

ADI method based on C^2 -continuous two-node integrated-RBF elements for viscous flows

D.-A. An-Vo¹, N. Mai-Duy¹, C.-D. Tran¹, T. Tran-Cong^{1,*}

*Corresponding author. Tel.: +61 7 4631 1332; fax: +61 7 4631 2110

Email addresses: duc-anh.an-vo@usq.edu.au (D.-A. An-Vo), nam.mai-duy@usq.edu.au (N. Mai-Duy), canh-dung.tran@usq.edu.au (C.-D. Tran), thanh.tran-cong@usq.edu.au (T. Tran-Cong)

¹Computational Engineering and Science Research Centre, Faculty of Engineering and Surveying, The University of Southern Queensland, Toowoomba, QLD 4350, Australia

Abstract: We propose a C^2 -continuous alternating direction implicit (ADI) method for the solution of the streamfunction-vorticity equations governing steady 2D incompressible viscous fluid flows. Discretisation is simply achieved with Cartesian grids. Local two-node integrated radial basis function elements (IRBFEs) [D.-A. An-Vo, N. Mai-Duy, T. Tran-Cong, A C^2 -continuous control-volume technique based on Cartesian grids and two-node integrated-RBF elements for second-order elliptic problems, CMES: Computer Modeling in Engineering & Sciences 72 (2011) 299-334] are used for the discretisation of the diffusion terms, and then the convection terms are incorporated into system matrices by treating nodal derivatives as unknowns. ADI procedure is applied for the time integration. Following ADI factorisation, the two-dimensional problem becomes a sequence of one-dimensional problems. The solution strategy consists of multiple use of a one-dimensional sparse matrix algorithm that helps saving the computational cost. High levels of accuracy and efficiency of the present methods are demonstrated with solutions of several benchmark problems defined on rectangular and non-rectangular domains.

Keywords: Navier-Stokes equations, integrated-radial-basis-function elements, C^2 -continuous solutions, Cartesian grid, ADI method, local RBF approximation.

1. Introduction

The dimensionless Navier-Stokes (N-S) equations for steady incompressible planar viscous flows, subject to negligible body forces, can be expressed in terms of the streamfunction ψ and the vorticity ω as follows.

$$\frac{\partial^2 \psi}{\partial x^2} + \frac{\partial^2 \psi}{\partial y^2} + \omega = 0, \quad (1)$$

$$\frac{\partial^2 \omega}{\partial x^2} + \frac{\partial^2 \omega}{\partial y^2} = Re \left(\frac{\partial \psi}{\partial y} \frac{\partial \omega}{\partial x} - \frac{\partial \psi}{\partial x} \frac{\partial \omega}{\partial y} \right), \quad (x, y)^T \in \Omega, \quad (2)$$

where $Re = UL/\nu$ is the Reynolds number, in which L is the characteristic length, U the characteristic speed of the flow and ν the kinematic viscosity of the fluid. The vorticity and streamfunction variables are defined by

$$\omega = \frac{\partial v}{\partial x} - \frac{\partial u}{\partial y}, \quad (3)$$

$$\frac{\partial \psi}{\partial y} = u, \quad \frac{\partial \psi}{\partial x} = -v, \quad (4)$$

where u and v are the x and y components of the velocity vector. In this study, the method of modified dynamics or false transients (e.g. [1, 2]) is applied to obtain the structure of a steady flow. The governing equations (1) and (2) are modified as

$$\frac{\partial^2 \psi}{\partial x^2} + \frac{\partial^2 \psi}{\partial y^2} + \omega = 0, \quad (5)$$

$$\frac{\partial \omega}{\partial t} + \frac{\partial^2 \omega}{\partial x^2} + \frac{\partial^2 \omega}{\partial y^2} = Re \left(\frac{\partial \psi}{\partial y} \frac{\partial \omega}{\partial x} - \frac{\partial \psi}{\partial x} \frac{\partial \omega}{\partial y} \right). \quad (6)$$

A steady state solution to (5) and (6), which is obtained by integrating the equations from a given initial condition up to the steady state, is also solution to (1) and (2).

Cartesian-grid-based methods for solving (1) and (2) can be very economical owing to the facts that (i) generating a grid is low-cost; and (ii) ADI procedure [3, 4] can be straightforwardly applied to accelerate computational processes. The approximations for the dependent variables and their spatial derivatives can be constructed globally on the whole grid or locally on small segments of the grid. A very prominent local approximation scheme is the finite-difference (FD) which can be based on two nodes (first-order accuracy) and three nodes. The three-node approximations can take the second-order central difference (CD) form, e.g. [5], or high-order compact (HOC) implicit forms, e.g. [6–8], where nodal values of the field variables and their derivatives are considered as unknowns. The three-node HOC implicit schemes can achieve higher order numerical accuracy and yield greater computational efficiency compared with CD schemes for the same level of the accuracy, e.g. [7]. However, the computational cost of these implicit schemes was quite high because of time consumption for solving (i) less-than-optimal banded matrices (block diagonal structure where each block corresponds to a grid line) [9–11]; or (ii) a larger number of equations per grid point [7], i.e. $3N$ equations for N grid points in 1D problems and $5N$ equations for N grid points in 2D problems. In addition, these finite difference schemes (i.e. the two-node and the three-node schemes) typically produce solutions which are continuous for the fields but not for their partial derivatives, i.e. C^0 -continuity. The grid thus needs to be sufficiently fine to mitigate the effects of discontinuity of partial derivatives.

On the other hand, the well-known alternating direction implicit (ADI) method proposed by Peaceman and Rachford (PR-ADI) was much more computationally cost-effective than the HOC implicit schemes. In the PR-ADI method, the solution is obtained by solving sets of equations defined on grid lines in x - and then y -direction sequentially. Each set

includes N_η equations (matrix dimensions are $N_\eta \times N_\eta$) for N_η grid points on a particular η -grid line where η represents x and y . These relatively small sets of equation (in tridiagonal matrix form) are solved separately and effectively by the Thomas algorithm that helps saving the computational cost. However, the numerical accuracy of PR-ADI is only second-order in space [3]. The combination of the ADI approach and the HOC schemes has been proposed by e.g. [7, 10] for solving fluid mechanics problems, by e.g. [12] for parabolic partial differential equations, and by e.g. [9, 11] for convection-diffusion problems. Hirsh [7] applied the ADI procedure to HOC implicit schemes for simulating a model square cavity flow through solving sets of $3N_\eta$ equations for N_η grid points on η -grid lines. Adam [12] further reduced the number of equations on each grid line to sets of $2N_\eta$ equations by means of the so-called implicit elimination. Recently, Karaa and Zhang [9] and Karaa [10] solve sets of N_η equations on η -grid lines through block matrices. However, as shown in [13], the solution quality of this ADI method is degraded for convection-diffusion equations with high Peclet numbers. Ma et al. [11] proposed to use fourth-order schemes for convection terms and second-order schemes for diffusion terms for convection-dominated diffusion problems and achieved very efficient sets of N_η equations in tridiagonal matrix form on η -grid lines. This ADI method hence becomes second-order accurate when diffusion terms are dominant.

Radial basis functions (RBFs) have recently emerged as an attractive tool for the solution of ordinary and partial differential equations (ODEs and PDEs), e.g. [14–16]. RBF-based approximants can be constructed through a conventional differentiation process, e.g. [17], or an integration process, e.g. [18–20]. RBF-based approximants can be global or local. Global RBF-based methods are very accurate, e.g. [21, 22]. However, they result in a system matrix that is dense and usually highly ill-conditioned for large problems. The use of RBF-approximants in local forms can help circumvent these difficulties, e.g. [23–25]. Recently, a local high order approximant based on 2-node elements and integrated RBFs (IRBFs) has been proposed by An-Vo et al. [26]. It was shown that such elements lead to a C^2 -continuous solution rather than the usual C^0 -continuous solution.

In this paper, we develop a high-order ADI method based on C^2 -continuous 2-node IRBFs [26] for solving the N-S equations in the form of streamfunction-vorticity formulation, discretised by Cartesian grids. Unlike finite difference methods, the proposed method can guarantee inter-element continuity of derivatives of the streamfunction and vorticity of or-

ders up to 2. 2-node IRBFEs are used for the discretisation of the diffusion terms, and then the convection terms are incorporated into system matrices by treating nodal first-derivatives as unknowns. By treating the convection terms as unknowns, we obtain matrices on grid lines that are always diagonally dominant. The matrix of each η -grid line includes $2N_\eta$ equations for N_η grid points as in Adam [12] without the need of implicit elimination. It is noted that in Adam [12], one has 6 nonzero entries for the governing equation and 5 nonzero entries for the equation of first-derivatives at a grid point. It will be shown later that the proposed C^2 -continuous ADI method yields 4 nonzero entries for the governing equation and 6 nonzero entries for the equation of first-derivatives (by imposing C^2 -continuity condition) at a grid point. Several viscous flows defined on rectangular and non-rectangular domains are considered to verify the proposed method in terms of computational cost and numerical accuracy on a wide range of Reynolds number.

The remainder of the paper is organised as follows. A brief review of integrated RBF elements is given in Section 2. Section 3 describes the proposed C^2 -continuous ADI method for the streamfunction-vorticity formulation. In section 4, viscous flows in square and triangular cavities are presented to demonstrate the attractiveness of the present method. Section 5 concludes the paper.

2. Two-node integrated-RBF elements (IRBFEs)

The problem domain is embedded in a Cartesian grid as shown in Figure 1. Grid points inside the problem domain are taken to be interior nodes, while boundary nodes are defined as the intersections of the grid lines and the boundaries. There are two types of elements, namely interior and semi-interior IRBFEs. An interior element is formed using two adjacent interior nodes while a semi-interior element is generated by an interior node and a boundary node.

2.1. Interior elements

Consider an interior element, $\eta \in [\eta_1, \eta_2]$, and its two nodes are locally named as 1 and 2. Let $\phi(\eta)$ be a function and $\phi_1, \partial\phi_1/\partial\eta, \phi_2$ and $\partial\phi_2/\partial\eta$ be the values of ϕ and $\partial\phi/\partial\eta$ at the two nodes, respectively (Figure 2a). The 2-node IRBFE scheme approximates the second-order

derivative of $\phi(\eta)$ using two multiquadric (MQ) functions whose centres are located at η_1 and η_2

$$\frac{\partial^2 \phi}{\partial \eta^2}(\eta) = \frac{d^2 \phi}{d\eta^2}(\eta) = w_1 \sqrt{(\eta - \eta_1)^2 + a_1^2} + w_2 \sqrt{(\eta - \eta_2)^2 + a_2^2} = w_1 I_1^{(2)}(\eta) + w_2 I_2^{(2)}(\eta), \quad (7)$$

where $I_i^{(2)}(\eta)$ conveniently denotes the MQ, w_i and a_i are the associated weight and MQ-width at node i ($i = \{1, 2\}$). We simply take $a_i = \beta h$, where h is a grid size and β is a factor.

First-order derivative of ϕ and the function ϕ are approximated by integrating (7) with respect to η

$$\frac{\partial \phi}{\partial \eta}(\eta) = \frac{d\phi}{d\eta}(\eta) = w_1 I_1^{(1)}(\eta) + w_2 I_2^{(1)}(\eta) + C_1, \quad (8)$$

$$\phi(\eta) = w_1 I_1^{(0)}(\eta) + w_2 I_2^{(0)}(\eta) + C_1 \eta + C_2, \quad (9)$$

where $I_i^{(1)}(\eta) = \int I_i^{(2)}(\eta) d\eta$, $I_i^{(0)}(\eta) = \int I_i^{(1)}(\eta) d\eta$, and C_1 and C_2 are the constants of integration. By collocating (9) and (8) at η_1 and η_2 , the relation between the physical space and the RBF coefficient space is obtained

$$\underbrace{\begin{pmatrix} \phi_1 \\ \phi_2 \\ \frac{\partial \phi_1}{\partial \eta} \\ \frac{\partial \phi_2}{\partial \eta} \end{pmatrix}}_{\widehat{\phi}} = \underbrace{\begin{pmatrix} I_1^{(0)}(\eta_1) & I_2^{(0)}(\eta_1) & \eta_1 & 1 \\ I_1^{(0)}(\eta_2) & I_2^{(0)}(\eta_2) & \eta_2 & 1 \\ I_1^{(1)}(\eta_1) & I_2^{(1)}(\eta_1) & 1 & 0 \\ I_1^{(1)}(\eta_2) & I_2^{(1)}(\eta_2) & 1 & 0 \end{pmatrix}}_{\mathcal{I}} \underbrace{\begin{pmatrix} w_1 \\ w_2 \\ C_1 \\ C_2 \end{pmatrix}}_{\widehat{w}}, \quad (10)$$

where $\widehat{\phi}$ is the nodal-value vector, \mathcal{I} the conversion matrix, and \widehat{w} the coefficient vector. It is noted that not only the nodal values of ϕ but also of $\partial\phi/\partial\eta$ are incorporated into the conversion system and this imposition is done in an exact manner owing to the presence of integration constants. Solving (10) yields

$$\widehat{w} = \mathcal{I}^{-1} \widehat{\phi}. \quad (11)$$

Substitution of (11) into (9), (8) and (7) leads to

$$\phi(\eta) = \left[I_1^{(0)}(\eta), I_2^{(0)}(\eta), \eta, 1 \right] \mathcal{I}^{-1} \widehat{\phi}, \quad (12)$$

$$\frac{\partial \phi}{\partial \eta}(\eta) = \left[I_1^{(1)}(\eta), I_2^{(1)}(\eta), 1, 0 \right] \mathcal{I}^{-1} \widehat{\phi}, \quad (13)$$

$$\frac{\partial^2 \phi}{\partial \eta^2}(\eta) = \left[I_1^{(2)}(\eta), I_2^{(2)}(\eta), 0, 0 \right] \mathcal{I}^{-1} \widehat{\phi}. \quad (14)$$

They can be rewritten in the form

$$\phi(\eta) = \varphi_1(\eta)\phi_1 + \varphi_2(\eta)\phi_2 + \varphi_3(\eta)\frac{\partial \phi_1}{\partial \eta} + \varphi_4(\eta)\frac{\partial \phi_2}{\partial \eta}, \quad (15)$$

$$\frac{\partial \phi}{\partial \eta}(\eta) = \frac{d\varphi_1(\eta)}{d\eta}\phi_1 + \frac{d\varphi_2(\eta)}{d\eta}\phi_2 + \frac{d\varphi_3(\eta)}{d\eta}\frac{\partial \phi_1}{\partial \eta} + \frac{d\varphi_4(\eta)}{d\eta}\frac{\partial \phi_2}{\partial \eta}, \quad (16)$$

$$\frac{\partial^2 \phi}{\partial \eta^2}(\eta) = \frac{d^2\varphi_1(\eta)}{d\eta^2}\phi_1 + \frac{d^2\varphi_2(\eta)}{d\eta^2}\phi_2 + \frac{d^2\varphi_3(\eta)}{d\eta^2}\frac{\partial \phi_1}{\partial \eta} + \frac{d^2\varphi_4(\eta)}{d\eta^2}\frac{\partial \phi_2}{\partial \eta}, \quad (17)$$

where $\{\varphi_i(\eta)\}_{i=1}^4$ is the set of basis functions in the physical space. These expressions allow one to compute the values of ϕ , $\partial\phi/\partial\eta$, and $\partial^2\phi/\partial\eta^2$ at any point η in $[\eta_1, \eta_2]$ in terms of four nodal unknowns, i.e. the values of the field variable and its first-order derivatives at the two extremes (also grid points) of the element.

For convenience, in the case of $\eta \equiv x$, we denote

$$\mu_i = \frac{d^2\varphi_i(x_1)}{dx^2}, \quad (18)$$

$$\nu_i = \frac{d^2\varphi_i(x_2)}{dx^2}, \quad (19)$$

and in the case of $\eta \equiv y$,

$$\theta_i = \frac{d^2\varphi_i(y_1)}{dy^2}, \quad (20)$$

$$\vartheta_i = \frac{d^2\varphi_i(y_2)}{dy^2}, \quad i = \{1, 2, 3, 4\}. \quad (21)$$

2.2. Semi-interior elements

As mentioned earlier, a semi-interior element is defined by two nodes: an interior node and a boundary node. The subscripts 1 and 2 are now replaced with b (b represents a boundary node) and g (g an interior grid node), respectively (Figure 2b). Assume that the value of ϕ

is given at η_b . The conversion system can be formed as

$$\begin{pmatrix} \phi_b \\ \phi_g \\ \frac{\partial \phi_g}{\partial \eta} \end{pmatrix} = \begin{pmatrix} I_b^{(0)}(\eta_b) & I_g^{(0)}(\eta_b) & \eta_b & 1 \\ I_b^{(0)}(\eta_g) & I_g^{(0)}(\eta_g) & \eta_g & 1 \\ I_b^{(1)}(\eta_g) & I_g^{(1)}(\eta_g) & 1 & 0 \end{pmatrix} \begin{pmatrix} w_b \\ w_g \\ C_1 \\ C_2 \end{pmatrix}. \quad (22)$$

(22) leads to

$$\phi(\eta) = \varphi_1(\eta)\phi_b + \varphi_2(\eta)\phi_g + \varphi_3(\eta)\frac{\partial \phi_g}{\partial \eta}, \quad (23)$$

$$\frac{\partial \phi}{\partial \eta}(\eta) = \frac{d\varphi_1(\eta)}{d\eta}\phi_b + \frac{d\varphi_2(\eta)}{d\eta}\phi_g + \frac{d\varphi_3(\eta)}{d\eta}\frac{\partial \phi_g}{\partial \eta}, \quad (24)$$

$$\frac{\partial^2 \phi}{\partial \eta^2}(\eta) = \frac{d^2\varphi_1(\eta)}{d\eta^2}\phi_b + \frac{d^2\varphi_2(\eta)}{d\eta^2}\phi_g + \frac{d^2\varphi_3(\eta)}{d\eta^2}\frac{\partial \phi_g}{\partial \eta}. \quad (25)$$

It can be seen that the conversion matrix in (22) is under-determined and its inverse can be obtained using the SVD technique (pseudo-inversion). Owing to the facts that point collocation is used and the RBF conversion matrix is not over-determined, the boundary condition ϕ_b is imposed in an exact manner. For other types of semi-interior elements, the reader is referred to An-Vo et al. [26] for details.

3. Derivation of C^2 -continuous ADI method

3.1. ADI scheme for N - S equations on a Cartesian grid

Consider a grid point P and its east, west, north and south neighbouring nodes denoted as E , W , N and S , respectively (Figure 3). Collocating (5) and (6) at P , one obtains

$$\frac{\partial^2 \psi_P}{\partial x^2} + \frac{\partial^2 \psi_P}{\partial y^2} + \omega_P = 0, \quad (26)$$

$$\frac{\partial \omega_P}{\partial t} + \frac{\partial^2 \omega_P}{\partial x^2} + \frac{\partial^2 \omega_P}{\partial y^2} = Re \left(\frac{\partial \psi_P}{\partial y} \frac{\partial \omega_P}{\partial x} - \frac{\partial \psi_P}{\partial x} \frac{\partial \omega_P}{\partial y} \right). \quad (27)$$

We now employ the ADI (Alternating-Direction Implicit) procedure [3, 4] to relax the time derivative term in (27) in two stages. At a time instant t^n , (41) and (27) become

$$\frac{\partial^2 \psi_P^n}{\partial x^2} + \frac{\partial^2 \psi_P^n}{\partial y^2} + \omega_P^{n-1} = 0, \quad (28)$$

$$\frac{\omega_P^{n-1/2} - \omega_P^{n-1}}{\Delta t/2} + \frac{\partial^2 \omega_P^{n-1/2}}{\partial x^2} + \frac{\partial^2 \omega_P^{n-1}}{\partial y^2} = Re \left(\frac{\partial \psi_P^n}{\partial y} \frac{\partial \omega_P^{n-1/2}}{\partial x} - \frac{\partial \psi_P^n}{\partial x} \frac{\partial \omega_P^{n-1}}{\partial y} \right), \quad (29)$$

$$\frac{\omega_P^n - \omega_P^{n-1/2}}{\Delta t/2} + \frac{\partial^2 \omega_P^{n-1/2}}{\partial x^2} + \frac{\partial^2 \omega_P^n}{\partial y^2} = Re \left(\frac{\partial \psi_P^n}{\partial y} \frac{\partial \omega_P^{n-1/2}}{\partial x} - \frac{\partial \psi_P^n}{\partial x} \frac{\partial \omega_P^n}{\partial y} \right). \quad (30)$$

It can be seen that in the first stage, i.e. (29), $\partial^2 \omega_P^{n-1/2} / \partial x^2$ and $\partial \omega_P^{n-1/2} / \partial x$ are treated implicitly and $\partial^2 \omega_P^n / \partial y^2$ and $\partial \omega_P^n / \partial y$ are treated implicitly in the second stage, i.e. (30). These derivatives and the second-order derivatives of streamfunction in (28) are typically approximated by a second-order CD scheme, e.g. [5], or HOC implicit schemes, e.g. [6, 7, 12, 27]. For instance in x -direction, one has

$$\frac{\partial \omega_P^{n-1/2}}{\partial x} = \frac{\omega_E^{n-1/2} - \omega_W^{n-1/2}}{2h} + O(h^2), \quad (31)$$

$$\frac{\partial^2 \omega_P^{n-1/2}}{\partial x^2} = \frac{\omega_E^{n-1/2} - 2\omega_P^{n-1/2} + \omega_W^{n-1/2}}{h^2} + O(h^2), \quad (32)$$

$$\frac{\partial^2 \psi_P^n}{\partial x^2} = \frac{\psi_E^n - 2\psi_P^n + \psi_W^n}{h^2} + O(h^2), \quad (33)$$

or

$$\frac{1}{6} \frac{\partial \omega_W^{n-1/2}}{\partial x} + \frac{2}{3} \frac{\partial \omega_P^{n-1/2}}{\partial x} + \frac{1}{6} \frac{\partial \omega_E^{n-1/2}}{\partial x} = \frac{\omega_E^{n-1/2} - \omega_W^{n-1/2}}{2h} + O(h^4), \quad (34)$$

$$\frac{1}{12} \frac{\partial^2 \omega_W^{n-1/2}}{\partial x^2} + \frac{10}{12} \frac{\partial^2 \omega_P^{n-1/2}}{\partial x^2} + \frac{1}{12} \frac{\partial^2 \omega_E^{n-1/2}}{\partial x^2} = \frac{\omega_E^{n-1/2} - 2\omega_P^{n-1/2} + \omega_W^{n-1/2}}{h^2} + O(h^4), \quad (35)$$

$$\frac{1}{12} \frac{\partial^2 \psi_W^n}{\partial x^2} + \frac{10}{12} \frac{\partial^2 \psi_P^n}{\partial x^2} + \frac{1}{12} \frac{\partial^2 \psi_E^n}{\partial x^2} = \frac{\psi_E^n - 2\psi_P^n + \psi_W^n}{h^2} + O(h^4). \quad (36)$$

In (31), (32) and (33) the derivatives are explicitly given in term of nodal values of the field variable while one has to solve tridiagonal systems to obtain the derivative values in (34), (35) and (36). If one makes use of (31) or (34) to approximate the convection term of (29), the obtained system matrix might not be diagonally dominant at high values of the Re . As a result, they suffer from the so-called cell Reynolds number limitation (e.g. [28]). Similar discussions can be made for equation (30) in y -direction.

3.2. Proposed C^2 -continuous IRBFE-ADI method

As in Figure 3, one can form 4 two-node IRBFEs associated with P , namely WP , PE , SP and PN , assumed to be interior elements. To approximate $\partial^2\psi_P^n/\partial x^2$ and $\partial^2\omega_P^{n-1/2}/\partial x^2$, $\partial^2\psi_P^n/\partial y^2$ and $\partial^2\omega_P^n/\partial y^2$ via (17), we propose to use the elements WP , SP , respectively, with abbreviations (19) and (21),

$$\frac{\partial^2\psi_P^n}{\partial x^2} = \nu_1\psi_W^n + \nu_2\psi_P^n + \nu_3\frac{\partial\psi_W^n}{\partial x} + \nu_4\frac{\partial\psi_P^n}{\partial x}, \quad (37)$$

$$\frac{\partial^2\omega_P^{n-1/2}}{\partial x^2} = \nu_1\omega_W^{n-1/2} + \nu_2\omega_P^{n-1/2} + \nu_3\frac{\partial\omega_W^{n-1/2}}{\partial x} + \nu_4\frac{\partial\omega_P^{n-1/2}}{\partial x}, \quad (38)$$

$$\frac{\partial^2\psi_P^n}{\partial y^2} = \vartheta_1\psi_S^n + \vartheta_2\psi_P^n + \vartheta_3\frac{\partial\psi_S^n}{\partial y} + \vartheta_4\frac{\partial\psi_P^n}{\partial y}, \quad (39)$$

$$\frac{\partial^2\omega_P^n}{\partial y^2} = \vartheta_1\omega_S^n + \vartheta_2\omega_P^n + \vartheta_3\frac{\partial\omega_S^n}{\partial y} + \vartheta_4\frac{\partial\omega_P^n}{\partial y}. \quad (40)$$

It will be shown later that C^2 -continuous conditions are imposed at P in both x - and y -grid lines. As a result, instead of using element WP to give approximations of $\partial^2\psi_P^n/\partial x^2$ and $\partial^2\omega_P^{n-1/2}/\partial x^2$, we are able to use element PE as a replacement. Similarly, element PN can replace SP to give approximations for $\partial^2\psi_P^n/\partial y^2$ and $\partial^2\omega_P^n/\partial y^2$. These possibilities will give the same results. Substituting (37) and (39) into (28), (38) into (29), and (40) into (30), we have

$$\nu_1\psi_W^n + \vartheta_1\psi_S^n + (\nu_2 + \vartheta_2)\psi_P^n + \nu_3\frac{\partial\psi_W^n}{\partial x} + \vartheta_3\frac{\partial\psi_S^n}{\partial y} + \nu_4\frac{\partial\psi_P^n}{\partial x} + \vartheta_4\frac{\partial\psi_P^n}{\partial y} = \omega_P^{n-1}, \quad (41)$$

$$\begin{aligned} \nu_1\omega_W^{n-1/2} + (\nu_2 + \frac{1}{\Delta t/2})\omega_P^{n-1/2} + \nu_3\frac{\partial\omega_W^{n-1/2}}{\partial x} + (\nu_4 - Re\frac{\partial\psi_P^n}{\partial y})\frac{\partial\omega_P^{n-1/2}}{\partial x} = \\ \frac{\omega_P^{n-1}}{\Delta t/2} - \frac{\partial^2\omega_P^{n-1}}{\partial y^2} - Re\frac{\partial\psi_P^n}{\partial x}\frac{\partial\omega_P^{n-1}}{\partial y}, \end{aligned} \quad (42)$$

$$\begin{aligned} \vartheta_1\omega_S^n + (\vartheta_2 + \frac{1}{\Delta t/2})\omega_P^n + \vartheta_3\frac{\partial\omega_S^n}{\partial y} + (\vartheta_4 + Re\frac{\partial\psi_P^n}{\partial x})\frac{\partial\omega_P^n}{\partial y} = \\ \frac{\omega_P^{n-1/2}}{\Delta t/2} - \frac{\partial^2\omega_P^{n-1/2}}{\partial x^2} + Re\frac{\partial\psi_P^n}{\partial y}\frac{\partial\omega_P^{n-1/2}}{\partial x}. \end{aligned} \quad (43)$$

Thus, at a nodal point P in (41) there are three unknowns, namely ψ_P^n , $\partial\psi_P^n/\partial x$ and $\partial\psi_P^n/\partial y$. To solve (41), two additional equations are needed and devised here by imposing C^2 -continuous conditions at P in x - and y -directions, i.e.

$$\left(\frac{\partial^2\psi_P^n}{\partial x^2}\right)_L = \left(\frac{\partial^2\psi_P^n}{\partial x^2}\right)_R, \quad (44)$$

$$\left(\frac{\partial^2\psi_P^n}{\partial y^2}\right)_B = \left(\frac{\partial^2\psi_P^n}{\partial y^2}\right)_T, \quad (45)$$

where $(\cdot)_L$ indicates that the computation of (\cdot) is based on the element to the left of P , i.e. element WP , and similarly subscript R, B, T denotes the right (PE), bottom (SP) and top (PN) elements. The left of equations (44) and (45) are replaced by (37) and (39) and the right by similar expressions obtained via (17), noting (18) and (20) respectively, yielding

$$\nu_1\psi_W^n + \nu_2\psi_P^n + \nu_3\frac{\partial\psi_W^n}{\partial x} + \nu_4\frac{\partial\psi_P^n}{\partial x} = \mu_1\psi_P^n + \mu_2\psi_E^n + \mu_3\frac{\partial\psi_P^n}{\partial x} + \mu_4\frac{\partial\psi_E^n}{\partial x}, \quad (46)$$

$$\vartheta_1\psi_S^n + \vartheta_2\psi_P^n + \vartheta_3\frac{\partial\psi_S^n}{\partial y} + \vartheta_4\frac{\partial\psi_P^n}{\partial y} = \theta_1\psi_P^n + \theta_2\psi_N^n + \theta_3\frac{\partial\psi_P^n}{\partial y} + \theta_4\frac{\partial\psi_N^n}{\partial y}. \quad (47)$$

At the nodal point P and for the vorticity field, in the first relaxation stage in the x -direction, there are two unknowns in (42), namely $\omega_P^{n-1/2}$ and $\partial\omega_P^{n-1/2}/\partial x$ and in the second stage of relaxation in the y -direction, two unknowns in (43), namely ω_P^n and $\partial\omega_P^n/\partial y$. To solve (42), one additional equation is needed and also devised by imposing C^2 -continuity condition at P in x -direction, i.e.

$$\left(\frac{\partial^2\omega_P^{n-1/2}}{\partial x^2}\right)_L = \left(\frac{\partial^2\omega_P^{n-1/2}}{\partial x^2}\right)_R. \quad (48)$$

The left of equation (48) is replaced by (38) and the right by a similar expression obtained via (17), noting (18), yielding

$$\nu_1\omega_W^{n-1/2} + \nu_2\omega_P^{n-1/2} + \nu_3\frac{\partial\omega_W^{n-1/2}}{\partial x} + \nu_4\frac{\partial\omega_P^{n-1/2}}{\partial x} = \mu_1\omega_P^{n-1/2} + \mu_2\omega_E^{n-1/2} + \mu_3\frac{\partial\omega_P^{n-1/2}}{\partial x} + \mu_4\frac{\partial\omega_E^{n-1/2}}{\partial x}. \quad (49)$$

In a similar manner, to solve (43), one additional equation is created by imposing C^2 -continuity condition at P in y -direction, i.e.

$$\left(\frac{\partial^2\omega_P^n}{\partial y^2}\right)_B = \left(\frac{\partial^2\omega_P^n}{\partial y^2}\right)_T, \quad (50)$$

The left of equation (50) is replaced by (40) and the right by a similar expression obtained via (17), noting (20), yielding

$$\vartheta_1\omega_S^n + \vartheta_2\omega_P^n + \vartheta_3\frac{\partial\omega_S^n}{\partial y} + \vartheta_4\frac{\partial\omega_P^n}{\partial y} = \theta_1\omega_P^n + \theta_2\omega_N^n + \theta_3\frac{\partial\omega_P^n}{\partial y} + \theta_4\frac{\partial\omega_N^n}{\partial y}. \quad (51)$$

(41-43), (46), (47), (49) and (51) will be slightly different at the domain boundary where (25) for semi-interior elements is used instead of (17). To solve for the streamfunction field we collect equations (41), (46) and (47) at all nodal points which leads to a large system matrix of size $3N \times 3N$ where N is the total number of grid points of the problem domain. Nonetheless, this system is sparse and banded. As a result, the LU technique is very efficient for factorisation. It is noted that the factorisation needs to be done only once.

For the vorticity field, it can be seen from (42), (43), (49) and (51) that there are 4 nonzero entries for the governing equations, i.e. (42) and (43), and 6 nonzero entries for the C^2 -continuity conditions at the grid point P , i.e. (49) and (51). At the first relaxation stage, collection of equations (42) and (49) at nodal points on each and every x -grid line leads to N_y independent sets of equations. Each set contains $2N_x$ equations for $2N_x$ unknowns associated with an x -grid line with N_x nodes. At the second stage, collection of equations (43) and (51) at nodal points on each and every y -grid line leads to N_x independent sets of equations. Each set contains $2N_y$ equations for $2N_y$ unknowns associated with a y -grid line with N_y nodes. In contrast to the direct solution approaches in [26, 29] where a system of $3N$ equations for $3N$ unknowns are required, the current approach results in considerable savings in terms of both storage and computational time. The latter is significantly reduced further when parallelisation is implemented to independently solve these relatively small sets of $2N_\eta$ equations.

At high values of the Re , the fourth terms on the LHS of (42) and (43) are dominant, which guarantees diagonally dominant system matrices. Owing to the fact that two-node IRBFs are used, the proposed method also leads to very sparse systems and its solution is a C^2 function across IRBFs.

4. Numerical examples

The performance of the proposed C^2 -continuous IRBFE-ADI method is studied through the simulation of flows in square and triangular cavities. For all numerical examples presented in this study, the MQ shape parameter a is simply chosen proportionally to the element length h by a factor β . $\beta = 1$ is used throughout the computations. In the case of non-rectangular domains, there may be some nodes that are too close to the boundary. If an interior node falls within a distance of $h/4$ to the boundary, such a node is removed from the set of nodal points. A steady solution is obtained with a time marching approach starting from a computed solution at a lower Reynolds number. For the special case of Stokes equation, the starting condition is the rest state.

The solution procedure involves the following steps

- (1) Guess the initial distributions of the streamfunction and vorticity in the case of Stokes

flow. Otherwise, take the solution of a lower Reynolds number as an initial guess.

(2) Discretise the streamfunction equation at a time instant t^n (28) by means of C^2 -continuous IRBFEs, i.e. (41), (46) and (47), and then apply the LU technique to factorise the system matrix into two triangular matrices. It is noted that the factorisation needs to be done only once.

(3) Solve (28) subjects to boundary conditions for the new streamfunction field.

(3) Derive a computational boundary condition for the vorticity from the updated streamfunction field.

(4) Solve for the new vorticity field in two stages by using (42) and (49), (43) and (51) in x - and y -directions respectively.

(5) Check to see whether the solution has reached a steady state through a condition on convergence measure

$$CM(\psi) = \frac{\sqrt{\sum_{i=1}^N (\psi_i - \psi_i^0)^2}}{\sqrt{\sum_{i=1}^N \psi_i^2}} < 10^{-9}, \quad (52)$$

where N is the total number of grid nodes.

(6) If CM is not satisfactorily small, advance pseudo-time and repeat from step (3). Otherwise, stop the computation and output the results.

4.1. Square cavity

Square cavity flow is the most studied case in the literature of internal flows. This type of flow is important firstly in its own right as a basic physical model. Then, owing to its simple geometry and rich flow physics at different Reynolds numbers, the problem also serves as a useful test for numerical algorithms in CFD. The cavity is taken to be a unit square, with the lid sliding from left to right at a unit velocity. The boundary conditions can be specified as

$$\begin{aligned} \psi = 0, \quad \partial\psi/\partial x = 0 \quad \text{on} \quad x = 0, \quad x = 1, \\ \psi = 0, \quad \partial\psi/\partial y = 0 \quad \text{on} \quad y = 0, \\ \psi = 0, \quad \partial\psi/\partial y = 1 \quad \text{on} \quad y = 1. \end{aligned}$$

We take Dirichlet boundary conditions, $\psi = 0$, on all walls for solving (5). The Neumann boundary conditions, $\partial\psi/\partial n$ (i.e. $\partial\psi/\partial n = \nabla\psi \cdot \hat{n}$, where \hat{n} is the outward unit vector

normal to the boundary), are used to derive computational boundary conditions for ω in solving (6). Making use of (5), the values of ω on the boundaries are computed by

$$\omega_b = -\frac{\partial^2 \psi_b}{\partial x^2} \quad \text{on } x = 0, \quad x = 1, \quad (53)$$

$$\omega_b = -\frac{\partial^2 \psi_b}{\partial y^2} \quad \text{on } y = 0, \quad y = 1. \quad (54)$$

In computing (53) and (54), one needs to incorporate $\partial\psi_b/\partial x$ into $\partial^2\psi_b/\partial x^2$, and $\partial\psi_b/\partial y$ into $\partial^2\psi_b/\partial y^2$, respectively. A simple technique to derive ω_b in the context of 2-node IRBFs can be found in [29]. It will be briefly reproduced here for the sake of completeness. Assuming that node 1 and 2 of an IRBFE are a boundary node and an interior grid node respectively (i.e. $1 \equiv b$ and $2 \equiv g$). Boundary values of the vorticity are obtained by applying (17) as

$$\omega_b = -\frac{\partial^2 \psi_b}{\partial \eta^2} = -\left(\frac{d^2 \varphi_1(\eta_b)}{d\eta^2} \psi_b + \frac{d^2 \varphi_2(\eta_b)}{d\eta^2} \psi_g + \frac{d^2 \varphi_3(\eta_b)}{d\eta^2} \frac{\partial \psi_b}{\partial \eta} + \frac{d^2 \varphi_4(\eta_b)}{d\eta^2} \frac{\partial \psi_g}{\partial \eta} \right), \quad (55)$$

where η represents x and y ; ψ_b and $\partial\psi_b/\partial\eta$ are the Dirichlet and Neumann boundary conditions for ψ , and ψ_g and $\partial\psi_g/\partial\eta$ are known values taken from the solution of the stream-function equation (5). It is noted that (i) all given boundary conditions are imposed in an exact manner; and (ii) this technique only requires the local values of ψ and $\partial\psi/\partial\eta$ at the boundary node and its adjacent grid node to estimate the Dirichlet boundary conditions for the vorticity equation (6).

It can be seen that there are two values of u at each top corner of the cavity making the solution singular. In the well-known paper by Ghia et al. [30], the flow was simulated by the finite-difference scheme and a multigrid method using very fine grids (i.e. 129×129 and 257×257). The obtained results are very accurate and they have been considered as a benchmark of finite-difference methods. In the later work by Botella and Peyret [31], the regular and singular parts of the solution are handled by a Chebyshev collocation and an analytic method respectively. Benchmark spectral results for the flow at $Re = 100$ and $Re = 1000$ were reported. In the present study, the set of 2-node IRBFs is generated from grid lines that pass through interior grid nodes. As a result, the set of interpolation points does not include the top corners of the cavity and hence corner singularities do not explicitly enter the discrete system.

Simulation is carried out for a wide range of Re , namely (100, 400, 1000, 3200, 5000, 7500). Grid convergence is also studied. Results obtained are compared with the benchmark so-

lutions, i.e. Ghia et al. [30] and Botella and Peyret [31], and with the global 1D-IRBF collocation (1D-IRBF-C) results recently given in Mai-Duy and Tran-Cong [32]. These comparisons aim to assess the accuracy of the present method. To assess the efficiency and stability, an ADI method where streamfunction and vorticity are discretised by a three-node CD is also implemented. We denote this method as CD-ADI. It is noted that the same method of deriving computational vorticity boundary conditions is used in both IRBFE-ADI and CD-ADI methods.

Time-stepping convergence: The convergence behaviours of IRBFE-ADI and CD-ADI with respect to time are shown in Figures 4-6 and Table 1. It can be seen that solutions converge faster and larger time steps can be used for the present IRBFE-ADI method. The numbers of iterations are about 2.8×10^3 and 5.8×10^3 to reach $CM < 10^{-9}$ for IRBFE-ADI and CD-ADI respectively in the case of $Re = 1000$ and a grid of 51×51 (Figure 4). In the case of $Re = 3200$ and a grid of 91×91 (Figure 5), IRBFE-ADI takes about 2.1×10^4 to have $CM < 10^{-9}$ while CD-ADI requires about 7.4×10^4 to have the same CM . At $Re = 7500$, in contrast to the IRBFE-ADI method ($\Delta t = 1 \times 10^{-6}$), the CD-ADI method diverges even with a smaller time step of 5×10^{-7} as shown in Figure 6. The numbers of iterations in IRBFE-ADI method are generally lower than in CD-ADI method, yielding shorter computational time (Table 1). It is noted that the Thomas algorithm is used to solve tridiagonal systems in CD-ADI method and CPU seconds are associated with a computer which has 3.25 GB of RAM and one Intel(R) Core(TM)2 Duo CPU of 3.0 GHz. All codes are written in MATLAB[®] language.

Grid-size convergence: The convergence of extrema of the vertical and horizontal velocity profiles along the centrelines of the cavity with respect to grid refinement is presented in Table 2. Benchmark results by Ghia et al. [30] and Botella and Peyret [31] are also included for comparison purposes. It can be seen that errors relative to the benchmark results are consistency reduced as the grid is refined ($Re = 100, 1000$); and (ii) extrema values very close to the benchmark values are obtained with relatively coarse grids (e.g. 21×21 for $Re = 100$, 41×41 for $Re = 400$ and 61×61 for $Re = 1000$).

Solution quality: The solution qualities of IRBFE-ADI are shown in Table 2 and Figures 7-12. Table 2 reveals that the IRBFE-ADI results are the closest to the benchmark spectral solutions in comparison with the CD-ADI, the benchmark finite-difference and the global

1D-IRBF-C results. Errors relative to the benchmark spectral results are less than 1% for $Re = 100$ using a grid of 21×21 and for $Re = 1000$ using a grid of 61×61 . It can be seen from Figs 7-11 that smooth contours are obtained in the present IRBFE-ADI method for both the streamfunction and vorticity fields at relatively coarse grids. In Figure 7, the IRBFE-ADI method captures the primary vortex and the bottom-right corner eddy better than the CD-ADI method at $Re = 100$ and a grid of 11×11 . With the same grids, CD-ADI method yields oscillatory contours especially for the vorticity field as shown in Figures 8-10. Converged velocity profiles at $Re = 1000$ and $Re = 3200$ are obtained by IRBFE-ADI method with grids of 51×51 and 91×91 , respectively, as shown in Figure 12.

4.2. *Triangular cavity*

The proposed method is further verified through the simulation of steady recirculating flow in an equilateral triangle cavity. This is an example that presents a severe test for structured grid-based numerical methods [33, 34]. Figure 13 shows the geometry of the triangular cavity with the boundary conditions and the coordinate system. As in the square cavity flow problem, no-slip boundary condition is imposed on the left and right boundaries, while a unit horizontal velocity is prescribed on the top boundary. Numerical studies of this problem can be categorised into structured and unstructured grid/mesh-based methods. The former includes e.g. [33, 35, 36] where a finite difference method (FDM) was employed and the equilateral triangle had to be transformed to a computational domain on an isosceles right triangle. In the latter, Jyotsna and Vanka [34] used a multigrid procedure and a control volume formulation on triangular grids. They numerically verified interesting features of the flow in the Stokes regime. Kohno and Bathe [37] presented a flow-condition-based interpolation finite element scheme on triangular meshes to achieve solutions for high Reynolds numbers.

The imposition of boundary conditions for ω on the top is similar to that used in the square cavity flow, i.e. (55). On the left and right sides, analytic formulae for computing the vorticity boundary condition on a non-rectangular boundary [38] are utilised here

$$\omega_b = - \left[1 + \left(\frac{t_x}{t_y} \right)^2 \right] \frac{\partial^2 \psi_b}{\partial x^2}, \quad (56)$$

for a x -grid line, and

$$\omega_b = - \left[1 + \left(\frac{t_y}{t_x} \right)^2 \right] \frac{\partial^2 \psi_b}{\partial y^2}, \quad (57)$$

for a y -grid line. In (56) and (57), t_x and t_y are the x - and y -components of the unit vector tangential to the boundary. The approximations in (56) and (57) require information about ψ in one direction only and they are constructed here by means of 2-node IRBFEs, i.e. (17). No exterior/fictitious points as in [33] are involved here.

Four Cartesian grids, namely Grid 1 (1952 interior points), Grid 2 (2680 points), Grid 3 (3526 points) and Grid 4 (4486 points) as shown in Figure 14, are employed to study the convergence of the solution. Unlike FDMs [33, 35, 36], the present method does not require any coordinate transformation, making modelling simple. The flow is simulated at $Re = (0, 100, 200, 500)$ where $Re = U(H/3)/\nu$, U the lid velocity and H the cavity height (i.e. length AD in Figure 13). An alternative definition of Reynolds number was $Re_s = US/\nu$ where S is the cavity side length. We have $Re_s = 2\sqrt{3}Re$. For example, $Re = 500$ here is equivalent to $Re_s = 1732$.

Figures 15 and 16 present contour plots of the streamfunction and vorticity fields, the stream and iso-vorticity lines look comparable to those available in the literature (e.g. [33, 37]).

Figure 17 shows the profiles of u along the vertical centreline $x = 0$ and v along the horizontal line $y = 2$. Results obtained in [37] are also included for comparison purposes. It can be seen that the velocity profiles obtained by Grid 1 and Grid 2 at $Re = 100$, and by Grid 2 and Grid 3 at $Re = 200$ are almost identical. The present results are in good agreement with those by the flow-conditioned-based interpolation FEM for all values of Re . The profiles of v near the the stagnant corner at different Reynolds numbers also confirm the Stokes flow assumption of the flow field in this region (i.e. [39]).

4.3. Discussion

4.3.1. Comparison with other RBF techniques

To the best of our knowledge, the present two-node RBF scheme yields the smallest RBF set ever used for constructing approximation. When compared with other local RBF techniques, the present method results in remarkably sparse and banded system matrices and C^2 -

continuous solutions rather than the usual C^0 -continuous solutions. C^2 -continuous stream-function field leads to smooth and highly accurate velocity field.

Unlike other conventional RBF techniques, the present technique considers both the field variables and their partial derivatives in Cartesian directions in the formulation. As a result, the convection terms are naturally incorporated into the system matrices as unknowns and diagonally dominant systems are always guaranteed. Numerical results show that the present technique is very stable for high Re flows without recourse to up-winding schemes. Although the present system matrices are much larger, bigger time steps can be used and hence a smaller number of iterations are required to obtain a steady state solution. The computational time hence becomes competitive to those required by the conventional techniques as shown in Table 1.

4.3.2. Comparison with other conventional discretisation techniques

In terms of geometric modelling, unlike pseudo-spectral and finite-difference methods, the present Cartesian-grid technique can handle irregular domains well. In contrast to finite element and finite volumes, the pre-processing here is much more economical. Non-boundary grid points are trivially generated and the intersections between grid lines and the domain boundary can be determined as in a typical FE mesh generation [40]. For example, the intersections of an x -grid line with the boundary can be found as follows. Either an xy -plane or an xz -plane passing through the x -grid line is used and the intersection between this plane and the boundary (a curve or curves) is determined by analytic geometric methods. The intersection between the grid line and the curve(s) can then be easily determined.

In terms of solution, the governing equations are collocated at nodal points without the need for numerical integration. ADI solution strategies are conveniently applied to accelerate the computational process. When compared with low-order techniques, the present technique can produce accurate results on a relatively coarse mesh, therefore has the ability to reduce computational effort for a given accuracy. On the other hand, with more information (first derivatives are also involved in the interpolation) the cost to construct the approximation is quite higher than the standard CD scheme. In addition, the system matrix on each grid line of the present ADI method is not as sparse as and twice the size of those obtained by the PR-ADI method, requiring more than twice the memory storage. However, works on optimising

the proposed ADI method are under investigation. It is pointed out that we can approach tridiagonal system matrices with the same size as those in PR-ADI for diffusion-convection type equations.

5. Concluding remarks

We propose a C^2 -continuous alternating direction implicit solution method for solving the streamfunction-vorticity formulation governing fluid flows. Numerical experiments are conducted with problems on rectangular and non-rectangular domains. The method successfully simulates the fluid flows considered in a wide range of Reynolds numbers. Attractive features of the proposed methods include (i) simple preprocessing (Cartesian grids); (ii) a sparse system matrix (2-node approximations); and a higher order of continuity across grid nodes (C^2 -continuous elements). Numerical results show that (i) larger time steps can be used and smaller numbers of iterations are required in comparison with the classical CD-ADI method; and (ii) smooth solutions and high levels of accuracy are achieved using relatively coarse grids.

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Table 1: Lid-driven cavity flow: computational times.

Re	Method	Grid	Δt	Cycle seconds	Number of cycles	CPU seconds
100	CD-ADI	11x11	4.0E-3	2.7E-4	198	5.3E-2
	IRBFE-ADI	11x11	4.0E-3	1.7E-4	110	1.9E-2
	CD-ADI	21x21	1.0E-3	9.5E-4	413	0.393
	IRBFE-ADI	21x21	1.0E-3	8.9E-4	276	0.246
400	CD-ADI	21x21	2.0E-4	9.3E-4	1144	1.066
	IRBFE-ADI	21x21	4.0E-4	9.6E-4	452	0.432
	CD-ADI	31x31	2.0E-4	2.1E-3	922	1.945
	IRBFE-ADI	31x31	3.0E-4	2.8E-3	588	1.654
	CD-ADI	41x41	2.5E-4	4.1E-3	1568	6.378
	IRBFE-ADI	41x41	2.5E-4	5.8E-3	693	3.997
1000	CD-ADI	31x31	2.0E-5	2.2E-3	8884	19.853
	IRBFE-ADI	31x31	6.0E-5	2.7E-3	2851	7.785
	CD-ADI	41x41	2.0E-5	3.9E-3	8681	33.510
	IRBFE-ADI	41x41	6.0E-5	5.4E-3	2853	15.451
	CD-ADI	51x51	2.0E-5	6.6E-3	8447	55.910
	IRBFE-ADI	51x51	6.0E-5	1.2E-2	2807	33.880

Table 2: Lid-driven cavity flow: extrema of the vertical and horizontal velocity profiles along the centrelines of the cavity. % denotes percentage errors relative to the benchmark spectral results [31]. Results of the global 1D-IRBF-C, FDM and Benchmark are taken from [41], [30] and [31] respectively.

Re	Method	Grid	u_{min}	%	v_{max}	%	v_{min}	%
100	CD-ADI	11x11	-0.15242	28.79	0.10823	39.73	-0.14355	43.44
	IRBFE-ADI	11x11	-0.19916	6.95	0.14276	20.50	-0.18436	27.36
	CD-ADI	21x21	-0.19725	7.84	0.16069	10.51	-0.22401	11.74
	IRBFE-ADI	21x21	-0.21537	0.62	0.17932	0.14	-0.24960	0.78
	FDM ($\psi - \omega$)	129x129	-0.21090	1.47	0.17527	2.40	-0.24533	3.34
	Benchmark		-0.21404		0.17957		-0.25380	
400	CD-ADI	21x21	-0.20572		0.16693		-0.25885	
	IRBFE-ADI	21x21	-0.29015		0.24953		-0.36327	
	CD-ADI	31x31	-0.27258		0.24391		-0.37199	
	IRBFE-ADI	31x31	-0.32166		0.29581		-0.43622	
	CD-ADI	41x41	-0.29689		0.26978		-0.40808	
	IRBFE-ADI	41x41	-0.32780		0.30305		-0.44986	
FDM($\psi - \omega$)	129x129	-0.32726		0.30203		-0.44993		
1000	CD-ADI	31x31	-0.26073	32.90	0.24723	34.41	-0.36708	30.36
	1D-IRBF-C	31x31	-0.34791	10.46	0.33580	10.91	-0.46765	11.27
	IRBFE-ADI	31x31	-0.33775	13.08	0.32592	13.54	-0.44434	15.70
	CD-ADI	41x41	-0.30741	20.89	0.29382	22.05	-0.42451	19.46
	1D-IRBF-C	41x41	-0.37122	4.47	0.35910	4.73	-0.50168	4.82
	IRBFE-ADI	41x41	-0.37334	3.92	0.36193	3.98	-0.49863	5.40
	CD-ADI	51x51	-0.33242	14.45	0.31932	15.29	-0.45556	13.57
	1D-IRBF-C	51x51	-0.37985	2.25	0.36781	2.42	-0.51469	2.35
	IRBFE-ADI	51x51	-0.38482	0.97	0.37336	0.95	-0.51831	1.66
	CD-ADI	61x61	-0.34772	10.51	0.33502	11.12	-0.47488	9.90
	1D-IRBF-C	61x61	-0.38366	1.26	0.37173	1.38	-0.52029	1.29
	IRBFE-ADI	61x61	-0.38886	0.08	0.37719	0.07	-0.52537	0.32
	FDM ($\psi - \omega$)	129x129	-0.38289	1.46	0.37095	1.59	-0.51550	2.20
	Benchmark		-0.38857		0.37694		-0.52708	

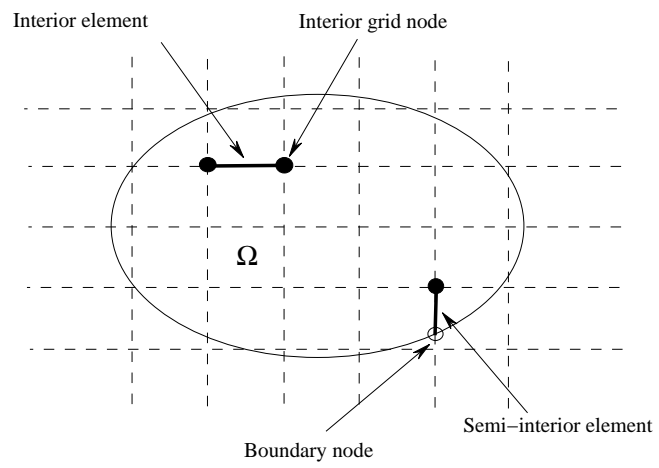
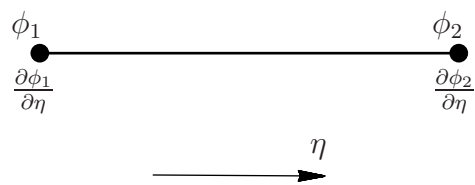


Figure 1: A domain is embedded in a Cartesian grid with interior and semi-interior elements.

(a) Interior element



(b) Semi-interior element

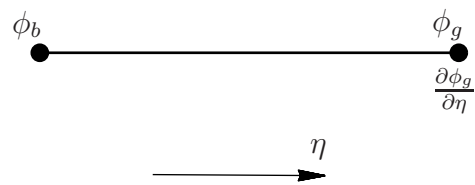


Figure 2: Schematic outline for 2-node IRBFs.

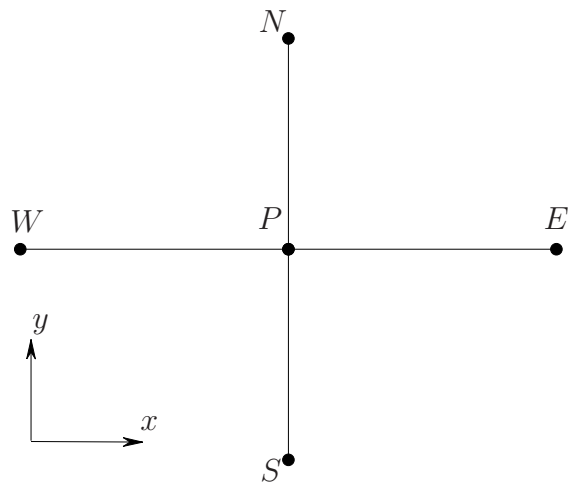


Figure 3: A grid point P and its neighbouring points on a Cartesian grid.

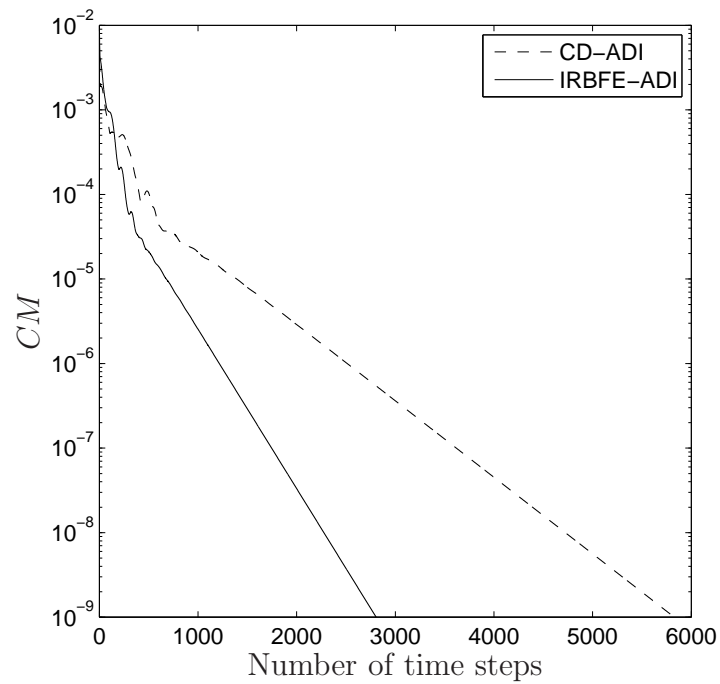


Figure 4: Lid-driven cavity flow, $Re = 1000$, 51×51 : convergence behaviour. IRBFE-ADI method using a time step of 6×10^{-5} converges faster than the CD-ADI method using a time step of 3×10^{-5} . It is noted that the latter diverges for time steps greater than 3×10^{-5} . CM denotes the relative norm of the difference of the streamfunction fields between two successive time levels.

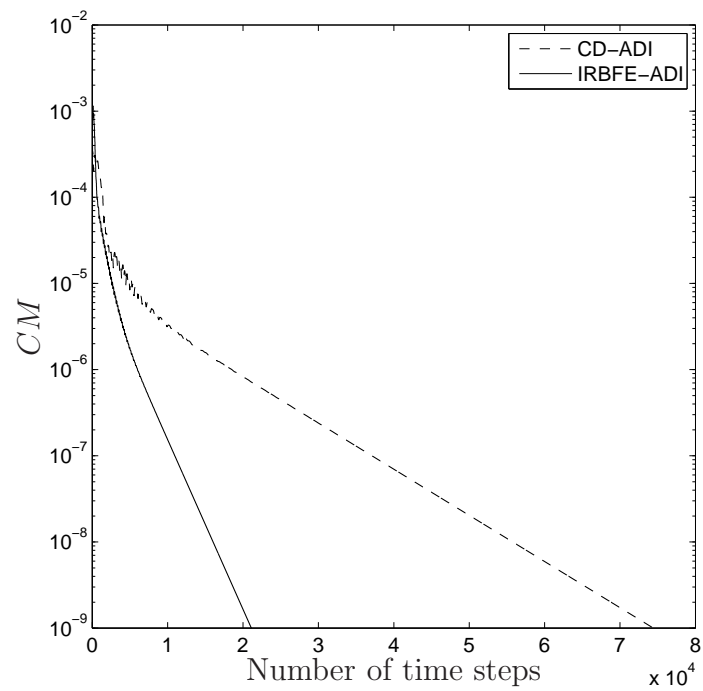


Figure 5: Lid-driven cavity flow, $Re = 3200$, 91×91 , solution at $Re = 1000$ used as initial guess: convergence behaviour. IRBFE-ADI method using a time step of 7×10^{-6} converges faster than the CD-ADI method using a time step of 2×10^{-6} . It is noted that the latter diverges for time steps greater than 2×10^{-6} . CM denotes the relative norm of the difference of the streamfunction fields between two successive time levels.

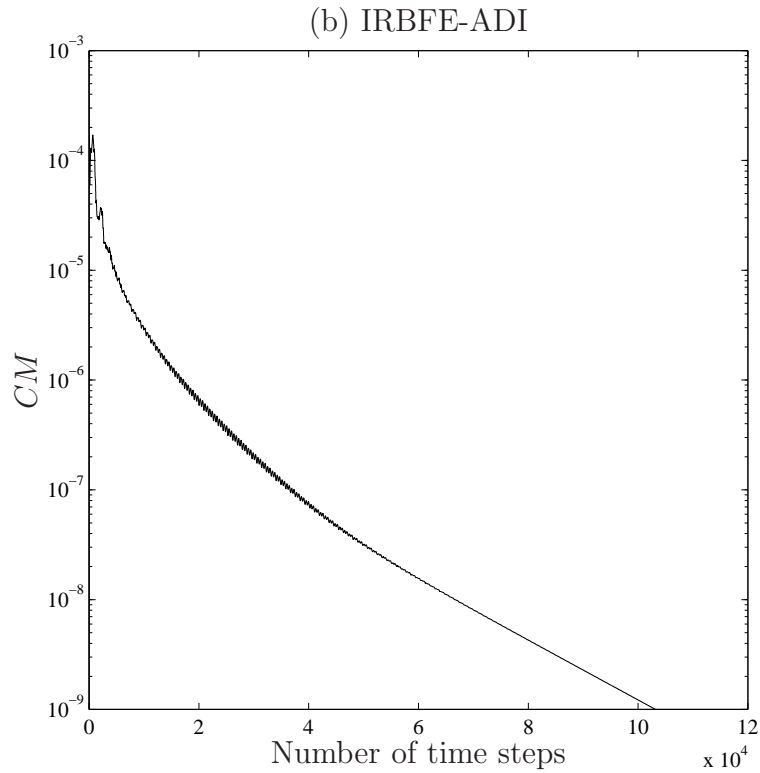
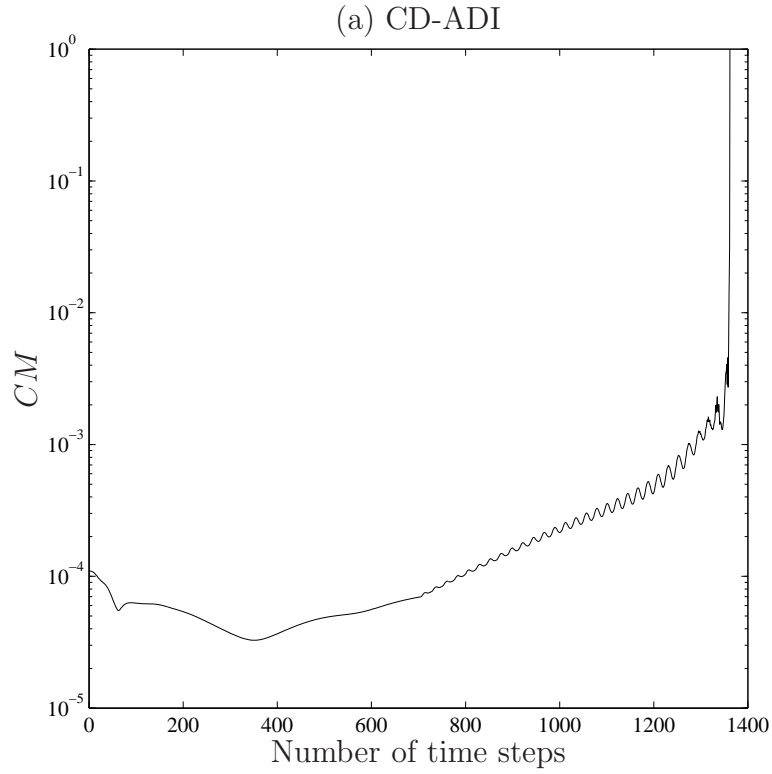


Figure 6: Lid-driven cavity flow, $Re = 7500$, 131×131 , solution at $Re = 5000$ used as initial guess: convergence behaviour. CD-ADI method uses a time step of 5×10^{-7} and IRBFE-ADI method uses a time step of 1×10^{-6} . CM denotes the relative norm of the difference of the streamfunction fields between two successive time levels.

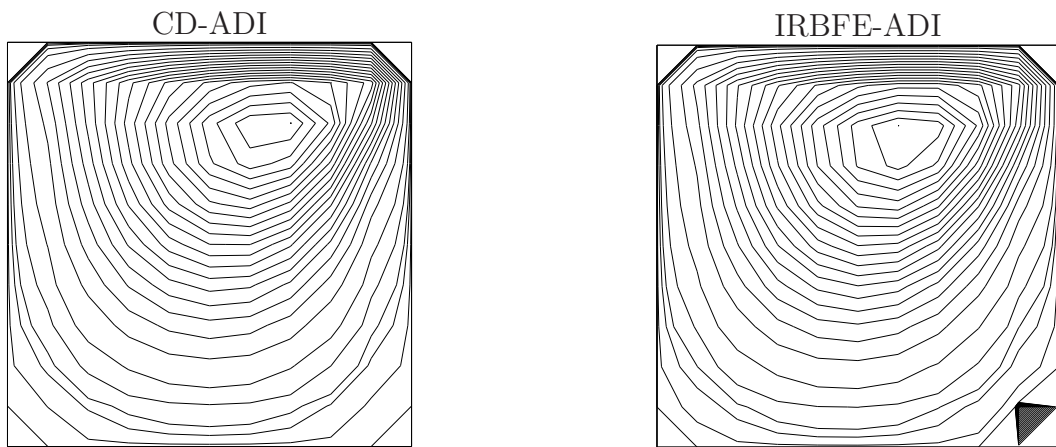


Figure 7: Lid-driven cavity flow, $Re = 100$, grid = 11×11 : streamlines. The contour values for CD-ADI and IRBFE-ADI plots are the same.

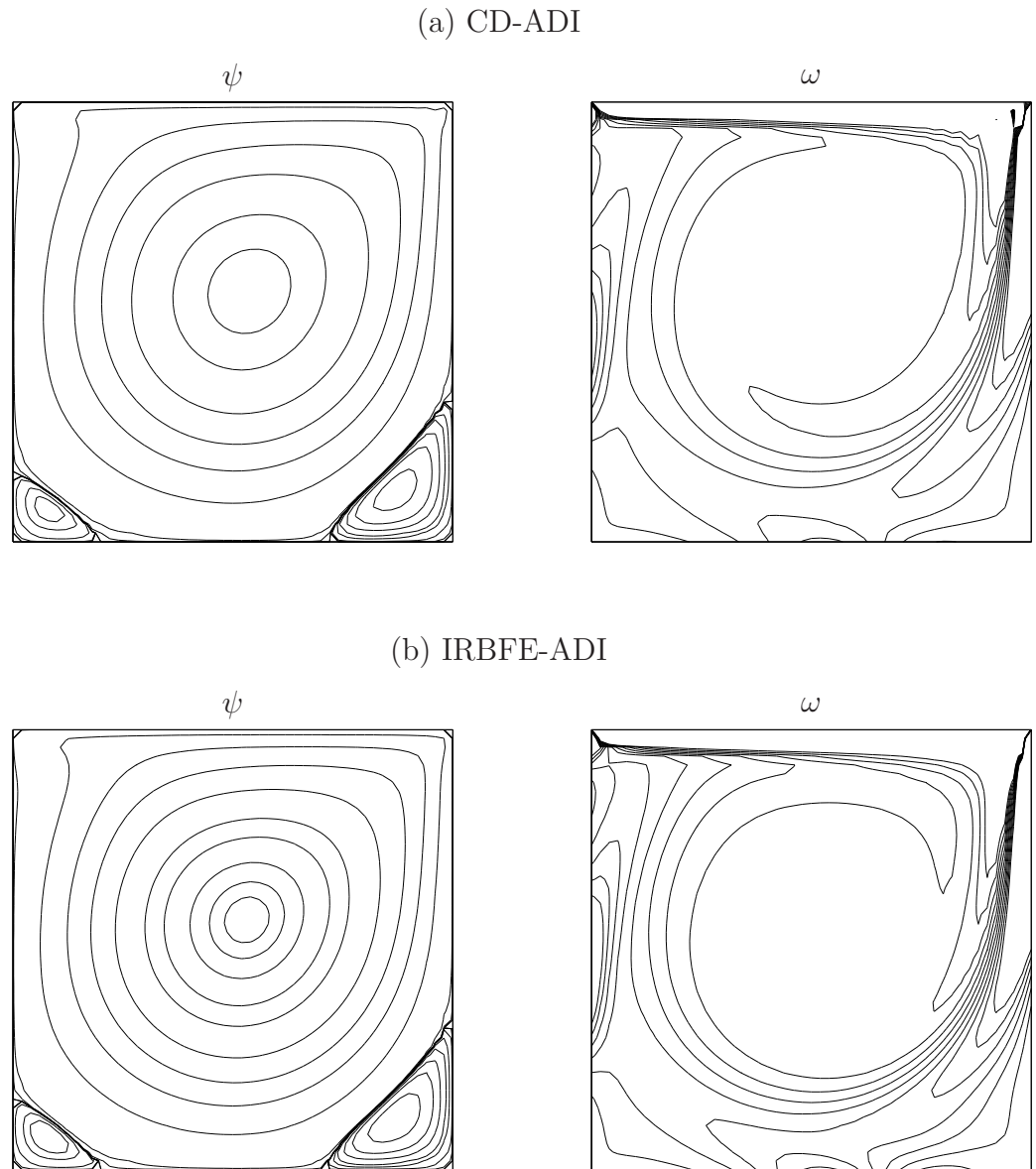


Figure 8: Lid-driven cavity flow, $Re = 1000$, grid = 51×51 : stream and iso-vorticity lines. The contour values are taken to be the same as those in [30] and [42] respectively. Note the oscillatory behaviour near the top right corner in the case of CD-ADI method.

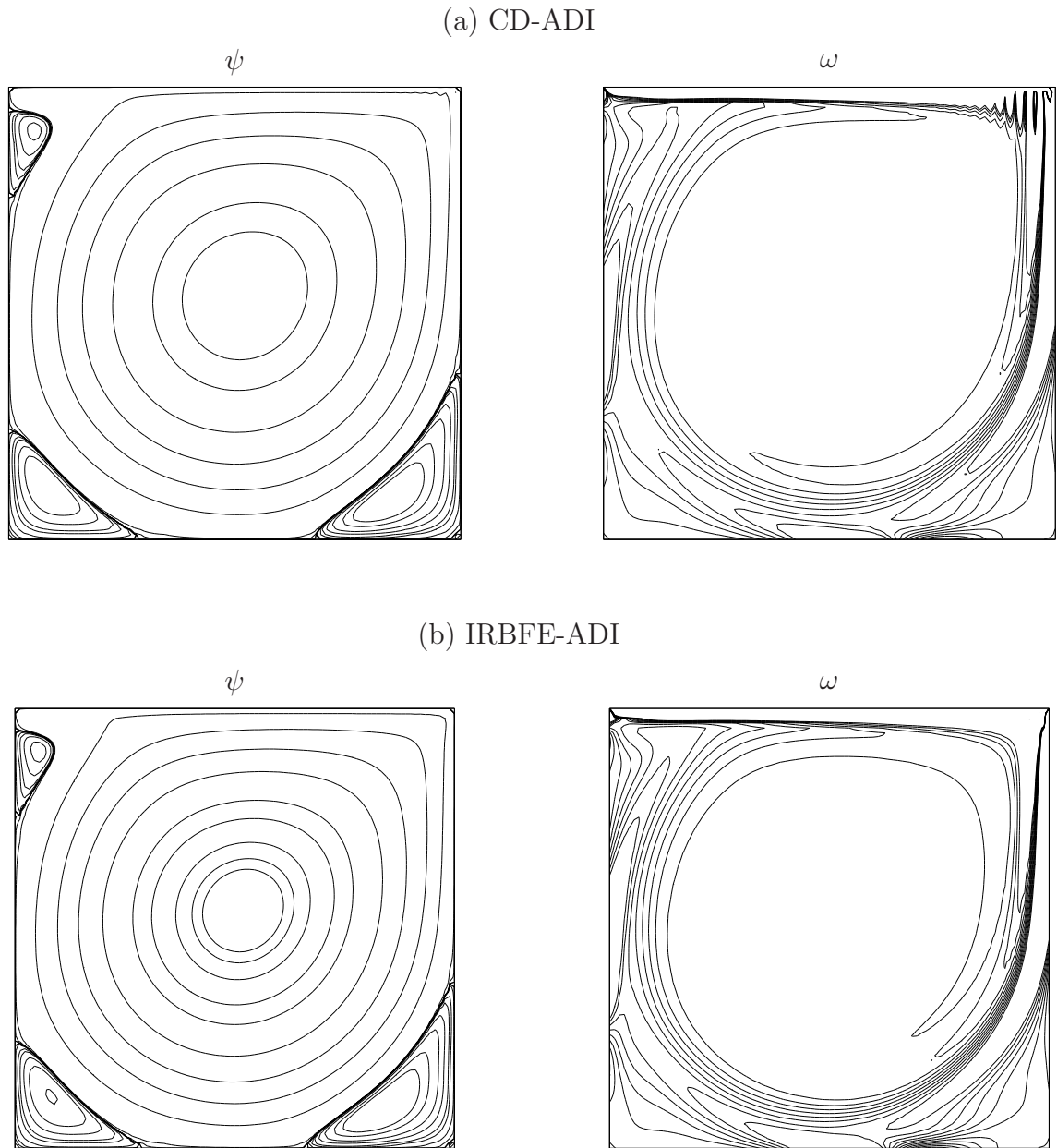


Figure 9: Lid-driven cavity flow, $Re = 3200$, grid = 91×91 : stream and iso-vorticity lines. The contour values are taken to be the same as those in [30] and [42] respectively. Note the oscillatory behaviour near the top right corner in the case of CD-ADI method.

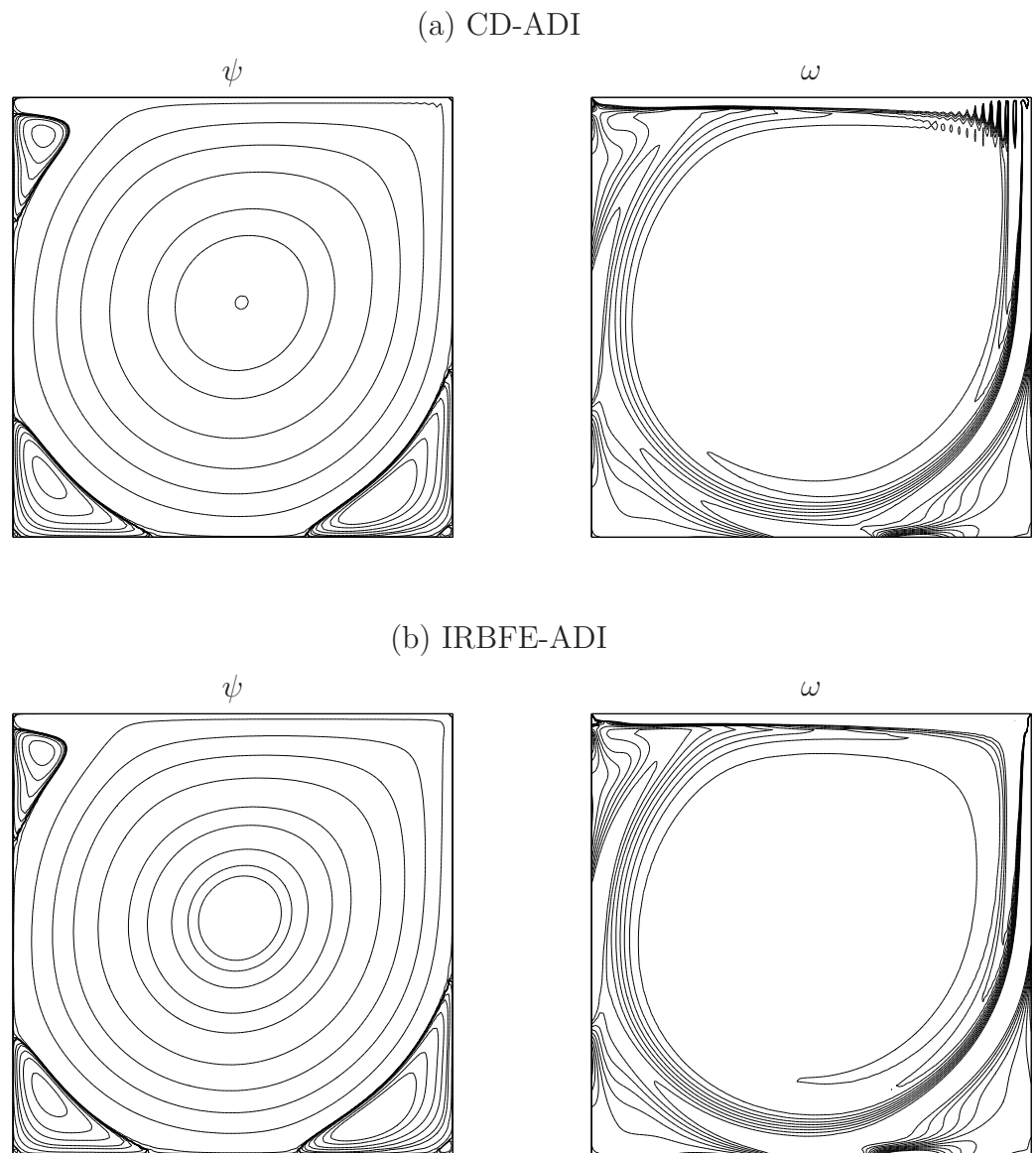


Figure 10: Lid-driven cavity flow, $Re = 5000$, grid = 111×111 : stream and iso-vorticity lines. The contour values are taken to be the same as those in [30] and [42] respectively. Note the oscillatory behaviour near the top right corner in the case of CD-ADI method.

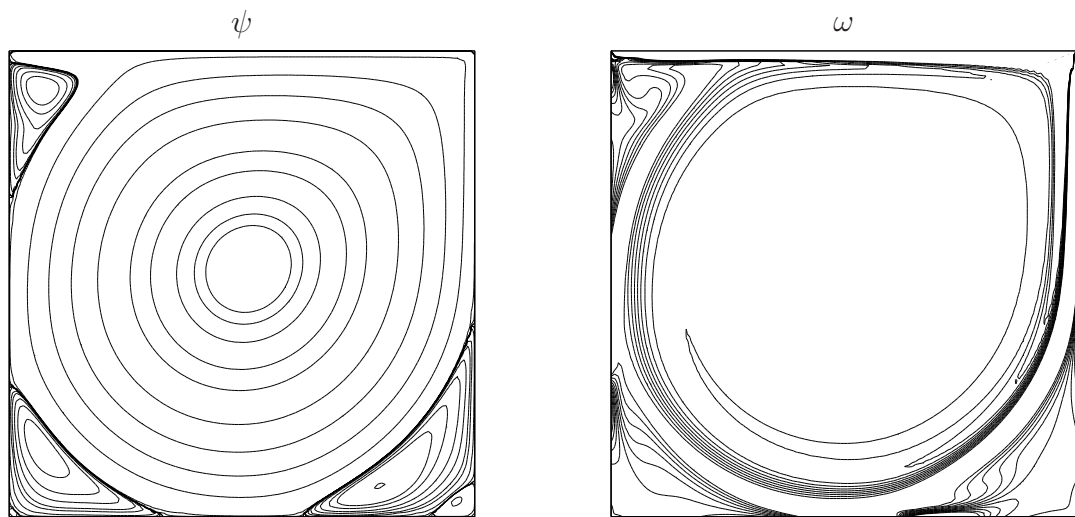
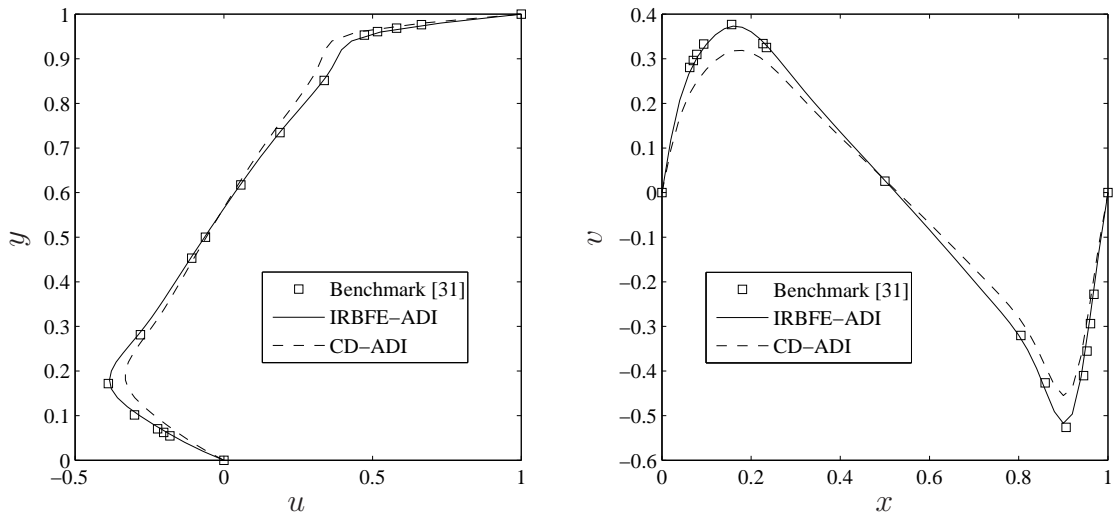


Figure 11: Lid-driven cavity flow, IRBFE-ADI, $Re = 7500$, grid = 131×131 : stream and iso-vorticity lines. The contour values are taken to be the same as those in [30] and [42] respectively.

(a) $Re = 1000$, grid= 51×51



(b) $Re = 3200$, grid= 91×91

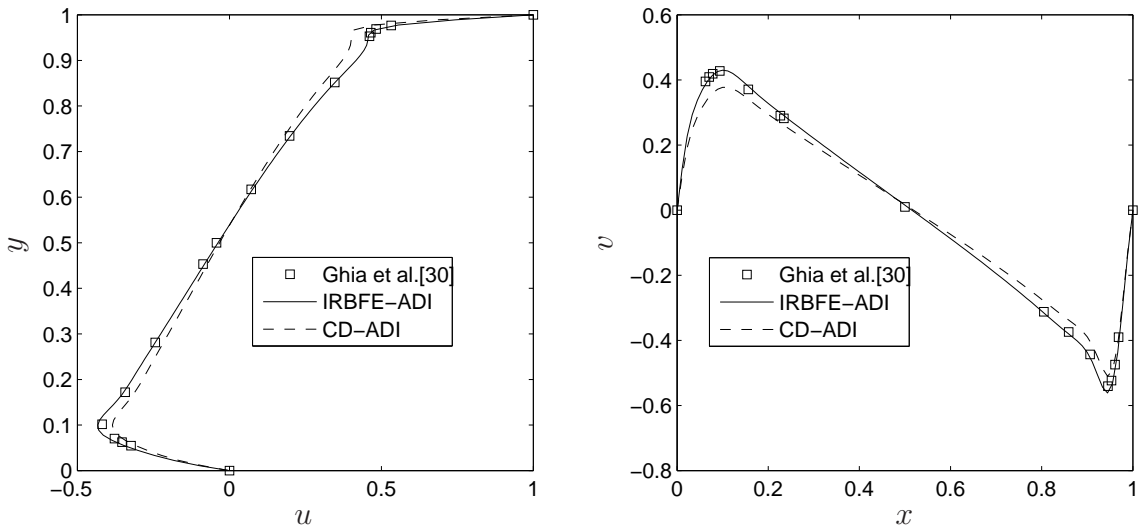


Figure 12: Lid-driven cavity flow: velocity profiles along the vertical and horizontal centrelines of the cavity.

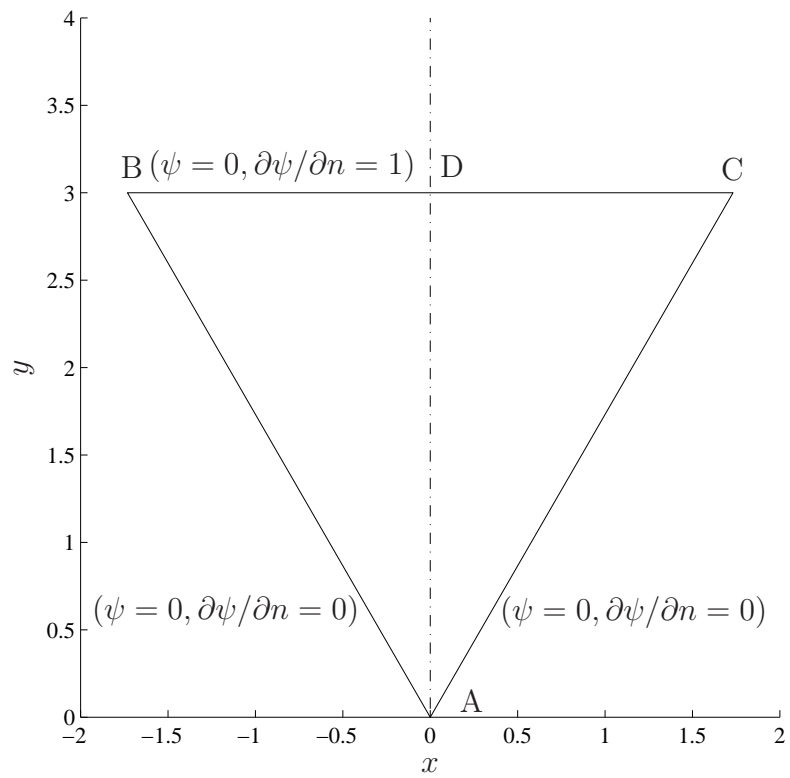


Figure 13: Triangular cavity flow: schematic outline of the computational domain and boundary conditions. Note that the characteristic length is chosen to be $AD/3$ to facilitate comparison with other published results [33, 37].

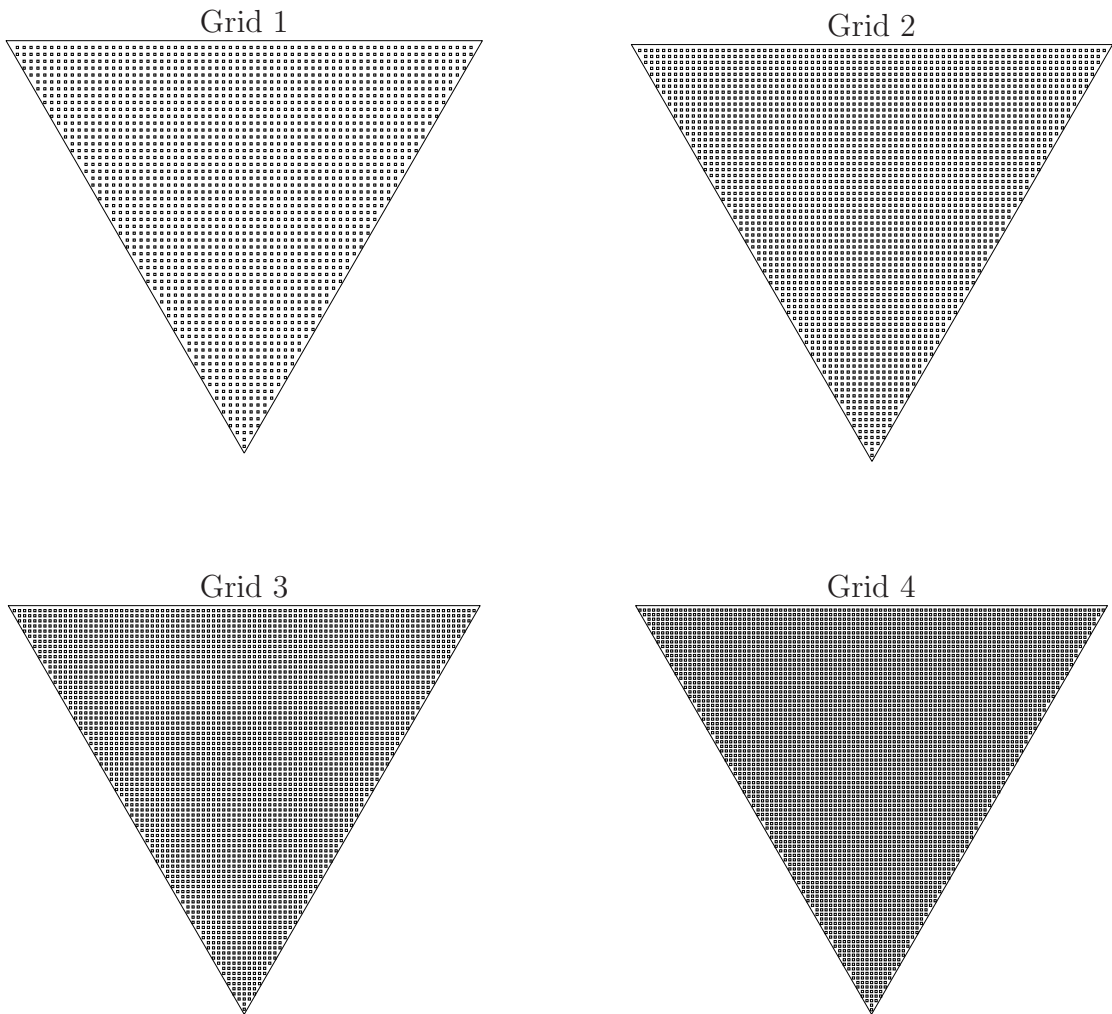


Figure 14: Triangular cavity flow: the computational domain is discretised by four Cartesian grids.

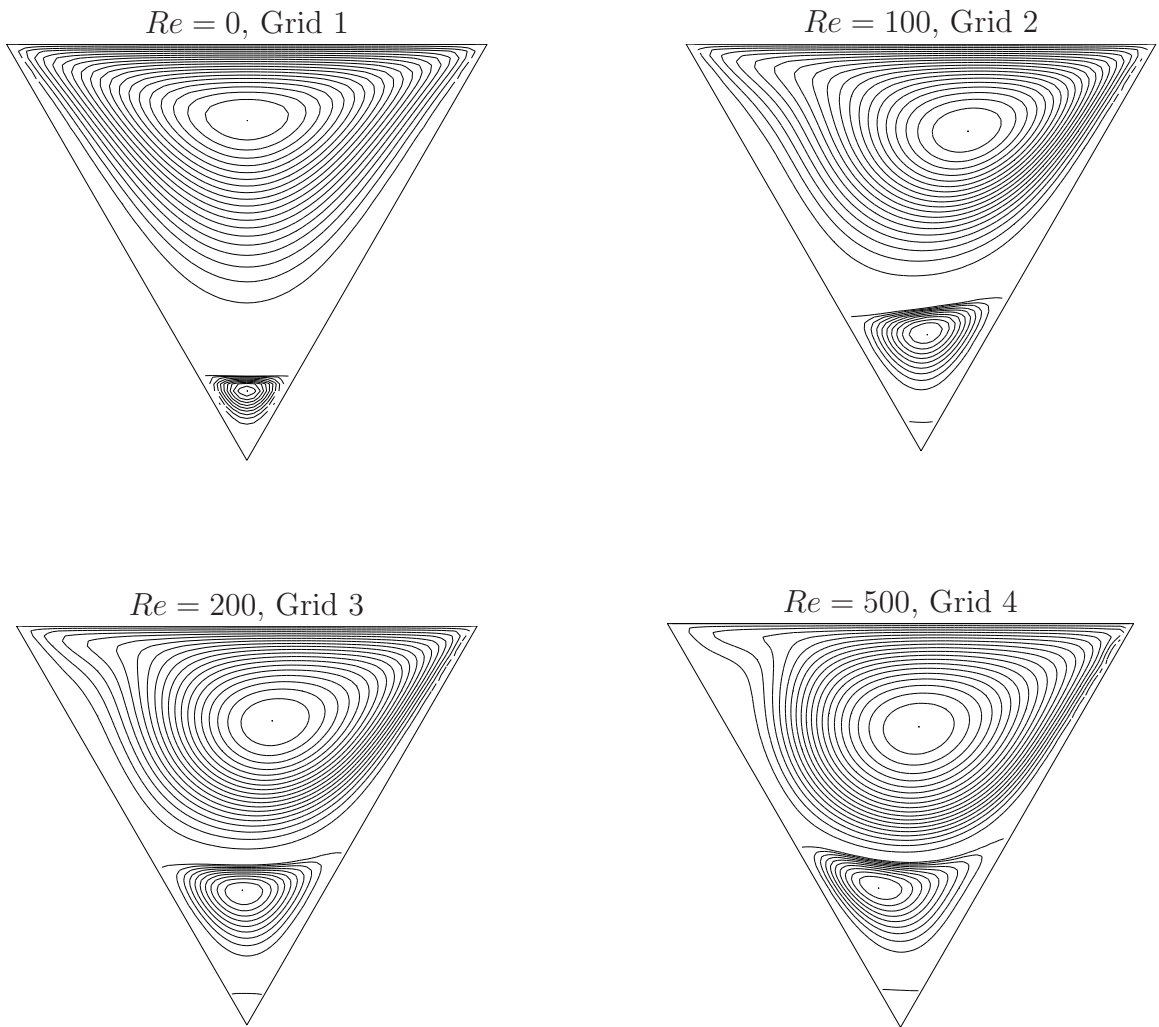


Figure 15: Triangular cavity flow: streamlines which are drawn using 21 equi-spaced levels between the minimum and zero values, and 11 equi-spaced levels between the zero and maximum values.

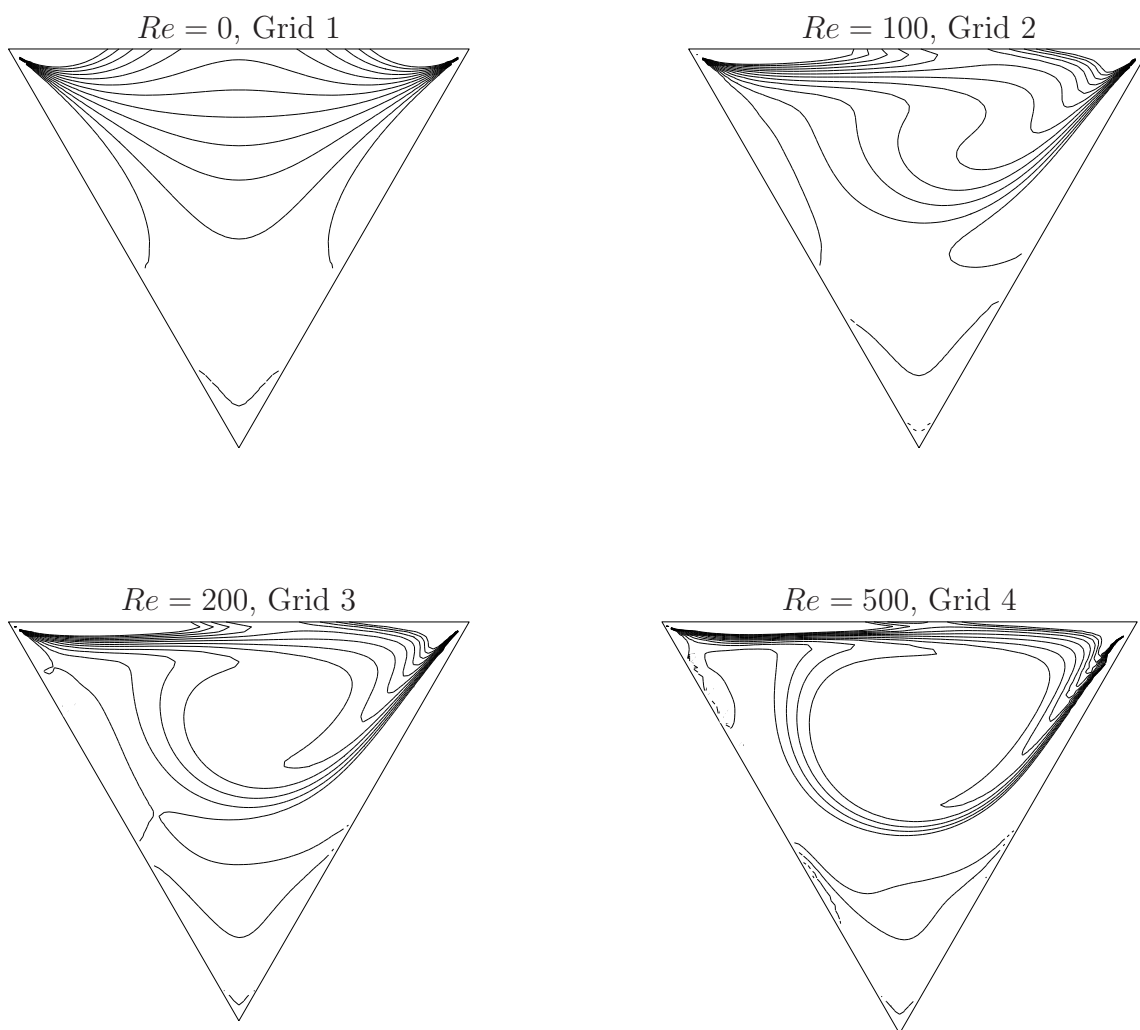


Figure 16: Triangular cavity flow: iso-vorticity lines which are drawn at intervals of $\Delta\omega = 0.5$ for a range of $-5 \leq \omega \leq 0.5$.

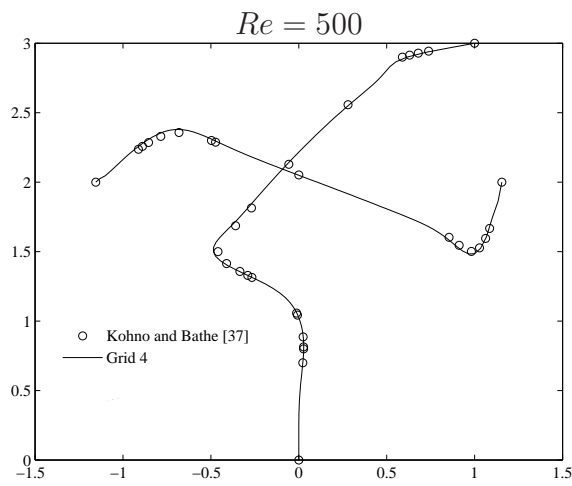
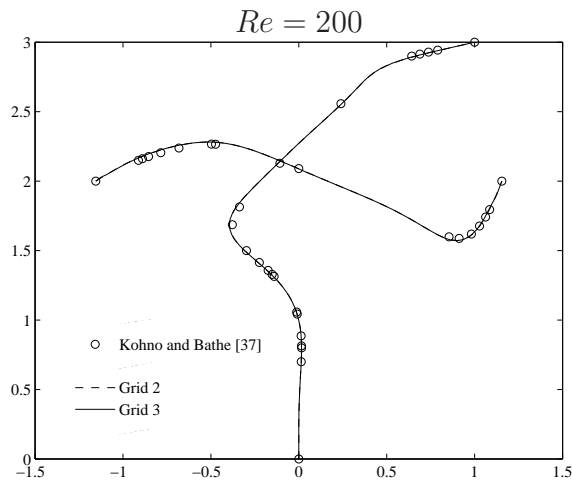
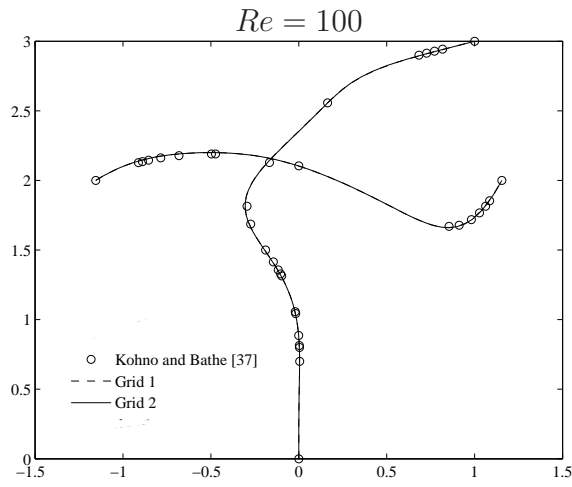


Figure 17: Triangular cavity flow: velocity profiles by the present method and the flow condition-based interpolation FEM [37].