Novel local smoothness indicators for improving the third-order WENO scheme

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SUMMARY

The local smoothness indicators play an important role in the performance of a weighted essentially non-oscillatory (WENO) scheme. Due to having only two points available on each sub-stencil, the local smoothness indicators calculated by conventional methods of Jiang and Shu[1] make the third-order WENO scheme too dissipative. In this paper, we propose a different method to calculate the indicators by using all the three points on the global stencil of the third-order WENO scheme. The numerical results demonstrate that the WENO scheme with the new indicators has less dissipation and better resolution than the ones of Jiang and Shu’s for both smooth and discontinuous solutions. Copyright © 2017 John Wiley & Sons, Ltd.

KEY WORDS: third-order WENO scheme; local smoothness indicator; numerical dissipation

1. INTRODUCTION

The weighted essentially non-oscillatory (WENO) scheme was first proposed by Liu et al.[2]. The main idea is that, instead of choosing the smoothest stencil in the ENO (essentially non-oscillatory) scheme[3], the WENO scheme uses a convex combination of all candidate stencils (sub-stencils) to obtain high order accuracy in smooth regions while keeping the ENO property near shock waves. The analysis of Jiang and Shu[1] showed that, with the local smoothness indicators introduced by Liu et al.[2], the WENO scheme constructed from the rth order ENO scheme is only (r + 1)th order accurate. Then, Jiang and Shu[1] proposed a classical way to calculate the local smoothness indicators, with which the WENO (WENO-JS) scheme can achieve (2r − 1)th order accuracy.
order. Balsara and Shu[4] extended the WENO schemes up to eleventh-order and designed a class of monotonicity preserving WENO schemes. However, Henrick et al.[5] pointed out that the indicators of Jiang and Shu[1] failed to improve the accuracy of WENO scheme at critical points and then proposed a mapping function to modify the weights. The new WENO (WENO-M) scheme can achieve the optimal convergence order at critical points. Gerolymos et al.[6] further extended both WENO-JS and WENO-M schemes up to the seventeenth-order. In [7], Borges et al. designed a global smoothness indicator (WENO-Z scheme) to improve the accuracy of the fifth-order WENO-JS scheme directly. Later, Castro et al.[8] developed higher order WENO-Z schemes. Ha et al.[9] derived a different method that measures the local smoothness of the numerical solution inside a stencil and devised a sixth-order global smoothness indicator to construct the fifth-order WENO-Z-type scheme. Recently, Fan et al.[10] constructed higher order (up to eighth-order) global smoothness indicators, with which the associated WENO-Z-type schemes can obtain fifth convergence order in smooth regions, even at the second order critical points where both the first and second derivatives vanish.

Compared with higher order WENO schemes, the third-order one has several advantages, for example, it is more robust for shock problems, it uses less grid points and hence it reduces the difficulty of boundary treatment and can be easily generalized to unstructured meshes, and also it provides a suitable compromise of the computation cost and the accuracy in some cases. However, the third-order WENO scheme with the smoothness indicators of Jiang and Shu[1] is too dissipative. And those techniques, such as the mapping function of WENO-M scheme[5] and the global stencil indicator of WENO-Z scheme[7], can effectively improve the accuracy of higher order WENO schemes[5, 6, 7, 8], but they cannot get satisfactory improvement on the third-order one. In [11], Yamaleev and Carpenter developed a third-order energy stable WENO (ESWENO) scheme, which can decrease the dissipation of the third-order WENO scheme apparently. However, the two tuning parameters $\epsilon$ and $\delta$, which are designed for calculating the weight function and the artificial dissipation term of ESWENO scheme, are both dependent on the mesh size (or the total number of grid cells), and $\epsilon$ is also determined by the initial solution. Hence, the applications of the ESWENO scheme are limited. In [12], Wu et al. proposed a new global smoothness indicator and developed a less dissipative third-order WENO-Z-type (WENO-N3) scheme. Because WENO-N3 cannot reach the expected goal of high accuracy and low-dissipation, later, in [13], they suggested using the power function of the global indicator as the new one to construct the WENO-NP3 scheme. However, the new global smoothness indicator and the local smoothness indicators have different dimensions, this makes the results of the WENO-NP3 scheme uncertain in applications, in addition, WENO-NP3 generates apparent oscillatory solutions in many problems with shock waves.

As discussed in [1], the local smoothness indicators play a key role in the performance of a WENO scheme. Those works in [11, 12, 13] paid attention mainly to design the global smoothness indicators based on the two local smoothness indicators, which are calculated on each of sub-stencils. Hence, those improved third-order WENO schemes are still not so satisfactory. In this paper, we propose a novel way to calculate the two local smoothness indicators by using all three points on the global stencil. Analysis shows that, in monotonic smooth regions, the two new local indicators presented in this paper are the same, hence the calculated weights are equal to the optimal weights. Because of this important advantage, the new WENO scheme can greatly decrease the numerical dissipation.

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2. THIRD-ORDER WENO SCHEMES

In this section, we briefly describe the WENO schemes by using the one-dimensional scalar conservative law equation[1]
\[
\frac{\partial u}{\partial t} + \frac{\partial f}{\partial x} = 0
\]
where \( u(x, t) \) is a conserved quantity, \( f(u) \) describes its flux, and \( x \) and \( t \) denote space and time, respectively. By defining the points \( x_i = i\Delta x, (i = 0, \ldots, N) \), where \( \Delta x \) is the uniform grid spacing, the (1) can be approximated by a conservative finite difference formula
\[
\frac{du_i}{dt} = -\frac{\hat{f}_{i+1/2} - \hat{f}_{i-1/2}}{\Delta x}
\]
where \( \hat{f}_{i+1/2} \) is the numerical flux.

The flux \( \hat{f}_{i+1/2} \) of a third-order WENO scheme can be written as
\[
\hat{f}_{i+1/2} = \omega_0 q_0 + \omega_1 q_1
\]
where, \( q_k \) is the second-order flux on the sub-stencil \( S_k = (i - 1 + k, i + k) \) and is given by
\[
\begin{cases}
q_0 = \frac{1}{2} f_{i-1} + \frac{3}{2} f_i \\
q_1 = \frac{1}{2} f_{i} + \frac{1}{2} f_{i+1}
\end{cases}
\]
the weight \( \omega_k \) is constructed as
\[
\omega_k = \frac{\alpha_k}{\alpha_0 + \alpha_1}, \quad \alpha_k = \frac{c_k}{(IS_k + \epsilon)^p}, \quad k = 0, 1
\]
where, \( IS_k \) is the local smoothness indicator that measures the smoothness of numerical solution on the sub-stencil \( S_k \). \( c_0 = 1/3 \) and \( c_1 = 2/3 \) are called the optimal weights since they generate the third-order upstream scheme. \( \epsilon \) is a positive real number introduced to avoid the denominator becoming zero. \( \epsilon = 10^{-40} \) is recommended by Henrick et al.[5] and used in this paper. The larger power \( p \) can make the weight assigned to the non-smooth sub-stencil approach to zero faster, and also result in more dissipative WENO schemes[8]. Numerical results in [11, 12, 13] showed that \( p = 1 \) is adequate for a third-order WENO scheme to preserve ENO property and also is used in all tested third-order WENO schemes in this paper.
2.1. The WENO-JS scheme

In [1], Jiang and Shu proposed the classical smoothness indicator as

\[ IS_k = \sum_{l=1}^{r-1} \int_{x_{i-l/2}}^{x_{i+l/2}} (\Delta x)^{2l-1} (q_k^{(l)})^2 dx \]  

(6)

where, \( q_k^{(l)} \) is the \( l \)th order derivative of \( q_k(x) \), and \( q_k(x) \) is the interpolation polynomial on sub-stencil \( S_k \). For the third-order WENO scheme, \( r = 2 \), and eqrefeq:jiang gives the same formula as those of Liu et al.[2]

\[
\begin{align*}
IS_0 & = (f_i - f_{i-1})^2 \\
IS_1 & = (f_{i+1} - f_i)^2
\end{align*}
\]

(7)

In smooth regions, the Taylor series expansions of (7) at \( x_i \) give

\[
\begin{align*}
IS_0 & = (f'_i \Delta x - \frac{1}{2} f''_i \Delta x^2 + O(\Delta x^3))^2 \\
IS_1 & = (f'_i \Delta x + \frac{1}{2} f''_i \Delta x^2 + O(\Delta x^3))^2
\end{align*}
\]

(8)

(8) can be written as

\[ IS_k = A(1 + O(\Delta x)) \]  

(9)

where, \( A \) is independent of \( k \). Substituting (9) into (5), we have

\[ \omega_k = c_k + O(\Delta x) \]  

(10)

Similar to the work of Henrick et al.[5], Yamaleev and Carpenter[11] derived the necessary and sufficient conditions that provide third-order convergence of a third-order WENO scheme. A sufficient condition is given by

\[ \omega_k = c_k + O(\Delta x^2) \]  

(11)

Note that (11) imposes a more severe constraint on the weights compared with (10). And numerical experiments[11, 12] also showed that the WENO-JS scheme with the indicators (7) is too dissipative.

2.2. The WENO-M scheme

In [5], Henrick et al. proposed a mapping function to improve the approximation accuracy of \( \omega_k \) to \( c_k \). The mapping function is defined as

\[ g_k(\omega_k) = \frac{\omega_k (c_k + c_k^2 - 3 c_k \omega_k + \omega_k^2)}{c_k^2 + \omega_k (1 - 2 c_k)} \]  

(12)

Applying the mapping function, the new weights for the third-order WENO-M scheme are

\[ \omega_k^M = \frac{\alpha_k}{\alpha_0 + \alpha_1}, \quad \alpha_k = g_k(\omega_k) \]  

(13)

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Using the similar analysis in [5], it is easy to find that

\[ \omega^M_k = c_k + O(\Delta x^3) \] (14)

(14) indicates theoretically that the weights (13) go beyond the sufficient condition (11) for a third-order WENO scheme. However, the numerical results (See Section 4.1) showed that the WENO-M scheme cannot reach the third-order accuracy. And hence the issue of improving the third-order WENO scheme is still open.

2.3. The WENO-Z-type schemes

As pointed out by Borges et al.[7], the mapping function used to construct the WENO-M scheme is computationally expensive. Hence, similar as the fifth-order WENO-Z scheme[7], one can easily construct a third-order WENO-Z-type scheme by introducing a global smoothness indicator \( \tau \), and the weights are calculated as

\[ \omega^Z_k = \frac{\alpha_k}{\alpha_0 + \alpha_1}, \quad \alpha_k = c_k \left( 1 + \frac{\tau}{IS_k + \epsilon} \right) \] (15)

For example, by directly generalizing the method of Borges et al., one can obtain a global smoothness indicator as

\[ \tau_3 = |IS_1 - IS_0| \] (16)

Yamaleev and Carpenter[11] proposed a different indicator for the ESWENO scheme as

\[ \tau_E = (f_{i-1} - 2f_i + f_{i+1})^2 \] (17)

and Wu and Zhao[12] constructed a indicator for the WENO-N3 scheme as

\[ \tau_N = \left| \frac{IS_0 + IS_1}{2} - IS_3 \right| \] (18)

where, \( IS_3 = \frac{13}{12}(f_{i-1} - 2f_i + f_{i+1})^2 + \frac{1}{4}(f_{i-1} - f_{i+1})^2 \) is the local smoothness indicator at the second sub-stencil \( S_1 = (x_{i-1}, x_i, x_{i+1}) \) in the fifth-order WENO scheme[1]. In fact, the two indicators (17) and (18) have the same expression as \( \tau = \gamma(f_{i-1} - 2f_i + f_{i+1})^2 \), in which, \( \gamma = 1 \) and \( \gamma = 5/6 \) are taken in (17) and (18), respectively. Hence, in this paper, we only discuss the WENO-N3 scheme with \( \tau_N \).

In smooth regions, by means of the Taylor series expansion and simple derivation, one can get

\[ \omega^{Z3}_k = c_k + O(\Delta x) \] (19)

\[ \omega^{TN}_k = c_k + O(\Delta x^2), \quad f'_i \neq 0 \] (20)

It should be pointed out that, if \( f'_i = 0 \), the global smoothness indicator \( \tau_N/\tau_E \) and the local ones (7) have the same order of magnitude. The numerical results in [12] showed that the improved third-order WENO schemes above cannot reach the expected goal.
In [13], Wu et al. proposed a modified indicator $\tau_{NP} = (\tau_N)^p$ ($p = 3/2$) to construct the WENO-NP3 scheme to improve the accuracy of the WENO-N3 scheme in smooth regions. However, such a modification of $\tau_N$ destroys the requirement that the global smoothness indicator and the local ones should have the same dimension. The numerical testing also showed the WENO-NP3 scheme cannot preserve the ENO property. Hence, the third-order WENO-NP3 scheme constructed in [13] is unavailable in applications.

3. THE NEW LOCAL SMOOTHNESS INDICATORS

In this section, we present a new way to calculate the indicators, which are used to replace (7) to improve the third-order WENO scheme.

First, the first-order derivatives of flux $f$ at node $i+1$, $i$ and $i-1$ are approximated by

$$
\begin{align*}
F'_{i+1} &= \frac{3f_{i+1} + f_{i-1} - 4f_i}{2\Delta x} \\
F'_{i} &= \frac{f_{i+1} - 2f_{i-1}}{2\Delta x} \\
F'_{i-1} &= \frac{4f_i - 3f_{i-1} - f_{i+1}}{2\Delta x}
\end{align*}
$$

(21)

Then, the new indicators are constructed as follows

$$
\begin{align*}
IS_0 &= (|F'_{i+1}\Delta x| - |F'_{i-1}\Delta x|)^2 = \frac{1}{4}(|f_{i+1} - f_{i-1}|^2 - |4f_i - 3f_{i-1} - f_{i+1}|^2) \\
IS_1 &= (|F'_{i-1}\Delta x| - |F'_{i+1}\Delta x|)^2 = \frac{1}{4}(|f_{i+1} - f_{i-1}|^2 - |3f_{i+1} + f_{i-1} - 4f_i|^2)
\end{align*}
$$

(22)

Next, we discuss the important properties of the new indicators.

Firstly, in monotonic smooth regions, the derivatives $F'_{i-1}$, $F'_{i}$ and $F'_{i+1}$ in (21) have the same sign, thus (22) becomes

$$
\begin{align*}
IS_0 &= \frac{1}{4} [f_{i+1} - f_{i-1} - (4f_i - 3f_{i-1} - f_{i+1})]^2 = (f_{i+1} - 2f_i + f_{i-1})^2 \\
IS_1 &= \frac{1}{4} [f_{i+1} - f_{i-1} - (3f_{i+1} + f_{i-1} - 4f_i)]^2 = (f_{i+1} - 2f_i + f_{i-1})^2
\end{align*}
$$

(23)

Hence we have

$$
IS_0 = IS_1
$$

(24)

Substituting (24) into the weight function (5) of the third-order WENO scheme, there is

$$
\omega_0 = c_0, \quad \omega_1 = c_1
$$

(25)

That means, in monotonic smooth regions, the new weights are exactly equal to the optimal weights, i.e. the fully third-order upstream scheme is recovered. Hence the new indicators (22) do not bring any additional dissipation for the WENO scheme. This property makes the new WENO scheme superior to the others in reducing numerical dissipation.

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Secondly, if there is a discontinuity inside a sub-stencil, for example, sub-stencil $S_0 = (i - 1, i)$, and the other one $S_1 = (i, i + 1)$ is smooth. Let

$$\begin{cases}
\beta_0 = (f_i - f_{i-1}) \\
\beta_1 = (f_{i+1} - f_i)
\end{cases}$$

(26)

Since $|\beta_0|$ is much larger than $|\beta_1|$, i.e. $|\beta_0| \gg |\beta_1|$, the new indicators of (22) can be approximated by

$$\begin{cases}
IS_0 = \frac{1}{4}(|\beta_0 + \beta_1| - |3\beta_0 - \beta_1|^2) = (\beta_0 + \beta_1)^2 = O(\beta_0^2) \\
IS_1 = \frac{1}{4}(|\beta_0 + \beta_1| - |3\beta_1 - \beta_0|^2) = (2\beta_1)^2 = O(\beta_1^2)
\end{cases}$$

(27)

(27) shows that the new indicators approximate the ones of Jiang and Shu (7). Hence, the ENO property can be maintained for problems with discontinuities.

Finally, at a critical point, $f'_i = 0$, the Taylor series expansions of (22) give

$$\begin{cases}
IS_0 = (\frac{f'''_i \Delta x^3}{6} + O(\Delta x^5)) + |f''_i \Delta x^2 - \frac{f'''_i \Delta x^3}{6} + O(\Delta x^4)|^2 \\
IS_1 = (\frac{f'''_i \Delta x^3}{6} + O(\Delta x^5)) + |f''_i \Delta x^2 + \frac{f'''_i \Delta x^3}{6} + O(\Delta x^4)|^2
\end{cases}$$

(28)

Similarly, (28) can be expressed as

$$IS_k = B(1 + O(\Delta x))$$

(29)

where $B$ is independent of $k$. Although the weights with the new indicators (22) do not meet the sufficient condition (11) near the critical points, owing to the advantage in monotonic smooth regions, they can still greatly decrease the numerical dissipation of the third-order WENO scheme, even near critical points.

4. NUMERICAL EXAMPLES

In this section, several linear advection problems, one-dimensional (1-D) and two-dimensional (2-D) Euler problems are calculated to test the performance of the new third-order WENO scheme. The time derivative is approximated with the third-order TVD Runge-Kutta method[14].

4.1. Linear advection problems

In the following we test the accuracy of WENO schemes on the linear advection equation

$$\begin{cases}
\frac{\partial u}{\partial t} + \frac{\partial u}{\partial x} = 0, & x_0 \leq x \leq x_1 \\
u(x, t = 0) = u_0(x), & \text{periodic boundary}
\end{cases}$$

(30)

The exact solution of (30) is given by

$$u(x, t) = u_0(x - t)$$

(31)
4.1.1. Case 1

This particular initial condition \([5]\) has two critical points at which \(f' = 0\) and \(f'' \neq 0\). The errors and the convergence order in \(L_{\infty}\) and \(L_1\) norm of the different schemes at \(t = 2\) with \(\Delta t = 8\Delta x^{5/3}\) are given in Table I. It can be seen that, all improved third-order WENO schemes can not achieve the expected third-order accuracy, but the present one performs best in all tested schemes, the \(L_1\) error of the present scheme is only half of the one of WENO-N3. Fig.1 displays the behavior of the weight \(\omega_0\) of different WENO schemes at initial step \(t = 0\) with a grid of 100 and shows the weight calculated by present local smoothness indicators is the same as the value of the ideal weight in monotonic smooth regions.
4.1.2. Case 2

\[ u_0(x) = \begin{cases} 
1, & \text{if } |x| < 0.5 \\
0, & \text{otherwise} 
\end{cases} \]  

(33)

This case is used to test the shock capturing capability of the WENO schemes. The numerical solution at \( t=8 \) with \( N=200 \) and time step \( \Delta t = \Delta x/2 \) are shown in Figs.2 and 3. It can be seen that, as mentioned previously, the WENO-NP3 scheme[13] generates apparent oscillations near discontinuous regions. The present scheme keeps the best behavior.

4.1.3. Case 3

\[ u_0(x) = e^{-(x-90)^2/400} \left( \cos \left( \frac{\pi}{8} (x - 90) \right) + \cos \left( \frac{\pi}{4} (x - 90) \right) \right), \quad 50 \leq x \leq 130 \]  

(34)

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This case is used to test the performance of WENO schemes for a smooth solution with several critical points [15]. Fig. 4 shows the numerical solutions with grid number $N = 200$ and $\Delta t = \Delta x/2$ at $t = 160$. As mentioned above, although the present scheme cannot achieve optimal convergence order, it still can greatly decrease the dissipation of the WENO scheme, even near critical points.

This example is also calculated by the further improved schemes (WENO-M-type and WENO-Z-type schemes), which are based on the WENO scheme with the new local smoothness indicators (22) and by means of the mapping function (12) of the WENO-M scheme and the global smoothness indicator (16) of the WENO-Z scheme, respectively. The results shown in Fig. 5 indicate that the mentioned methods can also be directly used to improve the performance of the new scheme. Since the purpose of this paper is mainly to show the benefits from the new local smoothness indicators, both the further improved schemes are no longer discussed and compared in this paper.

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4.1.4. Case 4

\[
\begin{align*}
  u(x,0) = & \begin{cases} 
    \frac{1}{6}(G(x,\beta,z-\delta) + G(x,\beta,z+\delta) + 4G(x,\beta,z)), & -0.8 \leq x < -0.6 \\
    1, & -0.4 \leq x < -0.2 \\
    1 - |10(x-0.1)|, & 0 \leq x < 0.2 \\
    \frac{1}{6}(F(x,\alpha,a-\delta) + F(x,\alpha,a+\delta) + 4F(x,\alpha,a)), & 0.4 \leq x < 0.6 \\
    0, & \text{otherwise}
  \end{cases}
\end{align*}
\]

\[
G(x,\beta,z) = e^{-\beta(x-z)^2},
\]

\[
F(x,\beta,a) = \sqrt{\max(1-\alpha^2(x-a)^2,0)}
\]

This is a classical testing case, which contains a Gaussian, a triangle, a square-wave and a semi-ellipse. It is often used to test the robustness and dissipation of high order schemes. The constants are \(z = -0.7\), \(\delta = 0.005\), \(\beta = \log(2)/(36\delta^2)\), \(a = 0.5\) and \(\alpha = 10\). It is solved with \(N = 200\) and time step \(\Delta t = \Delta x/2\). Figs.6 and 7 display the results at \(t = 2\). Same as shown in the previous case, the present scheme resolves all four waves better than the others.

4.2. Nonlinear Burgers equation (Case 5)

The nonlinear Burgers equation with a periodic boundary condition

\[
\begin{align*}
  \frac{\partial u}{\partial t} + \frac{\partial (u^2/2)}{\partial x} = 0, & -1 \leq x \leq 1, \\
  u(x,t=0) = u_0(x), & \text{periodic boundary},
\end{align*}
\]

is solved as case 5. For the initial data \(u_0(x) = \frac{1}{2} + \sin(\pi x)\), the exact solution is smooth up to \(t = \frac{1}{\pi}\). We list the errors and orders in Table II. It can be seen that, the present scheme can decrease the errors by about 70% compared with the WENO-JS scheme and by about 40% compared with the other two schemes.

<table>
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<tr>
<th>(N)</th>
<th>(L_\infty) (order)</th>
<th>(L_1) (order)</th>
<th>(L_\infty) (order)</th>
<th>(L_1) (order)</th>
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<tr>
<td>50</td>
<td>8.193E-3 (—)</td>
<td>2.152E-3 (—)</td>
<td>6.219E-3 (—)</td>
<td>1.380E-3 (—)</td>
</tr>
<tr>
<td>100</td>
<td>3.377E-3 (1.28)</td>
<td>4.536E-4 (2.25)</td>
<td>2.527E-3 (1.30)</td>
<td>2.996E-4 (2.20)</td>
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<tr>
<td>200</td>
<td>1.298E-3 (1.38)</td>
<td>1.014E-4 (2.16)</td>
<td>1.954E-3 (1.53)</td>
<td>2.242E-4 (2.29)</td>
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<tr>
<td>400</td>
<td>4.890E-4 (1.41)</td>
<td>2.330E-5 (2.12)</td>
<td>3.266E-4 (1.50)</td>
<td>1.286E-5 (2.31)</td>
</tr>
<tr>
<td>800</td>
<td>1.835E-4 (1.41)</td>
<td>5.215E-6 (2.16)</td>
<td>1.156E-4 (1.50)</td>
<td>2.624E-6 (2.30)</td>
</tr>
</tbody>
</table>

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<thead>
<tr>
<th>(N)</th>
<th>(L_\infty) (order)</th>
<th>(L_1) (order)</th>
<th>(L_\infty) (order)</th>
<th>(L_1) (order)</th>
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<td>1.189E-4 (1.50)</td>
<td>2.720E-6 (2.29)</td>
<td>8.408E-5 (1.53)</td>
<td>1.614E-6 (2.43)</td>
</tr>
</tbody>
</table>
4.3. One-dimensional Euler problems

The governing equations are as follows

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x} = 0$$

(37)

where $U = (\rho, \rho u, E)^T$, $F(U) = (\rho u, \rho u^2 + p, u(E + p))^T$. $\rho$, $u$, $E$, $p$ are the density, velocity, total energy and pressure, respectively. For ideal gas, $E = p/(\gamma - 1) + \rho u^2/2$. $\gamma = 1.4$ is the ratio of specific heat. Time step is taken as

$$\Delta t = \frac{\sigma \Delta x}{\max_i (|u_i| + c_i)}$$

(38)

where, $\sigma$ is CFL number, and 0.5 is used for all four 1-D cases in this subsection. $c$, the speed of sound, is given by $c = \sqrt{\gamma p/\rho}$. The local LF flux-splitting method is used and the WENO reconstruction is carried out in local characteristic fields[11]. All the reference solutions were obtained by the WENO-JS scheme with 2000 points.

4.3.1. Case 6 The first 1-D case is Shu-Osher problem[14], its initial condition is given as follows

$$(\rho, u, p) = \begin{cases} 
(3.857143, 2.629369, 31/3) & -5 \leq x < -4 \\
(1 + 0.2 \sin(5x), 0, 1) & -4 \leq x \leq 5 
\end{cases}$$

(39)

Figs.8 and 9 give the density distribution at $t = 1.8$ by using $N = 800$. As shown in the scalar case 3, for the solution with high frequency, the present scheme shows apparent improvement over the others.

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4.3.2. Case 7 The second 1-D case is Sod problem, with initial condition

\[
(\rho, u, p) = \begin{cases} 
(0.125, 0, 0.1) & 0.5 \leq x \leq 1 \\
(1, 0, 1) & 0 \leq x < 0.5
\end{cases}
\] (40)

It is solved with \( N = 200 \) at \( t = 0.14 \). The results in Figs. 10 and 11 show that the new scheme can capture shocks and contact discontinuities well.

4.3.3. Case 8 The third 1-D case is Lax problem[7], with initial condition

\[
(\rho, u, p) = \begin{cases} 
(0.445, 0.698, 3.528) & -5 \leq x < 0 \\
(0.5, 0, 0.571) & 0 \leq x \leq 5
\end{cases}
\] (41)

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It is solved with $N = 200$ at $t = 1.3$. The density distribution is displayed in Figs. 12 and 13. The new scheme obtains sharper shock profiles than the others.

### 4.3.4. Case 9

The last one is the blast-waves interaction problem[7], with initial condition

$$
(\rho, \ u, \ p) = \begin{cases}
(1, \ 0, \ 1000) & 0 \leq x < 0.1 \\
(1, \ 0, \ 0.001) & 0.1 \leq x < 0.9 \\
(1, \ 0, \ 100) & 0.9 \leq x < 1
\end{cases}
$$

The numerical results at $t = 0.038$ with $N = 800$ are presented in Figs. 14 and 15. As previously discussed, the new scheme can obtain more accurate solutions near extrema.

### 4.4. Two-dimensional Euler problems

In this subsection, we further study the shock capturing ability and numerical dissipation of the new scheme in the 2-D problems. The 2-D Euler equations are solved,

$$\frac{\partial U}{\partial t} + \frac{\partial F}{\partial x} + \frac{\partial G}{\partial y} = 0$$

where the conserved variables $U$ and the inviscid flux vectors $F$ and $G$ are

$$
U = \begin{bmatrix}
\rho \\
\rho u \\
\rho v \\
E
\end{bmatrix}, \quad F = \begin{bmatrix}
\rho u \\
\rho u^2 + p \\
\rho uv \\
E u + pu
\end{bmatrix}, \quad G = \begin{bmatrix}
\rho v \\
\rho uv \\
\rho v^2 + p \\
E v + pv
\end{bmatrix}
$$

The energy is given by

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

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The Roe[16] flux-splitting method is used for the inviscid convective fluxes, and the time step is taken as follows:

\[
\Delta t = \sigma \frac{\Delta t_x \Delta t_y}{\Delta t_x + \Delta t_y}
\]

\[
\begin{align*}
\Delta t_x &= \frac{\Delta x}{\max_{i,j} (|u_{i,j}| + c_{i,j})} \\
\Delta t_y &= \frac{\Delta y}{\max_{i,j} (|v_{i,j}| + c_{i,j})}
\end{align*}
\]

(46)

where, \(\sigma\) is CFL number, and 0.5 is used for all the tests below.

4.4.1. Case 10

This is a 2-D periodic vortex propagation problem[17, 18] used to assess the numerical dissipation of different schemes. The vortex is described as a perturbation to the velocity

Figure 16. Density contours of Case 10, 30 contours from 0.63 to 0.99
(u, v), temperature (T = p/ρ) and entropy (S = p/ρ) of the mean flow (ρ, u, v, p) = (1, 1, 1, 1)

\[
\begin{align*}
\hat{u} &= -\frac{\epsilon e^{(1-r^2)/2}}{2\pi} \tilde{y} \\
\hat{v} &= \frac{\epsilon e^{(1-r^2)/2}}{2\pi} \tilde{x} \\
\hat{T} &= -\frac{(\gamma - 1)\epsilon^2 e^{(1-r^2)}}{8\gamma\pi^2} \\
\hat{S} &= 0
\end{align*}
\]

(47)

where, (\tilde{x}, \tilde{y}) = (x - 5, y - 5), r^2 = \tilde{x}^2 + \tilde{y}^2, and the vortex strength \( \epsilon = 5.0 \). The computational domain is \([0, 10] \times [0, 10]\), and periodic boundary condition is used in both directions. It can be readily verified that the Euler equations with the above initial conditions admit an exact solution that moves with the speed (1, 1) in the diagonal direction. The errors based on the density of the WENO-M, WENO-N3 and the present schemes are presented in Table III. The present scheme reduces errors by about 70% compared with WENO-JS and by about 45% compared with the other two schemes.

The results are also carried out until \( t = 100 \) with a grid of \( 80 \times 80 \), to further demonstrate the low dissipation of the new scheme. Fig.16 shows the density contours of the WENO schemes. Figs.17 and 18 show the pressure distribution along x-axis at \( y = 5 \) and the evolution of kinetic energy \( E_k \), respectively. The kinetic energy is calculated by \( E_k = \sum_{i,j=1}^{N,M} \rho_{i,j} (u_{i,j}^2 + v_{i,j}^2) / (2N \times M) \), where \( N \) and \( M \) are the total cells in \( x \) and \( y \) direction of the grid, respectively. From these figures, it can be seen that, the WENO-JS scheme is too dissipative, while the present scheme has lowest dissipation.

<table>
<thead>
<tr>
<th>Grid</th>
<th>WENO-JS</th>
<th>WENO-M</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>( L_\infty ) (order)</td>
<td>( L_1 ) (order)</td>
</tr>
<tr>
<td>20 \times 20</td>
<td>4.153E-1 ( — )</td>
<td>2.083E-2 ( — )</td>
</tr>
<tr>
<td>40 \times 40</td>
<td>1.834E-1 (1.18)</td>
<td>9.525E-3 (1.29)</td>
</tr>
<tr>
<td>80 \times 80</td>
<td>6.652E-2 (1.46)</td>
<td>2.937E-3 (1.70)</td>
</tr>
<tr>
<td>160 \times 160</td>
<td>4.165E-2 (0.66)</td>
<td>1.084E-3 (1.44)</td>
</tr>
<tr>
<td>320 \times 320</td>
<td>1.964E-2 (1.08)</td>
<td>2.965E-4 (1.87)</td>
</tr>
</tbody>
</table>

<table>
<thead>
<tr>
<th>Grid</th>
<th>WENO-N3</th>
<th>Present</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td>( L_\infty ) (order)</td>
<td>( L_1 ) (order)</td>
</tr>
<tr>
<td>20 \times 20</td>
<td>3.944E-1 ( — )</td>
<td>1.927E-2 ( — )</td>
</tr>
<tr>
<td>40 \times 40</td>
<td>1.527E-1 (1.37)</td>
<td>6.746E-3 (1.51)</td>
</tr>
<tr>
<td>80 \times 80</td>
<td>7.060E-2 (1.11)</td>
<td>1.972E-3 (1.77)</td>
</tr>
<tr>
<td>160 \times 160</td>
<td>3.453E-2 (1.03)</td>
<td>5.732E-4 (1.78)</td>
</tr>
<tr>
<td>320 \times 320</td>
<td>1.020E-2 (1.76)</td>
<td>1.130E-4 (2.34)</td>
</tr>
</tbody>
</table>

4.4.2. Case II The 2-D shock/vortex interaction problem describes an interaction between a stationary shock and a vortex[1]. The computational domain is taken to be \([0, 2] \times [0, 1]\). A stationary Mach 1.1 shock is positioned at \( x = 0.5 \) and normal to the \( x \)-axis. Its left state is

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Figure 17. Pressure distributed along $y = 5$ of Case 10

Figure 18. Evolution of $E_k$ of Case 10

$(\rho, u, v, p) = (1, 1.1\sqrt{\gamma}, 0, 1)$. A small vortex is superimposed to the flow on the left of the shock and centered at $(x_c, y_c) = (0.25, 0.5)$. The vortex is described as a perturbation to the velocity $(u, v)$, temperature $(T = p/\rho)$ and entropy $S = \ln(p/\rho\gamma)$ of the mean flow

\[
\begin{align*}
\hat{u} &= \epsilon \tau e^{a(1-\tau^2)} \sin \theta \\
\hat{v} &= -\epsilon \tau e^{a(1-\tau^2)} \cos \theta \\
\hat{T} &= -\frac{(\gamma - 1)\epsilon^2 e^{2a(1-\tau^2)}}{4a\gamma} \\
\hat{S} &= 0
\end{align*}
\] (48)

where, $\tau = r/r_c$, $r = \sqrt{(x-x_c)^2 + (y-y_c)^2}$, $\epsilon = 0.3$, $r_c = 0.05$ and $a = 0.204$ are the same as in Refs.[1, 19]. The upper and lower boundaries are set to be reflective. The results at $t = 0.6$ with a grid of $100 \times 50$ are solved. Fig. 19 is the pressure distribution along the center line of $y = 0.5$. The result obtained by WENO-JS with a refined mesh of $2000 \times 1000$ is given as the reference solution. It can be seen that the present scheme obtained most accurate solution near the center of the vortex.

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4.4.3. Case 12 This is a 2-D Riemann problem[20] used to test the shock capturing capability of the present scheme for calculating shock/shock interaction problems. It is solved on a unit square with initial condition

\[
(p, u, v, p) = \begin{cases} 
(1.0, -0.75, -0.5, 1.0), & 0.5 \leq x \leq 1, \ 0.5 \leq y \leq 1 \\
(2.0, -0.75, 0.5, 1.0), & 0 \leq x < 0.5, \ 0.5 \leq y \leq 1 \\
(1.0, 0.75, 0.5, 1.0), & 0 \leq x < 0.5, \ 0 \leq y < 0.5 \\
(3.0, 0.75, -0.5, 1.0), & 0.5 \leq x \leq 1, \ 0 \leq y < 0.5 
\end{cases}
\]  

Figure 19. Pressure distributed along \(y = 0.5\) of Case 11
The solution at $t = 0.23$ is solved with a grid of $400 \times 400$. The reference is calculated by WENO-JS with a grid of $800 \times 800$. Fig. 20 shows the density contours of the reference solution and the results of different WENO schemes. As the figure shows, the present scheme obtained rich structures similar as those obtained by WENO-JS with refined meshes.
5. CONCLUSION REMARKS

In this paper, a new method is proposed to calculate the local smoothness indicators for the third-order WENO scheme by using all the three points on the global stencil. Since the two new indicators are exactly the same in monotonic smooth regions, the weights of the new WENO scheme are equal to the optimal weights, and this property makes the new scheme recover the fully third-order upstream scheme. The numerical results demonstrate that the WENO scheme with the new indicators has less dissipation and better resolution than the ones of Jiang and Shu for both smooth and discontinuous solutions.

Although this paper only focuses on the advantages brought by the new indicators, those methods, such as the mapping function used in the WENO-M scheme and the global smoothness indicator used in the WENO-Z scheme, can also directly be used to further improve the performance of the new third-order WENO scheme.

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