



Numerical Solution of High-Dimensional Shockwave Equations by Bivariate Multi-Quadric Quasi-Interpolation

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Abstract: Radial basis function-based quasi-interpolation performs efficiently in high-dimensional approximation and its applications, which can attain the approximant and its derivatives directly without solving any large-scale linear system. In this paper, the bivariate multi-quadrics (MQ) quasi-interpolation is used to simulate two-dimensional (2-D) Burgers' equation. Specifically, the spatial derivatives are approximated by using the quasi-interpolation, and the time derivatives are approximated by forward finite difference method. One advantage of the proposed scheme is its simplicity and easy implementation. More importantly, the proposed scheme opens the gate to meshless adaptive moving knots methods for the high-dimensional partial differential equations (PDEs) with shock or soliton waves. The scheme is also applicable to other non-linear high-dimensional PDEs. Two numerical examples of Burgers' equation (shock wave equation) and one example of the Sine–Gordon equation (soliton wave equation) are presented to verify the high accuracy and efficiency of this method.

Keywords: high-dimensional PDE; shockwave; quasi-interpolation; meshless method

MSC: 41A25; 65D05; 65M06; 65M12

1. Introduction

Studying the numerical solution of high-dimensional non-linear Burgers' equation is of great significance. One reason is that it is a fundamental shockwave PDE from fluid mechanics which often occurs in applied and computational mathematics, such as modeling of dynamics, heat conduction, and acoustic wave [1,2]. More importantly, it often plays the role of a testing equation to check the feasibility of the scheme, i.e., once the scheme is efficient for Burgers' equation, there is great possibility that it is applicable to other equations with shock or soliton waves, such as non-linear Schrodinger equation [3], Sine–Gordon equation [4], and Klein–Gordon equation [5].

In this paper, the proposed scheme is applicable to other non-linear high-dimensional PDEs. Burgers' equation has been studied by many researchers not only because its shockwave behavior when the coefficient *R* is large, but also because it is the simplest form of non-linear advection equation. Hence we take the 2-D coupled Burgers' equation as one example:



$$u_t + uu_x + vu_y = \frac{1}{R}(u_{xx} + u_{yy}), v_t + uv_x + vv_y = \frac{1}{R}(v_{xx} + v_{yy}).$$
(1)

The initial conditions are:

$$u(x, y, 0) = f_1(x, y), (x, y) \in D, v(x, y, 0) = f_2(x, y), (x, y) \in D,$$
(2)

and the boundary conditions are:

$$u(x, y, t) = g_1(x, y, t), (x, y) \in \partial D, v(x, y, t) = g_2(x, y, t), (x, y) \in \partial D,$$
(3)

where $t \in [0, T]$, $D \subset \mathbb{R}^2$ is a bounded domain and ∂D is the boundary. u(x, y, t) and v(x, y, t) are the coupled solutions to be determined, $f_1(x, y)$, $f_2(x, y)$, $g_1(x, y, t)$ and $g_2(x, y, t)$ are given functions, and R is the Reynolds number. Fletcher [6] obtained its exact solutions, so that one could evaluate the quality of numerical solutions.

During the past decades, various numerical techniques have been developed for the solution of 2-D Burgers' equation. Bahadir [7] introduced one fully implicit finite difference scheme, then Zhu et al. [8] proposed the discrete Adomian decomposition method (ADM), Yu et al. [9] applied the bivariate quintic spline to solve numerically.

Compared with these mesh-based methods, which depend on a topology shape of the mesh, the meshless method performs satisfactorily for problems with complicated and irregular geometries [10]. In recent years, the meshless collocation method has drawn considerable attention for solving PDEs [11–13]. Mittal and Tripathi [14] proposed a collocation method based on Modified bi-cubic B-Spline functions. However, it requires to solve the inverse of a large-scale matrix on each time step, which costs computational time and might be ill-conditioned.

Considering the problems mentioned above, we try to apply the MQ quasi-interpolation method for the numerical solution of PDEs because of the following reasons:

- Quasi-interpolation could give the approximant and its derivatives directly needing no solution of any large-scale system, hence it is easy to be implemented by the engineers in applications.
- Quasi-interpolation could filter the noise of the data which often occurs in applications because of various reasons, e.g., measure error, data pollution, and so forth.
- Ma [15] proved the superiority of MQ quasi-interpolation over finite difference and other interpolation methods in both stability and accuracy, especially when approximating scattered data.

MQ function was firstly introduced by Hardy [16], then it has been investigated extensively [17–19] and successfully applied to solve one dimensional (1-D) PDEs [12,20–23]. Due to the extraordinary performance in 1-D, Ling [24] extended the univariate MQ quasi-interpolation to 2-D. Feng and Zhou [25], Wu et al. [26] and Wu et al. [27,28] constructed high-dimensional quasi-interpolations.

The ambition of this paper is to construct a numerical scheme for solving 2-D Burgers' equation applying the bivariate MQ quasi-interpolation. Firstly, we develop a bivariate MQ quasi-interpolation and analyze its approximation order to a given function and its high order derivatives. Then the numerical scheme for solving the 2-D coupled Burgers' equation is proposed applying the bivariate MQ quasi-interpolation. In the numerical scheme, the spatial derivatives are approximated by using the derivatives of the quasi-interpolation, and the time derivatives of the dependent variables are approximated by the forward finite difference method. The truncation error and the total error of our scheme are estimated to show the accuracy of the proposed scheme. At last, two numerical examples of Burgers' equation (shockwave equation) and one example of the Sine–Gordon equation (soliton wave equation) are presented to verify the efficiency of the proposed method.

One advantage of the proposed scheme is its simplicity, easy implementation, but high accuracy and stability. More importantly, since the proposed method requires no triangulation of the region, adaptive refinement can be added in the scheme directly. In other words, it allows more flexible knots movement and one need not to worry the problems of mesh over gathering or condition numbers. In addition, the proposed scheme is applicable to other high-dimensional PDEs with shock or even soliton waves.

The rest of the paper is organized as follows. In Section 2, the bivariate MQ quasi-interpolant is introduced. In Section 3, we present the numerical scheme to solve 2-D Burgers' equation and give the error estimations. In Section 4, three numerical examples are proposed to verify our method. Section 5 concludes the whole paper.

2. Bivariate MQ Quasi-Interpolations

In this section, we firstly introduce the univariate MQ quasi-interpolation. On the basis, we develop the bivariate MQ quasi-interpolation and show its approximation order to the function and its derivatives.

Given the data $\{x_j, f_j\}, f_j = f(x_j)$, the univariate MQ quasi-interpolation is

$$Qf(x) = \sum_{j} f_{j}\psi_{j}(x), \tag{4}$$

$$\psi_j(x) = \frac{\phi_{j+1}(x) - \phi_j(x)}{2(x_{j+1} - x_j)} - \frac{\phi_j(x) - \phi_{j-1}(x)}{2(x_j - x_{j-1})},\tag{5}$$

 $\phi_j(x) = \sqrt{(x - x_j)^2 + c_1^2}$ is named MQ function, c_1 is shape parameter, which is often a small constant. As c_1 goes to zero, $\phi_j(x)$ will converge to $|x - x_j|$, consequently the univariate MQ interpolation could be taken as piecewise linear interpolation but possesses smoothness property. Denoting by $h_1 = \max_j(x_{j+1} - x_j)$, the approximation error of Qf(x) to the original function f(x) and the *k*-th derivatives were given in [19]:

Theorem 1. If $f(x) \in \mathbb{C}^{(k+2)}(\mathbb{R})$ and $f^{(j)}(x)$ is bounded by a polynomial of degree k + 2 - j, then

$$|Qf^{(k)}(x) - f^{(k)}(x)| \le \mathcal{O}(h_1^{\frac{2}{k+1}})$$
(6)

holds, provided that $c_1 = \mathcal{O}(h_1^{\frac{1}{k+1}})$.

Besides the high accuracy, the MQ quasi-interpolation has been proven to possess many other favorable properties, such as monotonicity preservation, shape preservation, and variation diminishing [17,18]. Because of these positive properties, researchers began to extend the univariate quasi-interpolation to multi-dimension [24–26].

In this paper, we propose one kind of bivariate MQ quasi-interpolation using the tensor product technique. Specifically, given data (x_i, y_j, f_{ij}) and $f_{ij} = f(x_i, y_j)$, the bivariate MQ quasi-interpolation is defined as:

$$(Qf)(x,y) = \sum_{i} \sum_{j} f_{ij} \psi_i(x) \psi_j(y),$$
(7)

where $\psi_i(x)$ are described in (5) and $\psi_i(y)$ are represented as follows:

$$\begin{split} \phi_j(y) &= \sqrt{(y-y_j)^2 + c_2^2}, \\ \psi_j(y) &= \frac{\phi_{j+1}(y) - \phi_j(y)}{2(y_{j+1} - y_j)} - \frac{\phi_j(y) - \phi_{j-1}(y)}{2(y_j - y_{j-1})}, \end{split}$$

 c_2 is a small constant. Denoting $h_2 = \max_j (y_{j+1} - y_j)$, the error estimation of the proposed bivariate MQ quasi-interpolation (7) is given in Theorem 2.

Theorem 2. If $f(x,y) \in \mathbb{C}^{(k+2)}(\mathbb{R}^2)$ and $\frac{\partial (Qf)^{i+j}(x,y)}{\partial x^i \partial y^j}$ is bounded by a polynomial $p_1(x)p_2(y)$, $p_1(x)$ and $p_2(y)$ are polynomials of degree $\alpha_1 + 2 - i$ and $\alpha_2 + 2 - j$ respectively $(\alpha_1, \alpha_2, \alpha_1 + 2 - i, and \alpha_2 + 2 - j are all nonnegative integers)$, then the error of bivariate MQ quasi-interpolation (7) satisfies

$$\left\|\frac{\partial(Qf)^{\alpha_1+\alpha_2}(x,y)}{\partial x^{\alpha_1}\partial y^{\alpha_2}} - \frac{\partial f^{\alpha_1+\alpha_2}(x,y)}{\partial x^{\alpha_1}\partial y^{\alpha_2}}\right)\right\|_{\infty} \leqslant \mathcal{O}(h_1^{\frac{2}{\alpha_1+1}} + h_2^{\frac{2}{\alpha_2+1}}),\tag{8}$$

provided that $c_1 = O(h_1^{\frac{1}{\alpha_1+1}})$ and $c_2 = O(h_2^{\frac{1}{\alpha_2+1}})$.

Proof. By adding one middle term, the following inequalities are obtained:

$$\| (Qf)(x,y) - f(x,y) \|_{\infty}$$

$$= \| \sum_{i} \sum_{j} f_{ij} \psi_{i}(x) \psi_{j}(y) - \sum_{i} f(x_{i},y) \psi_{i}(x) + \sum_{i} f(x_{i},y) \psi_{i}(x) - f(x,y) \|_{\infty}$$

$$\le \| \sum_{i} \psi_{i}(x) (\sum_{j} f_{ij} \psi_{j}(y) - f(x_{i},y)) \|_{\infty} + \| \sum_{i} f(x_{i},y) \psi_{i}(x) - f(x,y) \|_{\infty} .$$

Since $\sum_j f_{ij}\psi_j(y)$ and $\sum_i f(x_i, y)\psi_i(x)$ are the approximation of $f(x_i, y)$ and f(x, y) separately, when $c_1 = O(h_1)$ and $c_2 = O(h_2)$, they are estimated as:

$$\| \sum_{j} f_{ij} \psi_{j}(y) - f(x_{i}, y) \|_{\infty} = \mathcal{O}(h_{2}^{2}), \\ \| \sum_{i} f(x_{i}, y) \psi_{i}(x) - f(x, y) \|_{\infty} = \mathcal{O}(h_{1}^{2}).$$

In addition, $|\sum_{i} \psi_{i}(x)| \leq 1$, the result are as follows:

$$\| Qf(x,y) - f(x,y) \|_{\infty} \le | \sum_{i} \psi_{i}(x) | \mathcal{O}(h_{1}^{2}) + \mathcal{O}(h_{2}^{2}) \le \mathcal{O}(h_{1}^{2} + h_{2}^{2}).$$

Similarly, the derivative errors of $(Qf)^{(k)}$ to $f^{(k)}$ could be estimated based on Theorem 1. This ends the proof. \Box

3. Numerical Scheme Applying Bivariate MQ Quasi-Interpolant

In this section, we develop a numerical scheme for solving 2-D Burgers' Equations (1)–(3) based on bivariate MQ quasi-interpolation in Section 3.1. Then we provide the error estimations of Algorithm 1 in Section 3.2.

3.1. The Main Algorithm

We represent the approximations of u(x, y, t) and v(x, y, t) at the knots $(x_i, y_j, t_k) = (ih_1, jh_2, k\tau)$ by u_{ij}^k and v_{ij}^k respectively, where h_1 and h_2 are the mesh sizes in x and y direction separately, τ is the time step sizes, i.e., $u_{ij}^k \approx u(x_i, y_j, t_k)$, $v_{ij}^k \approx v(x_i, y_j, t_k)$. We use the similar expressions for the derivative

approximations, e.g., $(u_x)_{ij}^k \approx u_x(x_i, y_j, t_k)$, $(u_y)_{ij}^k \approx u_y(x_i, y_j, t_k)$ and so forth. The numerical solution could be realized in the following Algorithm 1.

3.2. Error Estimations of Algorithm 1

In Section 3.2, the truncation error of Algorithm 1 is proposed in Lemma 1. Then based on Lemma 1, the total error of Algorithm 1 is provided in Theorem 3.

Algorithm 1 Numerical scheme of 2-D Burgers' equation applying the bivariate MQ quasi-interpolation.

Input: u_{ij}^k, v_{ij}^k Output: $u_{ij}^{k+1}, v_{ij}^{k+1}$

1: Approximate the solutions' (*u* and *v*) first and second space derivatives by using the bivariate MQ quasi-interpolation. For example, the first derivatives of function *u* could be approximated as follows,

$$(u_x)_{ij}^k = \sum_m \sum_n u_{ij}^k \psi'_m(x_i) \psi_n(y_j),$$

$$(u_y)_{ij}^k = \sum_m \sum_n u_{ij}^k \psi_m(x_i) \psi'_n(y_j).$$

the other derivatives can be approximated similarly based on the bivariate MQ quasi-interpolation.

2: Discrete the Burgers' equation in time and substitute the points (x_i, y_j) into the derivatives, compute u_{ij}^{k+1} and v_{ij}^{k+1} :

$$u_{ij}^{k+1} = u_{ij}^k - \tau u_{ij}^k (u_x)_{ij}^k - \tau v_{ij}^k (u_y)_{ij}^k + \frac{\tau}{R} ((u_{xx})_{ij}^k + (u_{yy})_{ij}^k),$$

$$v_{ij}^{k+1} = v_{ij}^k - \tau u_{ij}^k (v_x)_{ij}^k - \tau v_{ij}^k (v_y)_{ij}^k + \frac{\tau}{R} ((v_{xx})_{ij}^k + (v_{yy})_{ij}^k).$$

3: Return to Step 1.

For the error estimation, we denote the following operators L_1 , L_2 , $L_{1,d}$, $L_{2,d}$ as

$$\begin{split} &L_{1}(u,v) = u_{t} + uu_{x} + vu_{y} - \frac{1}{R}(u_{xx} + u_{yy}), \\ &L_{2}(u,v) = v_{t} + uv_{x} + vv_{y} - \frac{1}{R}(v_{xx} + v_{yy}), \\ &L_{1,d}(u_{ij}^{k}, v_{ij}^{k}) = \frac{u_{ij}^{k+1} - u_{ij}^{k}}{\tau} + u_{ij}^{k}(u_{x})_{ij}^{k} + v_{ij}^{k}(u_{y})_{ij}^{k} - \frac{1}{R}((u_{xx})_{ij}^{k} + (u_{yy})_{ij}^{k}), \\ &L_{2,d}(u_{ij}^{k}, v_{ij}^{k}) = \frac{v_{ij}^{k+1} - v_{ij}^{k}}{\tau} + u_{ij}^{k}(v_{x})_{ij}^{k} + v_{ij}^{k}(v_{y})_{ij}^{k} - \frac{1}{R}((v_{xx})_{ij}^{k} + (v_{yy})_{ij}^{k}). \end{split}$$

Lemma 1. For any $t \in [0, T]$, if $u(x, t) \in \mathbb{C}^4(\mathbb{R})$ and $v(x, t) \in \mathbb{C}^4(\mathbb{R})$, the truncation error of Algorithm 1 is

$$\begin{aligned} & (R_1)_{ij}^k = L_{1,d}(u_{ij}^k, v_{ij}^k) - L_1(u, v)(x_i, y_j, t_k) = \mathcal{O}(\tau + h_1^{2/3} + h_2^{2/3}), \\ & (R_2)_{ij}^k = L_{2,d}(u_{ij}^k, v_{ij}^k) - L_2(u, v)(x_i, y_j, t_k) = \mathcal{O}(\tau + h_1^{2/3} + h_2^{2/3}). \end{aligned}$$

Proof. Firstly, based on the Taylor expansion, we have:

$$\frac{\frac{u(x_i, y_j, t_{k+1}) - u(x_i, y_j, t_k)}{\tau}}{\frac{v(x_i, y_j, t_{k+1}) - v(x_i, y_j, t_k)}{\tau}} = u_t(x_i, y_j, t_k) + \mathcal{O}(\tau),$$

Then based on Theorem 2, we get the following error estimations:

$$\begin{split} & u(x_i, y_j, t_k)(u_x)_{ij}^k + v(x_i, y_j, t_k)(u_y)_{ij}^k \\ &= u(x_i, y_j, t_k)(u_x(x_i, y_j, t_k) + \mathcal{O}(h_1)) + v(x_i, y_j, t_k)(u_y(x_i, y_j, t_k) + \mathcal{O}(h_2)), \\ & u(x_i, y_j, t_k)(v_x)_{ij}^k + v(x_i, y_j, t_k)(v_y)_{ij}^k \\ &= u(x_i, y_j, t_k)(v_x(x_i, y_j, t_k) + \mathcal{O}(h_1)) + v(x_i, y_j, t_k)(v_y(x_i, y_j, t_k) + \mathcal{O}(h_2)), \end{split}$$

and

$$\begin{aligned} (u_{xx})_{ij}^k + (u_{yy})_{ij}^k &= u_{xx}(x_i, y_j, t_k) + u_{yy}(x_i, y_j, t_k) + \mathcal{O}(h_1^{\frac{2}{3}} + h_2^{\frac{2}{3}}), \\ (v_{xx})_{ij}^k + (v_{yy})_{ij}^k &= v_{xx}(x_i, y_j, t_k) + v_{yy}(x_i, y_j, t_k) + \mathcal{O}(h_1^{\frac{2}{3}} + h_2^{\frac{2}{3}}). \end{aligned}$$

Finally, we get the truncation errors:

$$\begin{split} (R_1)_{ij}^k &= L_{1,d}(u_{ij}^k, v_{ij}^k) - L_1(u, v)(x_i, y_j, t_k) \\ &= \mathcal{O}(\tau) + \mathcal{O}(h_1 + h_2) - \frac{1}{R}\mathcal{O}(h_1^{2/3} + h_2^{2/3}) \\ &= \mathcal{O}(\tau + h_1^{2/3} + h_2^{2/3}), \\ (R_2)_{ij}^k &= L_{2,d}(u_{ij}^k, v_{ij}^k) - L_2(u, v)(x_i, y_j, t_k) \\ &= \mathcal{O}(\tau) + \mathcal{O}(h_1 + h_2) - \frac{1}{R}\mathcal{O}(h_1^{2/3} + h_2^{2/3}) \\ &= \mathcal{O}(\tau + h_1^{2/3} + h_2^{2/3}). \end{split}$$

This ends the proof. \Box

Suppose $u_k(x, y)$ and $v_k(x, y)$ are the simulations of the exact solutions $u(x, y, t_k)$ and $v(x, y, t_k)$ using Algorithm 1, the total error of Algorithm 1 is provided in Theorem 3:

Theorem 3. Defining the total errors of Algorithm 1 as $e_u^k(x, y) = u_k(x, y) - u(x, y, t_k)$, $e_v^k(x, y) = v_k(x, y) - u(x, y, t_k)$. $v(x, y, t_k)$, if $u(x, y, t) \in \mathbb{C}^4(\mathbb{R})$, $v(x, y, t) \in \mathbb{C}^4(\mathbb{R})$, the error estimation of Algorithm 1 satisfies:

$$\begin{pmatrix} \|e_{u}^{K}\|_{2,\infty} \\ \|e_{v}^{K}\|_{2,\infty} \end{pmatrix} = \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}), K = T/\tau,$$
(9)

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where the Soblev norm for the function g(x, y) is defined as:

$$\|g\|_{2,\infty} = \max_{0 \le \alpha_1 + \alpha_2 \le 2} \|\frac{\partial^{\alpha_1 + \alpha_2}g}{\partial^{\alpha_1} x \partial^{\alpha_2} y}\|_{\infty}.$$

Proof. Since u(x, y, t), v(x, y, t) are the exact solutions of (1), they satisfy:

$$\frac{u(x, y, t_{k+1}) - u(x, y, t_k)}{\tau} = -u(x, y, t_k)u'_x(x, y, t_k) - v(x, y, t_k)u'_y(x, y, t_k)
+ \frac{1}{R}(u''_{xx}(x, y, t_k) + u''_{yy}(x, y, t_k)) + \mathcal{O}(\tau),
\frac{v(x, y, t_{k+1}) - v(x, y, t_k)}{\tau} = -u(x, y, t_k)v'_x(x, y, t_k) - v(x, y, t_k)v'_y(x, y, t_k)
+ \frac{1}{R}(v''_{xx}(x, y, t_k) + v''_{yy}(x, y, t_k)) + \mathcal{O}(\tau).$$
(10)

According to Lemma 1, $u_k(x, y)$, $v_k(x, y)$ satisfy

$$\frac{u_{k+1}(x,y) - u_k(x,y)}{\tau} = -u_k(x,y)(u_k)'_x(x,y) - v_k(x,y)(u_k)'_y(x,y)
+ \frac{1}{R}((u_k)''_{xx}(x,y) + (u_k)''_{yy}(x,y)) + \mathcal{O}(h_1^{2/3} + h_2^{2/3}),
\frac{v_{k+1}(x,y) - v_k(x,y)}{\tau} = -u_k(x,y)(v_k)'_x(x,y) - v_k(x,y)(v_k)'_y(x,y)
+ \frac{1}{R}m((v_k)''_{xx}(x,y) + (v_k)''_{yy}(x,y)) + \mathcal{O}(h_1^{2/3} + h_2^{2/3}).$$
(11)

Subtracting (10) from (11), one can get the following equation by adding some middle terms:

$$\frac{e_{u}^{k+1}(x,y) - e_{u}^{k}(x,y)}{\tau} = -e_{u}^{k}(x,y)(u_{k})_{x}'(x,y) - u_{k}(x,y)(e_{u}^{k})_{x}'(x,y)
- e_{v}^{k}(x,y)(u_{k})_{y}'(x,y) - v_{k}(x,y)(e_{u}^{k})_{y}'(x,y)
+ \frac{1}{R}((e_{u}^{k})_{xx}''(x,y) + (e_{u}^{k})_{yy}''(x,y)) + \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}),
\frac{e_{v}^{k+1}(x,y) - e_{v}^{k}(x,y)}{\tau} = -e_{u}^{k}(x,y)(v_{k})_{x}'(x,y) - u_{k}(x,y)(e_{v}^{k})_{x}'(x,y)
- e_{v}^{k}(x,y)(v_{k})_{x}'(x,y) - v_{k}(x,y)(e_{v}^{k})_{x}'(x,y)
+ \frac{1}{R}((e_{v}^{k})_{xx}''(x,y) + (e_{v}^{k})_{yy}''(x,y)) + \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}).$$
(12)

Therefore, the Soblev norm of the errors satisfy

$$\frac{\|e_{u}^{k+1}\|_{2,\infty} - \|e_{u}^{k}\|_{2,\infty}}{\tau} \leq C(\|e_{u}^{k}\|_{2,\infty} + \|e_{v}^{k}\|_{2,\infty}) + \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}),$$
$$\frac{\|e_{v}^{k+1}\|_{2,\infty} - \|e_{v}^{k}\|_{2,\infty}}{\tau} \leq C(\|e_{u}^{k}\|_{2,\infty} + \|e_{v}^{k}\|_{2,\infty}) + \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}),$$

where $C = 2max\{||u||_{1,\infty}, ||v||_{1,\infty}, \frac{1}{R}\}.$

Hence we have

$$\begin{pmatrix} \|e_u^{k+1}\|_{2,\infty} \\ \|e_v^{k+1}\|_{2,\infty} \end{pmatrix} = B \begin{pmatrix} \|e_u^k\|_{2,\infty} \\ \|e_v^k\|_{2,\infty} \end{pmatrix} + \mathcal{O}(\tau^2 + \tau h_1^{2/3} + \tau h_2^{2/3})$$

...
$$= B^{k+1} \begin{pmatrix} \|e_u^0\|_{2,\infty} \\ \|e_v^0\|_{2,\infty} \end{pmatrix} + (B^k + B^{k-1} + \dots + B^0) \mathcal{O}(\tau^2 + \tau h_1^{2/3} + \tau h_2^{2/3}),$$

where $B = \begin{pmatrix} 1 + C\tau & C\tau \\ C\tau & 1 + C\tau \end{pmatrix}$. B is a positive definite matrix, hence there exists a nonsingular matrix *X* satisfying

$$B = X^{-1} \left(\begin{array}{cc} 1 & 0 \\ 0 & 1 + 2C\tau \end{array} \right) X.$$

Since $||e_u^0||_{2,\infty} = 0$, $||e_v^0||_{2,\infty} = 0$, we can get

$$\begin{pmatrix} \|e_{u}^{K}\|_{2,\infty} \\ \|e_{v}^{K}\|_{2,\infty} \end{pmatrix} = X^{-1} \begin{pmatrix} K & 0 \\ 0 & \frac{1 - (1 + 2C\tau)^{K}}{-2C\tau} \end{pmatrix} X \mathcal{O}(\tau^{2} + \tau h_{1}^{2/3} + \tau h_{2}^{2/3})$$
$$= \mathcal{O}(\tau + h_{1}^{2/3} + h_{2}^{2/3}).$$

This completes the proof. \Box

Remark 1. In this section, we apply the bivariate MQ quasi-interpolation for solving the 2-D coupled Burgers' equation. As is shown in Ma [15], the MQ method is more accurate, more stable, and simpler for simulating the high order derivatives than finite difference method, especially for scatted data or the data with noise. In addition, the proposed method is simple and easy to implement, since one need not to solve any large-scale linear system and could get the original function and derivatives' approximation directly.

4. Numerical Examples

2-D Burgers' Equation

In this section, we select two presentative test examples to illustrate the performance of the method described in the previous section. There are several ways to process the boundary, here the popular bivariate operator which appears in Ling [24] is employed. All these experiments are carried out on a computer with an Intel (R) Core (TM) i7-7500 CPU, 3.20 GHZ processor and 8.00 GB RAM. To check the validity of the scheme, the root mean square (RMS) and L_{∞} error norms are applied to make comparisons with the previous methods [8]. The error norms are defined as

$$\begin{aligned} \text{RMS} &= \sqrt{\frac{\sum_i \sum_j (u_{ij}^{exact} - u_{ij}^{num})^2}{(N_1 + 1)(N_2 + 1)}}, \\ L_{\infty} &= \max_{ij} |u_{ij}^{exact} - u_{ij}^{num}|. \end{aligned}$$

where u_{ij}^{exact} and u_{ij}^{num} are the exact and numerical results of *u* at the knot (x_i, y_j) .

Example 1. In this problem, we choose the computational domain as $D = \{(x, y) : 0 \le x \le 0.5, 0 \le y \le 0.5\}$, and the spatial knots distribute uniformly on the computational domain with a mesh width $h_1 = h_2 = 0.025$. For the Burgers' Equations (1) and (2), the initial conditions at t = 0 is

$$f_1(x,y) = x + y, \qquad (x,y) \in D, f_2(x,y) = x - y, \qquad (x,y) \in D.$$
(13)

Under the above conditions, the exact solutions could be given [6]:

$$u(x,y,t) = \frac{x+y-2xt}{1-2t^2},$$
(14)

$$v(x,y,t) = \frac{x - y - 2yt}{1 - 2t^2}.$$
(15)

In this example, we observe the performance of our method with R = 100 for the convenience of comparison, and the numerical solutions are simulated using uniform knots, with a space width $h_1 = h_2 = 0.025$, the time step is $\tau = 10^{-4}$. The parameter is set as $c = 2.65 \times 10^{-6}$.

The study compares the errors of the presented method and Zhu et al.'s method [8]. The numerical errors at time t = 0.1 and t = 0.4 are listed in Tables 1 and 2 respectively. It is observed that at time t = 0.1, our errors are similar(even a little larger) than Zhu et al.'s method [8]. However, at time t = 0.4, our method is far smaller than Zhu et al.'s method [8]. That is to say, our method could maintain a longer time simulation.

Mesh Grid	Error of <i>u</i>	Error of <i>u</i> Error of <i>v</i>		
	Zhu et al. [8]	The Present	Zhu et al. [8]	The Present
(0.1, 0.1)	$3.3075 imes 10^{-6}$	$3.3081 imes 10^{-6}$	$1.0538 imes 10^{-6}$	1.0524×10^{-6}
(0.3, 0.1)	5.5616×10^{-6}	$7.6686 imes 10^{-6}$	$3.3077 imes 10^{-5}$	$7.5178 imes 10^{-6}$
(0.2, 0.2)	$6.6152 imes10^{-6}$	6.6161×10^{-6}	$2.1077 imes10^{-6}$	$2.1049 imes10^{-6}$
(0.4, 0.2)	8.8694×10^{-6}	1.0977×10^{-5}	2.2540×10^{-6}	$8.5703 imes10^{-6}$
(0.1, 0.3)	$7.6693 imes 10^{-6}$	5.8456×10^{-6}	7.5234×10^{-6}	3.3081×10^{-6}
(0.3, 0.3)	$9.9233 imes 10^{-6}$	9.9242×10^{-6}	3.1615×10^{-6}	$3.1573 imes 10^{-6}$
(0.2, 0.4)	$1.0977 imes 10^{-5}$	1.0977×10^{-6}	$8.5770 imes 10^{-6}$	$2.2556 imes 10^{-6}$
(0.3, 0.4)	1.2104×10^{-5}	1.1052×10^{-5}	$6.3960 imes 10^{-6}$	$9.7708 imes 10^{-7}$
(0.5, 0.5)	1.6539×10^{-5}	1.6540×10^{-5}	$5.2692 imes 10^{-6}$	5.2622×10^{-6}

Table 1. Example 1: Comparisons of numerical errors by Zhu et al. [8] and the present method with R = 100, $\tau = 10^{-4}$, at t = 0.1.

Table 2. Example 1: Comparison of the numerical errors by Zhu et al. [8] and the present method with R = 100, $\tau = 10^{-4}$ at t = 0.4.

Mesh Grid	Error of <i>u</i>	<i>u</i> Error of <i>v</i>		
	Zhu et al. [8]	The Present	Zhu et al. [8]	The Present
(0.1, 0.1)	1.0195×10^{-4}	2.2716×10^{-5}	$3.5483 imes10^{-4}$	4.2353×10^{-5}
(0.3, 0.1)	5.5872×10^{-4}	$8.7786 imes 10^{-5}$	1.0195×10^{-4}	1.9213×10^{-4}
(0.2, 0.2)	2.0389×10^{-4}	4.5433×10^{-5}	7.0967×10^{-4}	8.4705×10^{-5}
(0.4, 0.2)	6.6067×10^{-4}	1.0502×10^{-5}	4.5678×10^{-4}	2.3448×10^{-4}
(0.1, 0.3)	1.5094×10^{-4}	3.0802×10^{-6}	1.3174×10^{-3}	2.2716×10^{-5}
(0.3, 0.3)	$3.0584 imes10^{-4}$	6.8149×10^{-6}	1.0645×10^{-3}	$1.2706 imes10^{-4}$
(0.2, 0.4)	$4.8996 imes 10^{-5}$	$2.5797 imes 10^{-5}$	1.6722×10^{-3}	$1.9636 imes 10^{-5}$
(0.3, 0.4)	1.7939×10^{-4}	1.0068×10^{-4}	1.5458×10^{-3}	9.4524×10^{-5}
(0.5, 0.5)	5.0973×10^{-4}	1.1358×10^{-4}	1.7742×10^{-3}	2.1176×10^{-4}

In Table 3, the L_{∞} errors and the CPU time of the present method at t = 0.1 and t = 0.4 are listed. One can conclude that our method is efficient in sense that it could get a satisfying error with rather little CPU time. Figure 1 describes the numerical solutions of u(x, y, t), v(x, y, t) by the present method, as expected, the present method could simulate Example 1 quickly and rather precisely.

Table 3. Example 1: The L_{∞} errors and CPU time of the present method at t = 0.1 and t = 0.4 with $\tau = 10^{-4}$.

	$L_{\infty}(u)$	$L_\infty(v)$	CPU Time (s)
t = 0.1s	1.6540	1.6163×10^{-5}	6.728
t = 0.4s	1.6267×10^{-4}	$3.7444 imes10^{-4}$	25.658

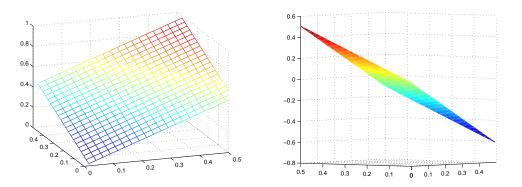


Figure 1. Example 1: Numerical solutions of u(x, y, t), and v(x, y, t) by the present method with $\tau = 10^{-4}$, $h_1 = h_2 = 0.025$, at t = 0.1.

Example 2. When applying the Hopf-Cole transformation in [6], one could get the exact solutions of Burgers' *Equations* (1) *and* (2)

$$u(x,y,t) = \frac{3}{4} - \frac{1}{4(1 + exp((-4x + 4y - t)R/32))},$$
(16)

$$v(x,y,t) = \frac{3}{4} + \frac{1}{4(1+exp((-4x+4y-t)R/32))}.$$
(17)

In this example, we choose the computational domain as $D = \{(x, y) : 0 \le x \le 1, 0 \le y \le 1\}$, and we take the initial and boundary conditions from the above exact solution. The knots are distributed uniformly on the computational domain with a width $h_1 = h_2 = 0.05$, and R = 100. However the time step sizes of our method ($\tau = 10^{-3}$) in this example is chosen to be ten times of the method in Zhu et al. [8], and Bahadir [7] ($\tau = 10^{-4}$). The parameter *c* is chosen as $c = 1.53 \times 10^{-5}$. The numerical results will show that we could get the similar accuracy while with little CPU time.

In Table 4, the L_{∞} numerical errors of the solutions at some typical mesh points at time t = 0.01, 0.05 and 0.2 with $\tau = 10^{-3}$. It is observed that our method could get satisfactory accuracy. In addition, the CPU time is rather little. The numerical results using the proposed algorithm are in good agreement with the exact solutions.

Mesh Grid	<i>t</i> = 0.01	t = 0.05	t = 0.2
(0, 1)	1.09714×10^{-8}	3.01044×10^{-7}	$1.39170 imes 10^{-5}$
(0.1, 0.9)	$5.21864 imes 10^{-8}$	$2.88925 imes 10^{-7}$	$1.12905 imes 10^{-5}$
(0.2, 0.8)	$6.05494 imes 10^{-7}$	$3.45039 imes 10^{-6}$	$1.52312 imes 10^{-5}$
(0.3, 0.7)	$6.52640 imes 10^{-6}$	$3.53596 imes 10^{-5}$	$1.70559 imes 10^{-4}$
(0.4, 0.6)	$6.56867 imes 10^{-6}$	$1.00372 imes 10^{-6}$	$5.22325 imes 10^{-4}$
(0.5, 0.5)	$5.55844 imes 10^{-5}$	$1.67227 imes 10^{-4}$	$6.56682 imes 10^{-4}$
(0.6, 0.4)	$3.08342 imes 10^{-5}$	$1.27268 imes 10^{-4}$	$2.60973 imes 10^{-4}$
(0.7, 0.3)	$2.12905 imes 10^{-7}$	$1.01505 imes 10^{-6}$	$7.31540 imes 10^{-6}$
(0.8, 0.2)	$1.47614 imes 10^{-8}$	$6.45096 imes 10^{-8}$	$9.81073 imes 10^{-6}$
(0.9, 0.1)	$2.52669 imes 10^{-9}$	$4.25191 imes 10^{-8}$	$1.28665 imes 10^{-5}$
(1, 0)	$3.35256 imes 10^{-8}$	$4.05295 imes 10^{-7}$	$1.00646 imes 10^{-5}$
CPU time (s)	0.130	0.404	1.418

Table 4. Example 2: The L_{∞} numerical errors and CPU time by the present method with R = 100, $\tau = 10^{-3}$ at different time *t*.

In Table 5, we compare the numerical errors of the present method with time step $\tau = 10^{-3}$, and that of Zhu et al. [8], and Bahadir [7] with time step size $\tau = 10^{-4}$. It is observed that even using the time step that is ten times of the method in Zhu et al. [8], and Bahadir [7], our method could get similar accuracy.

Table 5. Example 2: The error of numerical results by Zhu et al. [8], Bahadir[7] with R = 100, $\tau = 10^{-4}$, at t = 0.01, and the present method with $\tau = 10^{-3}$.

Mesh Grid	Error of <i>u</i>			Error of v		
	Zhu et al. [8]	Bahadir [7]	The Present	Zhu et al. [8]	Bahadir [7]	The Present
(0.1, 0.1)	5.91368×10^{-5}	7.24132×10^{-5}	5.44503×10^{-5}	5.91368×10^{-5}	$8.35601 imes 10^{-5}$	$5.44503 imes 10^{-5}$
(0.5, 0.1)	$4.84030 imes 10^{-6}$	$2.42869 imes 10^{-5}$	$6.49632 imes 10^{-6}$	$4.84030 imes 10^{-6}$	$5.13642 imes 10^{-5}$	$6.49632 imes 10^{-6}$
(0.9, 0.1)	$3.41000 imes 10^{-8}$	$8.39751 imes 10^{-6}$	$5.21864 imes 10^{-8}$	$3.41000 imes10^{-8}$	$7.03298 imes 10^{-6}$	$5.21864 imes 10^{-8}$
(0.3, 0.3)	$5.91368 imes 10^{-5}$	$8.25331 imes 10^{-5}$	$5.55844 imes 10^{-5}$	$5.91368 imes 10^{-5}$	$6.15201 imes 10^{-5}$	5.55844×10^{-5}
(0.7, 0.3)	$4.84030 imes 10^{-6}$	$3.43163 imes 10^{-5}$	$6.52640 imes 10^{-6}$	$4.84030 imes 10^{-6}$	$5.41000 imes 10^{-5}$	$6.52640 imes 10^{-6}$
(0.1, 0.5)	$1.64290 imes 10^{-6}$	$5.62014 imes 10^{-5}$	$2.71889 imes 10^{-7}$	$1.64290 imes 10^{-6}$	$7.35192 imes 10^{-5}$	$2.71889 imes 10^{-7}$
(0.5, 0.5)	$5.91368 imes 10^{-5}$	$7.32522 imes 10^{-5}$	5.55844×10^{-5}	5.91368×10^{-5}	$8.51040 imes 10^{-5}$	5.55844×10^{-5}

The numerical solutions of u(x, y, t), v(x, y, t) by the present method are shown in Figure 2. In conclusion, Example 2 shows that the present method is easy to implement, accurate, and rather time saving.

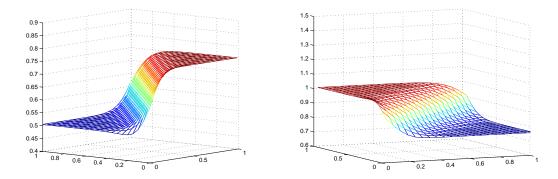


Figure 2. Example 2: The numerical solutions of u(x, y, t), and v(x, y, t) by the present method with $\tau = 10^{-3}$, $h_1 = h_2 = 0.05$ at t = 0.1

Example 3. In this example, we will show that the proposed scheme is also applicable to other high-dimensional PDEs with shock waves or even soliton waves. We take the typical soliton equation Sine–Gordon equation as the example:

$$u_t = v,$$

$$v_t = u_{xx} + u_{yy} - sinu,$$
(18)

The initial conditions are:

$$u(x, y, 0) = 4arctan(exp(x)) + 4arctan(exp(y)), -6 \le x, y \le 6, v(x, y, 0) = 0,$$

and the boundary conditions are:

$$u_x = 0, \text{ for } x = -6, 6, -6 \le y \le 6, u_y = 0, \text{ for } y = -6, 6, -6 \le x \le 6.$$

The example depicts the superposition of two orthogonal line solitons, which is also used by many researchers to verify their schemes [4,29]. The numerical computations are simulated using equi-distributed knots, with a space width $h_1 = h_2 = 0.2$ and time steps $\tau = 0.01$. In this example, the parameter *c* is chosen as c = 0.0075. Figure 3 shows that our methods could simulate the solitons rather exactly. That is to say, our scheme can solve a wild range of high-dimensional PDEs efficiently.

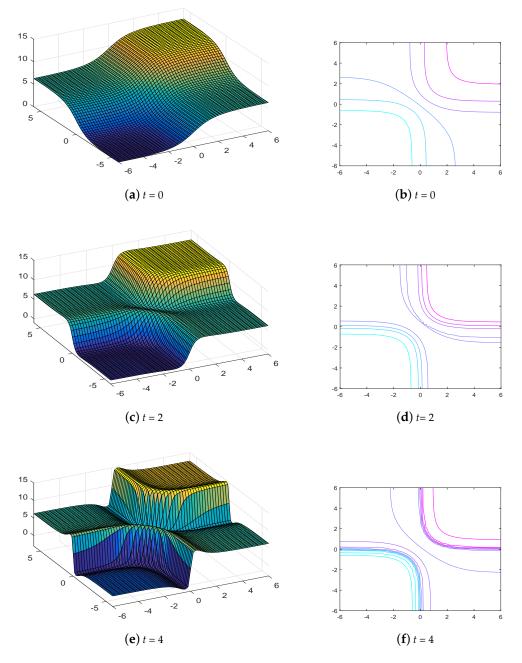


Figure 3. Example 3: The numerical solutions and contours by the present method with $\tau = 10^{-2}$, $h_1 = h_2 = 0.2$ at t = 0, 2, 4.

5. Conclusions

In this paper, we present a numerical scheme for 2-D Burgers' equation by bivariate MQ quasi-interpolation. The error estimations of the algorithm are also provided. Compared with the previous method (such as Zhu et al. [8]), the proposed method owns similar accuracy, while using far less computational effort. The present algorithm owns the following advantages: Firstly, it is very simple and easy to be implemented by the engineers in the applications. Secondly, it is of satisfactory accuracy and efficient for the time dependent PDEs with shock waves. Thirdly, it uses less computational effort, because it need not to solve any large-scale linear system. The method is also applicable to other high-dimensional time dependent PDEs [22,30], such as the Burgers-Fisher, the Klein–Gordon, and so forth.

Remark 2. Since the quasi-interpolation method have no requirement of the knots' topology structure and allows flexible knots moving, there are many works could be done to improve the efficiency and accuracy of the proposed algorithm with similar computational effort. One method is to introduce the moving knots strategy into the algorithm and let the nodes concentrated in the areas with shockwave, such as [31]. Hence better approximation accuracy could be achieved with the same computational effort. Another method is using the symmetric quasi-interpolation to construct the algorithm that preserve the energy or some important property of the original PDE, e.g., [12,22]. In this way, one could simulate the PDE for a longer time.

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Abbreviations

The following abbreviations are used in this manuscript:

- PDE Partial differential equation
- MQ Multi-quadric

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