme (2.3), $w_j^n, w_{j+1}^n \in G$ implies $w_j^{n+1} \in G$ under the CFL condition $\lambda a \leq 1$. Other examples of positivity preserving fluxes include the Godunov flux, the Boltzmann type flux [14], and the HLLE flux.

### 2.2 Discontinuous Galerkin method

We review the formulation of DG method in [2] briefly. Assume $T_h$ is a triangulation of the spatial domain $\Omega$. For simplicity, we use $x$ to denote $(x, y)$ and $dx$ to denote $dxdy$ in this subsection. For each $t \geq 0$, we seek the approximation $w_h(x, t)$ in the piecewise polynomial space

$$V_h = \{v_h \in L^\infty(\Omega) : v_h|_K \in P^k(K), \forall K \in T_h\},$$

where $P^k$ denotes all the polynomials of degree $k$. The weak formulation of DG method solving (2.1) is, find $w_h \in V_h$ satisfying $\forall v_h \in V_h$,

$$\frac{d}{dt} \int_K w_h v_h dx + \sum_{e \in \partial K} \int_e \mathbf{h}(w_{h}^{int}, w_{h}^{ext}, \nu_e) v_h d\Gamma - \int_K \mathbf{F}(w_h) \cdot \nabla v_h dx = \int_K s(w_h) v_h dx,$$

(2.5)

where $\mathbf{F} = \langle f, g \rangle$, $\nu_e$ is the outward normal vector of the edge $e$ on the element $K$. We consider Lax-Friedrichs flux as an example throughout the rest of this paper,

$$\mathbf{h}(u, v, \nu) = \frac{1}{2} |\langle F(u) \cdot \nu + F(v) \cdot \nu - a(v - u)\rangle|, \quad a = ||(|u| + c)||_\infty.$$

Except the first one, the integrals in (2.5) can be approximated by proper quadrature rules. Time discretizations can be solved by the strong stability preserving (also called TVD) Runge-Kutta or multi-step time discretization. See [2] for more details.
To construct conservative positivity-preserving schemes, the most important step is to achieve positivity in the mean. We only need to discuss Euler forward time discretization because high order SSP time discretizations are convex combinations of Euler forward thus will keep the positivity due to the convexity of $G$. Take the test function as $v_h = 1$ in (2.5), we get the scheme satisfied by the cell average in the DG method. Consider the Euler forward time discretization,

$$
\frac{|K|}{\Delta t}(\bar{w}_K^{n+1} - \bar{w}_K^n) + \sum_{e \in \partial K} \int_e h(w_{h\text{int}}, w_{h\text{ext}}, \nu_e) d\Gamma = \int_K s(w_h) dx,
$$

where $\bar{w}_K^n$ denotes the cell average $w_h$ on $K$ at time level $n$ and $|K|$ is the area of $K$.

### 2.3 Rectangular meshes

For simplicity we assume we have a uniform rectangular mesh. At time level $n$, we have the DG polynomials of degree $k$, $w_{ij}(x, y) = (\rho_{ij}(x, y), m_{ij}(x, y), n_{ij}(x, y), E_{ij}(x, y), \rho Y_{ij}(x, y))^T$ with the cell average $\bar{w}_{ij}$ on the $(i, j)$ cell $[x_{i-\frac{1}{2}}, x_{i+\frac{1}{2}}] \times [y_{j-\frac{1}{2}}, y_{j+\frac{1}{2}}]$. Let $w_{i-\frac{1}{2}, j}^+(y), w_{i+\frac{1}{2}, j}^-(y), w_{i,j-\frac{1}{2}}^+(x), w_{i,j+\frac{1}{2}}^-(x)$ denote the traces of $w_{ij}(x, y)$ on the four edges respectively.

Assume that we use a $L$-point Gauss quadrature where $L \geq k + 1$ (see [2] for an analysis of the requirement of the numerical quadrature for the accuracy of the DG solution). Let $S_{i}^x = \{x_i^\beta : \beta = 1, \cdots, L\}$ denote the Gauss quadrature points on $[x_{i-\frac{1}{2}}, x_{i+\frac{1}{2}}]$, and $S_{j}^y = \{y_j^\beta : \beta = 1, \cdots, L\}$ denote the Gauss quadrature points on $[y_{j-\frac{1}{2}}, y_{j+\frac{1}{2}}]$. For instance, $(x_{i-\frac{1}{2}}, y_j^\beta)$ $(\beta = 1, \cdots, L)$ are the Gauss quadrature points on the left edge of the $(i, j)$ cell. The subscript $\beta$ will denote the values at the Gauss quadrature points, for instance, $w_{i-\frac{1}{2}, j}^+ = w_{i-\frac{1}{2}, j}^+(y_j^\beta)$. Also, $w_{\beta}$ denotes the corresponding quadrature weight on interval $[-\frac{1}{2}, \frac{1}{2}]$, so that $\sum_{\beta=1}^L w_{\beta} = 1$. We need to use the $N$-point Gauss-Lobatto quadrature rule where $N$ is the smallest integer such that $2N - 3 \geq k$, and we distinguish the two quadrature rules by adding hats to the Gauss-Lobatto points, i.e., $\hat{S}_{i}^x = \{\hat{x}_i^\alpha : \alpha = 1, \cdots, N\}$ will denote the Gauss-Lobatto quadrature points on $[x_{i-\frac{1}{2}}, x_{i+\frac{1}{2}}]$, and $\hat{S}_{j}^y = \{\hat{y}_j^\alpha : \alpha = 1, \cdots, N\}$ will denote the Gauss-Lobatto quadrature points on $[y_{j-\frac{1}{2}}, y_{j+\frac{1}{2}}]$. 

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Then (2.6) becomes

\[
\overline{w}_{ij}^{n+1} = \overline{w}_{ij}^n - \lambda_1 \sum_{\beta=1}^{L} w_{\beta} \left[ h_1 \left( w_{i+\frac{1}{2},\beta}, w_{i+\frac{1}{2},\beta+1} \right) - h_1 \left( w_{i-\frac{1}{2},\beta}, w_{i-\frac{1}{2},\beta+1} \right) \right] \\
- \lambda_2 \sum_{\beta=1}^{L} w_{\beta} \left[ h_2 \left( w_{\beta,j+\frac{1}{2}}, w_{\beta,j+\frac{1}{2}} \right) - h_2 \left( w_{\beta,j-\frac{1}{2}}, w_{\beta,j+\frac{1}{2}} \right) \right] \\
+ \Delta t \sum_{\alpha=1}^{L} \sum_{\beta=1}^{L} w_{\alpha} w_{\beta} s(w(x_i^\alpha, y_j^\beta))
\]

(2.7)

where \( \lambda_1 = \frac{\Delta t}{\Delta x} \), \( \lambda_2 = \frac{\Delta t}{\Delta y} \) and

\[
h_1(u, v) = \frac{1}{2} [f(u) + f(v) - a(v - u)]
\]

\[
h_2(u, v) = \frac{1}{2} [g(u) + g(v) - a(v - u)].
\]

We use \( \otimes \) to denote the tensor product, for instance,

\[
S^x_i \otimes S^y_j = \{ (x, y) : x \in S^x_i, y \in S^y_j \}.
\]

Define the set \( S_{ij} \) as

\[
S_{ij} = (S^x_i \otimes \hat{S}^y_j) \cup (\hat{S}^x_i \otimes S^y_j) \cup (S^x_i \otimes S^y_j).
\]

(2.8)

**Theorem 2.1.** If the DG polynomial \( w_{ij}(x, y) \in G, \forall (x, y) \in S_{ij} \), then the scheme (2.7) is positivity-preserving, namely, \( \overline{w}_{ij}^{n+1} \in G \) under the time step restriction

\[
a(\lambda_1 + \lambda_2) \leq \frac{1}{2} \bar{w}_1, \quad \Delta t \bar{K} \min e^{-\bar{T}/T} \leq \frac{1}{2},
\]

where the minimum is taken over \( S^x_i \otimes S^y_j \) for all the rectangles.

**Proof.** We can rewrite (2.6) as \( w_{ij}^{n+1} = \frac{1}{2} C + \frac{1}{2} S \) where

\[
C = \overline{w}_{ij}^n - 2\lambda_1 \sum_{\beta=1}^{L} w_{\beta} \left[ h_1 \left( w_{i+\frac{1}{2},\beta}, w_{i+\frac{1}{2},\beta+1} \right) - h_1 \left( w_{i-\frac{1}{2},\beta}, w_{i-\frac{1}{2},\beta+1} \right) \right] \\
- 2\lambda_2 \sum_{\beta=1}^{L} w_{\beta} \left[ h_2 \left( w_{\beta,j+\frac{1}{2}}, w_{\beta,j+\frac{1}{2}} \right) - h_2 \left( w_{\beta,j-\frac{1}{2}}, w_{\beta,j+\frac{1}{2}} \right) \right]
\]

and

\[
S = \overline{w}_{ij}^n + 2\Delta t \sum_{\alpha=1}^{L} \sum_{\beta=1}^{L} w_{\alpha} w_{\beta} s(w(x_i^\alpha, y_j^\beta)).
\]
By Theorem 3.1 in [22], we have $C \in G$. So it suffices to prove $S \in G$. Gauss quadrature rule implies
\[ w_{ij}^n = \sum_{\alpha=1}^{L} \sum_{\beta=1}^{L} w_{\alpha\beta} w(x_i^\alpha, y_j^\beta). \] (2.9)

Thus,
\[ S = \sum_{\alpha=1}^{L} \sum_{\beta=1}^{L} w_{\alpha\beta} \left[ w(x_i^\alpha, y_j^\beta) + 2\Delta ts \left( w(x_i^\alpha, y_j^\beta) \right) \right]. \]

Given $w \in G$, then it is easy to check that $w + 2\Delta ts(w) \in G$ if $\Delta t \tilde{K} e^{-T/T} \leq \frac{1}{2}$. So $S \in G$. □

2.4 Triangular meshes

For each triangle $K$ we denote by $l_K^i$ ($i = 1, 2, 3$) the length of its three edges $e_K^i$ ($i = 1, 2, 3$), with outward unit normal vector $\nu^i$ ($i = 1, 2, 3$). Assume the line integrals in (2.6) are solved by $L$-point Gauss quadrature where $L \geq k + 1$. And the source integral is solved by the $M$-point Gauss quadrature on a triangle in which $x_K^\gamma$ and $\tilde{w}_\gamma$ denote the quadrature points and normalized weights with $\sum_{\gamma=1}^{M} \tilde{w}_\gamma = 1$.

Then (2.6) becomes
\[ \bar{w}_K^{n+1} = \bar{w}_K^n - \frac{\Delta t}{|K|} \sum_{i=1}^{3} \sum_{\beta=1}^{L} h w_{i,\beta}^{int(K)} w_{\beta}^{ext(K)} w_{\beta} l_K^i + \Delta t \sum_{\gamma=1}^{M} \tilde{w}_\gamma s(w(x_K^\gamma)), \] (2.10)

where and $w_{i,\beta}^{int(K)}$ and $w_{i,\beta}^{ext(K)}$ denote the values of $u$ evaluated at the $\beta$-th Gauss quadrature point on the $i$-th edge from the interior and exterior of the element $K$ respectively.

We need a special quadrature on a triangle introduced in [24]. In the barycentric coordinates, the set $S_K^k$ of quadrature points for polynomials of degree $k$ on a triangle $K$ can be written as
\[ S_K^k = \{ \left( \frac{1}{2} + v^\beta, \frac{1}{2} + \tilde{u}^\alpha \right), \left( \frac{1}{2} - \tilde{u}^\alpha, \frac{1}{2} - v^\beta \right), \left( \frac{1}{2} - v^\beta, \frac{1}{2} + \tilde{u}^\alpha \right), \left( \frac{1}{2} + \tilde{u}^\alpha, \frac{1}{2} - v^\beta \right), \left( \frac{1}{2} - v^\beta, \frac{1}{2} - \tilde{u}^\alpha \right), \left( \frac{1}{2} + v^\beta, \frac{1}{2} + \tilde{u}^\alpha \right) \}. \] (2.11)
where $u^\alpha (\alpha = 1, \cdots, N)$ and $v^\beta (\beta = 1, \cdots, L)$ are the Gauss-Lobatto and Gauss quadrature points on the interval $[-\frac{1}{2}, \frac{1}{2}]$ respectively. Define $S_K$ as the union of $S^k_K$ and the $M$-point Gauss quadrature on the triangle $K$. Following Theorem 5.1 in [24] and Theorem 2.1, we have

**Theorem 2.2.** If the DG polynomial $w_K(x, y) \in G$, $\forall (x, y) \in S_K$, then the scheme (2.10) is positivity-preserving, namely, $\overline{w}^{n+1}_K \in G$ under the time step restriction

$$a \frac{\Delta t}{|K|} \sum_{i=1}^{3} l_i^K \leq \frac{1}{3} \tilde{w}_1, \quad \Delta t \min e^{-\tilde{T}/T} \leq \frac{1}{2},$$

where the minimum is taken over $x^\gamma_K, \gamma = 1, \cdots, M$ for all the triangles.

**2.5 $L^1$ stability**

The limiter in [22] can be used to enforce the conditions in Theorem 2.1 and Theorem 2.2. We will describe an improved implementation of this limiter in the next section.

With the limiter added, the full high order DG scheme will keep density, pressure and reactant mass fraction non-negative in the mean during the time evolution.

**Theorem 2.3.** Assuming vanishing, reflective or periodic boundary conditions, suppose the DG polynomial satisfies $w_K(x) \in G, \forall x \in S_K$, then the scheme (2.10) satisfy the following $L^1$ stability:

$$\sum_K |\overline{p}^{n+1}_K| = \sum_K |\overline{p}^n_K|, \sum_K |\overline{E}^{n+1}_K| = \sum_K |\overline{E}^n_K|, \sum_K |\overline{\rho Y}^{n+1}_K| \leq \sum_K |\overline{\rho Y}^n_K|.$$

**Proof.** Take the summation of (2.10) for all $K$, then we have

$$\sum_K \overline{w}^{n+1}_K = \sum_K \overline{w}^n_K + \Delta t \sum_{\gamma=1}^{M} \tilde{w}_\gamma s(w(x^\gamma_K)).$$

The first component reads $\sum_K \overline{p}^{n+1}_K = \sum_K \overline{p}^n_K$. Theorem (2.2) and (2.9) imply that $\overline{w}^{n+1}_K, \overline{w}^n_K \in G$, thus $\overline{p}^{n+1}_K, \overline{p}^n_K \geq 0$. Therefore, $\sum_K |\overline{p}^{n+1}_K| = \sum_K |\overline{p}^n_K|$. Similarly, we get $\sum_K |\overline{E}^{n+1}_K| = \sum_K |\overline{E}^n_K|$. Notice that the fifth component of $s(w(x^\gamma_K))$ is non-positive, we get the last inequality. \qed
3 An improved implementation of the positivity-preserving limiter

3.1 Positivity-preserving limiter

At the time level \( n \), given the DG polynomial \( w_K(x) \) with the cell average \( \bar{w}_K \in G \), we would like to modify it into another polynomial

\[
\tilde{w}_K(x) = \theta_K(w_K(x) - \bar{w}_K) + \bar{w}_K \tag{3.1}
\]

where \( \theta_K \in [0,1] \) is to be determined, such that \( \tilde{w}_K(x) \in G, \forall x \in S_K \). If \( \theta_K \) is the largest such number (the smallest one is \( \theta_K = 0 \)), then this limiter will not destroy accuracy for smooth solutions, see [22]. Following [22], it can be implemented as:

1. First, enforce the positivity of density (and reactant mass fraction). For each element \( K \), compute

\[
\hat{\rho}_K(x) = \theta^1_K [\rho_K(x) - \overline{\rho}_K] + \overline{\rho}_K, \quad \theta^1_K = \min_{x \in S_K} \left\{ 1, \frac{\overline{\rho}_K}{\rho_K - \rho_K(x)} \right\}, \tag{3.2}
\]

\[
\hat{\rho}Y_K(x) = \theta^2_K [\rho Y_K(x) - \rho Y_K] + \rho Y_K, \quad \theta^2_K = \min_{x \in S_K} \left\{ 1, \frac{\rho Y_K}{\rho Y_K - \rho Y_K(x)} \right\}. \tag{3.3}
\]

2. Second, enforce the positivity of pressure. Define \( \tilde{w}_K = (\hat{\rho}_K, m_K, n_K, E_K, \hat{\rho}Y_K)^T \). For each \( x \in S_K \), if \( p(\tilde{w}_K(x)) > 0 \) define \( \theta_x = 1 \); otherwise, define \( \theta_x \) as the solution of

\[
p(\theta_x(\tilde{w}_K(x) - \bar{w}_K) + \bar{w}_K) = 0. \tag{3.4}
\]

Then get the limited polynomial

\[
\tilde{w}_K(x) = \theta_K(\tilde{w}_K(x) - \bar{w}_K) + \bar{w}_K, \quad \theta_K = \min_{x \in S_K} \theta_x. \tag{3.5}
\]

Even though we only need to solve a quadratic equation of \( \theta_x \) in (3.4), in practice \( \theta_K \) solved from (3.4) cannot guarantee the strict non-negativity of \( p(\tilde{w}_K(x)) \) numerically due to the round off errors for some wild data, e.g., blast waves. In [22], it was reported that for
problems with very strong shocks, positivity limiter itself implemented as above may not be stable, thus TVB limiter must be used.

Here we propose a slightly different but very robust implementation of (3.4) so that TVB limiter will no longer be needed. Notice that $p$ is a concave function of $w$, thus we have the Jensen’s inequality

$$p(\theta(w - \bar{w}) + \bar{w}) = p(\theta w + (1 - \theta)\bar{w}) \geq \theta p(w) + (1 - \theta)p(\bar{w}). \quad (3.6)$$

Therefore, if $\bar{w} \in G$ and $\rho(w) > 0$,

$$\theta = \frac{p(\bar{w})}{p(w) - p(\bar{w})} \quad (3.7)$$

satisfies that $p(\theta(w - \bar{w}) + \bar{w}) \geq 0$. Although $\theta$ defined in (3.7) is smaller than the real solution of $p(\theta(w - \bar{w}) + \bar{w}) = 0$, it is actually the similar type as $\theta_1$ in (3.2). It is straightforward to prove the accuracy for smooth solutions following the argument in [22].

We can formulate the new robust implementation of the limiter as, for each element $K$,

1. Compute (3.2) and (3.3).

2. Define $\hat{w}_K = (\hat{\rho}_K, m_K, n_K, E_K, \hat{Y}_K)^T$. For each $x \in S_K$, if $p(\hat{w}_K(x)) \geq 0$ set $\theta_x = 1$; otherwise, set $$\theta_x = \frac{p(\bar{w}_K)}{p(w_K) - p(\hat{w}_K(x))}.$$ Then get the limited polynomial (3.5).

### 3.2 The algorithm for SSP Runge-Kutta time discretizations

Theoretically, there is a complication regarding the time step restriction in Theorem 2.1 and Theorem 2.2 for a Runge-Kutta time discretization. Consider the third order SSP Runge-Kutta method. To enforce the CFL condition rigorously, we need to get an accurate estimation of $a = \| (|u| + c) \|_\infty$ for all the three stages based only on the numerical solution at time level $n$, which is highly nontrivial mathematically. An efficient solution is, if a preliminary calculation to the next time step produces negative density or pressure, then
reCalculate from the time step \( n \) with half the previous time step. This complication does not exist if we use a SSP multi-step time discretization.

The algorithm flow chart for the third order SSP Runge-Kutta method on triangular meshes is

1. Given the DG polynomials \( \mathbf{w}_K(x) \) at time step \( n \) satisfying the cell average \( \overline{\mathbf{w}}_K \in G \) and \( \mathbf{w}_K(x) \in G, \forall x \in S_K \), calculate \( a = \max \| \langle \langle u, v \rangle \rangle + c \|, \ b = \max \tilde{K} e^{-\tilde{T}/T} \) where the maximum is taken over \( S_K \) for all \( K \). Set the time step

\[
\Delta t = \min \left\{ \frac{1}{3} \overline{\mathbf{w}}_1 |K|, \ \frac{1}{2b} \right\}.
\]

2. Calculate the first stage with \( \mathbf{w}_k(x) \). Let \( \mathbf{w}^1_K(x) \) denote the solution of the first stage. Modify it by the limiter (3.5) into \( \tilde{\mathbf{w}}^1_K(x) \).

3. Calculate the second stage with \( \tilde{\mathbf{w}}^1_K(x) \). Let \( \mathbf{w}^2_K(x) \) denote the solution of the second stage. If its cell average is not in \( G \) (by Theorem 2.2, this means that \( a \) or \( b \) calculated based on \( \mathbf{w}_K(x) \) is smaller than the ones of \( \tilde{\mathbf{w}}^1_K(x) \)), then go back to step two and restart with half time step; otherwise, modify it by the limiter (3.5) into \( \tilde{\mathbf{w}}^2_K(x) \).

4. Calculate the third stage with \( \tilde{\mathbf{w}}^2_K(x) \). Let \( \mathbf{w}^3_K(x) \) denote the solution of the third stage. If its cell average is not in \( G \) (by Theorem 2.2, this means that \( a \) or \( b \) calculated based on \( \mathbf{w}_K(x) \) is smaller than the ones of \( \tilde{\mathbf{w}}^2_K(x) \)), then go back to step two and restart with half time step; otherwise, modify it by the limiter (3.5) into \( \tilde{\mathbf{w}}^3_K(x) \).

4 Numerical Tests

4.1 Euler equations

We test the robustness of the new implementation of the positivity-preserving limiter for non-reactive Euler equations by using the algorithm in the previous section.

In [1, 2], the TVB limiter was used to kill oscillations for high order DG schemes solving Euler equations with strong shocks. For smooth solutions, the TVB limiter will not destroy
the accuracy. However, TVB limiter is not sufficient for stabilizing high order schemes solving Euler equations when low density or low pressure emerges. In [22, 24, 23], the third order DG scheme with TVB limiter and positivity-preserving limiter performed very well for all test cases for which DG method with only TVB limiter will blow up due to the presence of negative density or negative pressure.

By the following numerical tests, we will see that the positivity-preserving limiter itself can stabilize the high order DG schemes without TVB limiter. We test the third order DG method and the third order SSP Runge-Kutta time discretization with only the positivity-preserving limiter (3.5), solving the one-dimensional compressible Euler equations for ideal gas with $\gamma = 1.4$.

**Example 4.1.** *Shock tube problems.*

See Figure 4.1 for the results of Sod and Lax problems using 100 cells, which are comparable to the results of RKDG method with characteristicwise TVB limiter in [1].

**Example 4.2.** *Interaction of blast waves.*

This example is the same as the one in [1]. See Figure 4.2 for the comparison of results of two limiters. As we expected, TVB can kill oscillations while the positivity limiter cannot, however, the positivity limiter alone will smear the discontinuity less.

**Example 4.3.** *Sedov point blast.*

The initial and boundary conditions are the same as in [22]. See Figure 4.3 for the result of DG with only positivity limiter.

### 4.2 The reactive Euler equations

In this subsection, we show the test results for the third order RKDG method with only positivity-preserving limiter solving (2.1). The parameters are $\gamma = 1.2$, $q = 50$, $\tilde{T} = 50$, $\tilde{K} = 2566.4$. For all the test cases in this subsection, the RKDG method with only TVB limiter may blow up at a certain time.
Figure 4.1: Shock Tube Problems. $P^2$ element DG with only positivity-preserving limiter. The solid lines are the exact solutions. The symbols are the numerical solutions.
Figure 4.2: Interaction of blast waves. $P^2$ element DG with two different limiters on 400 cells. The solid line is computed by the fifth order WENO on a very fine mesh. The symbols are the numerical solutions.
Example 4.4. Convergence study.

We test the convergence of our scheme in this example. The domain is \([0, 2] \times [0, 2]\). The initial condition is, if \(x^2 + y^2 \leq 0.36\), then \((\rho, u, v, p, Y) = (1, 0, 0, 80, 0)\); otherwise, \((\rho, u, v, p, Y) = (1, 0, 0, 10^{-9}, 1)\). The boundary conditions for the bottom and the left are reflective. The terminal time is \(t = 0.2\). The mesh is uniformly rectangular. See the comparison of results of \(\Delta x = \Delta y = \frac{1}{60}\) and \(\Delta x = \Delta y = \frac{1}{120}\) in Figure 4.4.

Example 4.5. Detonation diffraction problems.

The simulation of gaseous detonation waves through different geometries is numerically challenging especially for the high order schemes mainly because the pressure or density may drop very close to zero when the shock wave is diffracted. Here we test the detonation diffraction at three different angles.

The first one is ninety degrees. The initial conditions are, if \(x < 0.5\), \((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\); otherwise, \((\rho, u, v, E, Y) = (1, 0, 0, 55, 1)\). The boundary conditions are reflective except that at \(x = 0\), \((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\). The terminal time is \(t = 0.6\). The mesh is uniformly rectangular. See Figure 4.6 for the result of \(\Delta x = \Delta y = \frac{1}{48}\).
Figure 4.4: Convergence study. The colored contour and the solid line on the right is the result of $\Delta x = \Delta y = \frac{1}{120}$. The symbols on the right denote the result of $\Delta x = \Delta y = \frac{1}{60}$.
The second one is sixty degrees. The initial conditions are, if \( x < 0.6 \) and \( y \geq 2 \), 
\((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\); otherwise, 
\((\rho, u, v, E, Y) = (1, 0, 0, 55, 1)\). The boundary conditions are reflective except that at \( x = 0 \), 
\((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\). The terminal time is \( t = 0.68 \). The mesh is nonuniform, mixed with rectangles and triangles. See Figure 4.5(a) for the illustration of the mesh. See Figure 4.6 for the result where the length of the smallest edge in the mesh is \( \frac{1}{32\sqrt{3}} \).

The third one is forty-five degrees. The initial conditions are, if \( x < 1.5 \) and \( y \geq 2 \), 
\((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\); otherwise, 
\((\rho, u, v, E, Y) = (1, 0, 0, 55, 1)\). The boundary conditions are reflective except that at \( y = 0 \), 
\((\rho, u, v, E, Y) = (11, 6.18, 0, 970, 1)\). The terminal time is \( t = 0.68 \). The mesh is uniform, mixed with rectangles and triangles. See Figure 4.5(b) for the illustration of the mesh. See Figure 4.6 for the result of the mesh size \( \frac{1}{21} \).

**Example 4.6.** Multiple obstacles.

The initial condition is, if \( x^2 + y^2 \leq 0.36 \), then 
\((\rho, u, v, E, Y) = (7, 0, 0, 200, 0)\); otherwise, 
\((\rho, u, v, E, Y) = (1, 0, 0, 55, 1)\). The boundary conditions are reflective everywhere. The location of the first obstacle is \([1.3, 3.3] \times [0, 2.6]\) and the second one is \([5.1, 8.3] \times [0, 4.3]\). The terminal time is \( t = 1.4 \). The parameters are set as \( \gamma = 1.2, q = 50, \tilde{T} = 20, \tilde{K} = 2410.2 \). The mesh is nonuniformly rectangular. See Figure 4.9 for the result where the length of the smallest edge in the mesh is \( \frac{0.85}{32} \).

## 5 Concluding remarks

We have shown an extension of the positivity-preserving techniques in [22, 24, 23] to construct robust high order RKDG schemes for reactive Euler equations modelling gaseous detonations. Numerical tests suggest that positivity-preserving is sufficient to stabilize the high order DG method without the TVB limiter, and robust high order RKDG schemes can successfully simulate detonation diffraction cases in which the density or pressure of the numerical solutions may become negative easily. In the future work, we will use the RKDG
schemes to carry out numerical simulation on gaseous detonation in complex geometrical configurations in order to have more comprehensive insight of its propagation mechanism.
Figure 4.6: Detonation Diffraction at a Ninety-Degree Corner.
Figure 4.7: Detonation Diffraction at a Sixty-Degree Corner
Figure 4.8: Detonation Diffraction at a Forty-Five-Degree Corner
Figure 4.9: Multiple obstacles
References


