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**ORIGINAL PAPER**

# A pseudo-compressible variational multiscale solver for turbulent incompressible flows

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**Abstract** In this work, we design an explicit time-stepping solver for the simulation of the incompressible turbulent flow through the combination of VMS methods and artificial compressibility. We evaluate the effect of the artificial compressibility on the accuracy of the explicit formulation for under-resolved LES simulations. A set of benchmarks have been solved, e.g., the 3D Taylor–Green vortex problem in turbulent regimes. The resulting method is proven to be an effective alternative to implicit methods in some application ranges (in terms of problem size and computational resources), providing comparable results with very low memory requirements. As an example, with the explicit approach, we are able to solve accurately the Taylor-Green vortex benchmark in a fine mesh with $512^3$ cells on a 12 cores 64 GB ram machine.

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## 1 Introduction

The presence of turbulence, which is always produced at high Reynolds numbers, poses a challenge for the numerical simulation of the Navier–Stokes equations. Capturing all the scales of the turbulent flow, known as the "Direct Numerical Simulation" (DNS) [28], or even only the larger scales in "Large Eddy Simulation" (LES), involve large scale problems on very fine grids and small time steps [31]. Taking into account the small-scale flow features in both time and space, explicit methods with small time steps is an appealing choice. In this sense, it is reasonable to envisage an increasing interest on matrix-free schemes as researchers will struggle with the draconian memory requirements of the forthcoming supercomputers.

Grid-based methods can be combined with explicit time integration when supplemented with an artificial compressibility method. Intensive research in this area has been done for laminar flows, within the framework of finite difference methods [9], finite volume methods [26], or the characteristic-based splitting methods [29]. In this work, we will develop explicit finite element (FE) schemes for turbulent flows. For FE schemes, the first problem when discretising the Navier–Stokes equations using FE methods is the fact that traditional Galerkin-based FE schemes are not suitable for convection dominated flows when dealing with under-resolved simulations. The second problem is the compatibility requirement between the velocity and pressure FE spaces for stability purposes, which is known as the LBB condition. The variational multiscale method (VMS) framework is introduced to overcome these difficulties [22]. The VMS approach can be understood as a residual-based stabilised FE method, and so, the singularly perturbed nature of the Navier-Stokes equations is stabilised as the Reynolds number increases. On the other hand, it

allows one to avoid the compliance of the compatibility conditions (the so-called discrete inf-sup) between velocity and pressure spaces, and so, we are able to use equal-order interpolation.

The VMS concept is an LES turbulent model for the Navier–Stokes equations, as suggested in [10], exploited in [23], and analytically justified in [21]. One further development of the VMS framework was to consider the subscale proportional to the orthogonal component of the residual with respect to the FE spaces, leading to the orthogonal subscales VMS technique [14]. Another improvement was to consider a transient evolution equation for the subscale component, leading to the dynamic subscales in [13]. In these situations, the projection-based VMS method uses variational projections similar to the traditional filtered equations, which provides a numerically oriented turbulent modelling feature. Compared with classical LES-type filters, the VMS method does not have the problem associated to inhomogeneous non-commutative filters for wall-bounded flows. Further, the VMS method retains numerical consistency in the FE equations, whereas Smagorinsky-type models introduce an error of order $h^{\frac{4}{3}}$. Good numerical results have been obtained with this approach, which demonstrate its capability to compete with traditional LES turbulence modelling approaches (see e.g. [5,16]).

The numerical solution of the VMS discretisation of incompressible flow is traditionally obtained using either a pressure segregation scheme [1] or monolithic approaches [5], which require to solve linear systems and (possibly) nonlinear iterations. For large scale problems, the linear system is solved via Krylov iterative solvers. In order to solve large scale problems with implicit methods, the main requirement is the use of weakly scalable preconditioning techniques [3,4,30].

So far, VMS methods have been restricted to implicit time stepping techniques. However, explicit methods (or matrix-free lightweight implicit methods) are favoured in some situations, e.g., when one wants to tackle larger problems with limited computational resources. As a result, in this paper we investigate the possibility of imposing an explicit time-stepping approach for the simulation of the incompressible turbulent flow through the combination of VMS methods and artificial compressibility. We show that the resulting scheme, after the artificial compressibility perturbation, can still provide accurate under-resolved LES simulations of turbulent flows with a very limited memory consumption. The proposed scheme can be straightforwardly applied to structured and unstructured meshes. In particular, (multi-)linear equal-order velocity and pressure interpolation is used in this work. Compared to other explicit methods, the built-in sub-grid component acts as an implicit LES turbulence model and no extra turbulent modelling is required.

This paper is split into the following sections. In Sect. 2, the governing equations of the Navier–Stokes problem and its artificial compressibility formulation are presented. Section 3 combines the VMS formulation with the artificial compressibility technique. Section 4 presents a range of numerical results in order to demonstrate the aspects described in previous sections, including the 2-D cavity flow, 2-D plane mixing layer, 3-D cavity flow and 3-D Taylor Green vortex. Finally, some concluding remarks are made in Sect. 5.

## 2 Governing equations

### 2.1 Incompressible Navier–Stokes equations

Let $\Omega$ be a bounded domain of $\mathbb{R}^d$ where $d = 2$ or 3 is the spatial dimension, with boundary $\Gamma = \partial\Omega$, and let the time interval be $[0, T]$. The strong form of the incompressible Navier–Stokes problem consists of finding the velocity field $\boldsymbol{u}$ and pressure $p$ such that

$$\begin{aligned} \partial_t \boldsymbol{u} - \nu\nabla^2\boldsymbol{u} + \boldsymbol{u}\cdot\nabla\boldsymbol{u} + \nabla p &= \boldsymbol{f} \quad \text{in } \Omega\times(0,T), \\ \nabla\cdot\boldsymbol{u} &= 0 \quad \text{in } \Omega\times(0,T), \end{aligned} \tag{1}$$

where $\boldsymbol{f}$ is the force vector and $\nu > 0$ is the kinematic viscosity. System (1) has to be supplied with appropriate boundary and initial conditions. The boundary $\Gamma$ is divided into the Dirchilet boundary $\Gamma_D$ and Neumann boundary $\Gamma_N$, satisfy $\Gamma_D \cup \Gamma_N = \Gamma$ and $\Gamma_D \cap \Gamma_N = \varnothing$. Then, the boundary and initial conditions read:

$$\boldsymbol{u} = \bar{\boldsymbol{u}} \quad \text{on } \Gamma_D\times(0,T), \tag{2a}$$

$$(-p\boldsymbol{I} + \nu(\nabla u + (\nabla u)^T))\cdot\boldsymbol{n} = \bar{\boldsymbol{t}} \quad \text{on } \Gamma_N\times(0,T), \tag{2b}$$

$$\boldsymbol{u} = \boldsymbol{u}_0 \quad \text{in } \Omega\times\{0\}, \tag{2c}$$

where $\boldsymbol{n}$ is the unit outward vector normal to the surface $\Gamma$. To simplify the exposition, we consider the pure Dirichlet boundary $\Gamma_D \equiv \Gamma$.

### 2.2 Artificial compressibility method

The artificial compressibility formulation replaces the divergence free constraint by adding a pseudo-time derivative of the pressure. System (1) is modified to

$$\begin{aligned} \partial_t\boldsymbol{u} - \nu\nabla^2\boldsymbol{u} + \boldsymbol{u}\cdot\nabla\boldsymbol{u} + \nabla p &= \boldsymbol{f} \quad \text{in } \Omega\times(0,T) \\ \partial_t p + \varepsilon^{-1}\nabla\cdot\boldsymbol{u} &= 0 \quad \text{in } \Omega\times(0,T) \end{aligned} \tag{3}$$

In (3), $\varepsilon^{-1}$ is the artificial compressibility coefficient or pseudo compressibility coefficient, where an artificial sound speed can be evaluated as $c = \sqrt{\varepsilon^{-1}}$. As the pressure

wave speed $c$ reaches a very large number, (3) approximates (1). A relaxation of the incompressibility will allow one to use an explicit time stepping scheme. However, a large artificial pressure wave speed will be a limit factor of the explicit time stepping. The characteristic velocity $\varepsilon^{-1} = c^2 = (\text{ASS}||\boldsymbol{u}||_\infty)^2$, where $||\boldsymbol{u}||_\infty$ is the maximum norm of the velocity field, and ASS is the artificial sound speed coefficient usually takes a value from 1 to 10. For the steady state problem, the coefficient will not effect the final solution because the residual will convergence to 0. The effect of ASS for turbulent flows will be investigated in the transient numerical example, 2D plane mixing layer and 3D Taylor Green vortex. The numerical experiments indicate that values of ASS 5 provide a good enough satisfaction of the incompressibility constraint.

Notice that the pressure behaviour in (3) changes from elliptic to hyperbolic (wave-propagation) character with the artificial compressibility method, and the Courant–Friedrichs–Lewy (CFL) condition is a necessary condition for temporal stability. In particular, for the problem considered the time step size $\Delta t$ has to be such that

$$\Delta t \leq \alpha \frac{\Delta h}{c},$$

where $\alpha$ is a factor that depends on the time integration scheme being used and the safety margin desired, and $\Delta h$ is the minimum mesh size. In what follows, we will assume that the time discretisation is uniform, with time step size $\Delta t$. Time step levels will be denoted with a superscript.

### 2.3 Weak formulation and notation

Let us introduce some notation. $L^2(\Omega)$ denotes the spaces of functions that are square integrable in $\Omega$

$$L^2(\Omega) \doteq \left\{ u \colon \Omega \to \mathbb{R} \,\middle|\, \int_\Omega u^2 \, d\Omega < \infty \right\}.$$

It is a Hilbert space with scalar product

$$(u, v)_\Omega \equiv (u, v) \doteq \int_\Omega u(\boldsymbol{x}) v(\boldsymbol{x}) \, d\Omega. \tag{4}$$

When the integral is performed over a subdomain $\omega$, we will denote it by $(\cdot, \cdot)_\omega$. $H^1(\Omega)$ denotes the Sobolev space of square integrable functions with square integral derivative as

$$H^1(\Omega) \doteq \{ u \colon \Omega \to \mathbb{R} \mid u, \nabla u \in L^2(\Omega) \}.$$

We define the solution $\boldsymbol{H}^1_E(\Omega)$ space and the test space $\boldsymbol{H}^1_0(\Omega)$ by

$$\boldsymbol{H}^1_E(\Omega) \doteq \{ \boldsymbol{u} \in H^1(\Omega)^d \mid \boldsymbol{u} = \bar{\boldsymbol{u}} \text{ on } \Gamma \},$$
$$\boldsymbol{H}^1_0(\Omega) \doteq \{ \boldsymbol{u} \in H^1(\Omega)^d \mid \boldsymbol{u} = \boldsymbol{0} \text{ on } \Gamma \},$$

respectively. Moreover, $\boldsymbol{H}^{-1}(\Omega)$ denote the topological dual of $\boldsymbol{H}^1_0(\Omega)$ and $\langle \cdot, \cdot \rangle$ the duality pairing between $\boldsymbol{H}^{-1}(\Omega)$ and $\boldsymbol{H}^1_0(\Omega)$. Then, the standard weak form of (3) is the following: for each time $t$, find a velocity $\boldsymbol{u}(\cdot, t) \in \boldsymbol{H}^1_E(\Omega)$ and a pressure $p(\cdot, t) \in L^2(\Omega)$ such that

$$(\partial_t \boldsymbol{u}, \boldsymbol{v}) + \nu (\nabla \boldsymbol{u}, \nabla \boldsymbol{v}) - (p, \nabla \cdot \boldsymbol{v}) + \langle \boldsymbol{u} \cdot \nabla \boldsymbol{u}, \boldsymbol{v} \rangle = \langle \boldsymbol{f}, \boldsymbol{v} \rangle$$
$$\text{for all } \boldsymbol{v} \in \boldsymbol{H}^1_0(\Omega), \tag{5}$$
$$(\partial_t p, q) + \varepsilon^{-1} (\nabla \cdot \boldsymbol{u}, q) = 0 \quad \text{for all } q \in L^2(\Omega).$$

### 2.4 Galerkin FE method

Given a FE partition of the domain $\Omega$, the problem is to construct a velocity solution space $\boldsymbol{V}^E_h \subset \boldsymbol{H}^1_E(\Omega)$ and $L_h \subset L^2(\Omega)$ to set the discrete problem: for each time $t$, find a velocity $\boldsymbol{u}_h(\cdot, t) \in \boldsymbol{V}^E_h$ and a pressure $p(\cdot, t)_h \in L_h$ such that

$$(\partial_t \boldsymbol{u}_h, \boldsymbol{v}_h) + \nu (\nabla \boldsymbol{u}_h, \nabla \boldsymbol{v}_h) - (p_h, \nabla \cdot \boldsymbol{v}_h) + \langle \boldsymbol{u}_h \cdot \nabla \boldsymbol{u}_h, \boldsymbol{v}_h \rangle$$
$$= \langle \boldsymbol{f}, \boldsymbol{v}_h \rangle \tag{6}$$
$$(\partial_t p_h, q_h) + \varepsilon^{-1} (\nabla \cdot \boldsymbol{u}_h, q_h) = 0,$$

for all $\boldsymbol{v}_h \in \boldsymbol{V}^0_h$ and $q_h \in L_h$, where $\boldsymbol{V}^0_h$ is constructed as $\boldsymbol{V}^E_h$ but with homogeneous boundary conditions. As commented above, it is well known that (6) suffers from numerical instabilities for high Reynolds number problems. Also, the velocity and pressure solution space must satisfy the inf-sup condition in order to have a well-posed problem, with bounded pressure field. These difficulties can be handled by using the VMS method.

## 3 Explicit variational multiscale method

In general, the VMS method decomposes the solution $\boldsymbol{u}$ and $p$ on a large scale component and a subscale component. The large scale components $\boldsymbol{u}_h$ and $p_h$ are resolved by the FE mesh, whilst the subscale components $\tilde{\boldsymbol{u}}$ and $\tilde{p}$ are approximated by a certain analytical approach. Let us consider a scale decomposition of spaces $\boldsymbol{H}^1_E(\Omega)$ and $L^2(\Omega)$ such that $\boldsymbol{H}^1_E(\Omega) = \boldsymbol{V}_h \oplus \tilde{V}$ and $L^2(\Omega) = L_h \oplus \tilde{L}$, where $\tilde{V}$ and $\tilde{L}$ denote the infinite-dimensional spaces that complete the FE spaces $\boldsymbol{V}_h$ and $L_h$ to approximate the velocity and pressure in the standard Galerkin FE method. $(\cdot)_h$ and $\tilde{(\cdot)}$ denote the FE component and the sub-grid component, respectively. Thus, we approximate the velocity $\boldsymbol{u}$ and pressure $p$ by

$$\boldsymbol{u} \approx \boldsymbol{u}_h + \tilde{\boldsymbol{u}}, \quad p \approx p_h + \tilde{p},$$

where $\boldsymbol{u}_h \in \boldsymbol{V}_h$, $\tilde{\boldsymbol{u}} \in \tilde{\boldsymbol{V}}$, $p_h \in L_h$, $\tilde{p} \in \tilde{L}$ for each time $t$. Using this splitting in (5), yields

$$(\partial_t (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \boldsymbol{v}) + \nu (\nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \nabla \boldsymbol{v}) - (p_h + \tilde{p}, \nabla \cdot \boldsymbol{v})$$
$$+ \langle (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}_h) \cdot \nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}_h), \boldsymbol{v} \rangle = \langle \boldsymbol{f}, \boldsymbol{v} \rangle,$$
$$(\partial_t (p_h + \tilde{p}), q) + \varepsilon^{-1} (\nabla \cdot (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}_h), q) = 0. \quad (7)$$

### 3.1 FE scale problem

First we consider FE test function, i.e. $\boldsymbol{v} = \boldsymbol{v}_h \in \boldsymbol{V}_h^0$ and $q = q_h \in L_h$. The resolved large-scale component of (7) yields

$$(\partial_t (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \boldsymbol{v}_h) + \nu (\nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \nabla \boldsymbol{v}_h)$$
$$- (p_h + \tilde{p}, \nabla \cdot \boldsymbol{v}_h) + \langle (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \boldsymbol{v}_h \rangle = \langle \boldsymbol{f}, \boldsymbol{v}_h \rangle,$$
$$(\partial_t (p_h + \tilde{p}), q_h) + \varepsilon^{-1} (\nabla \cdot (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), q_h) = 0.$$

Let us elaborate on some of the terms of these equations:

- A very important ingredient of our approach is that *we choose the sub-grid spaces to be orthogonal in the $L^2(\Omega)$-sense to the FE spaces*. Therefore, the time derivative terms $(\partial_t \tilde{\boldsymbol{u}}, \boldsymbol{v}_h)$ and $(\partial_t \tilde{p}, q_h)$ vanish:

$$(\partial_t (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \boldsymbol{v}_h) = (\partial_t \boldsymbol{u}_h, \boldsymbol{v}_h);$$
$$(\partial_t (p_h + \tilde{p}), q_h) = (\partial_t p_h, q_h).$$

  We show below how to satisfy this orthogonality in practice.

- The viscous term $\nu (\nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \nabla \boldsymbol{v}_h)$ is simplified as

$$\nu (\nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \nabla \boldsymbol{v}_h) = \nu (\nabla \boldsymbol{u}_h, \nabla \boldsymbol{v}_h)$$
$$+ \nu \sum_K \left[ (-\tilde{\boldsymbol{u}}, \Delta \boldsymbol{v}_h)_K + (\tilde{\boldsymbol{u}}, \boldsymbol{n} \cdot \nabla \boldsymbol{v}_h)_{\partial K} \right]$$
$$= \nu (\nabla \boldsymbol{u}_h, \nabla \boldsymbol{v}_h).$$

  The term $(-\tilde{\boldsymbol{u}}, \Delta \boldsymbol{v}_h)_K$ vanishes for linear elements. The term $(\tilde{\boldsymbol{u}}, \boldsymbol{n} \cdot \nabla \boldsymbol{v}_h)_{\partial K}$ is usually neglected in practice, although it could be taken into account using the strategy proposed in [12].

- In order to avoid derivatives of the sub-grid component, the convective component can be re-written as

$$\langle (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), \boldsymbol{v}_h \rangle$$
$$= \langle (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla \boldsymbol{u}_h, \boldsymbol{v}_h \rangle - \langle \tilde{\boldsymbol{u}}, (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla \boldsymbol{v}_h \rangle,$$

  where we have used the fact that $\boldsymbol{u}_h + \tilde{\boldsymbol{u}}$ is divergence free when integrating by parts.

- Again to avoid derivatives of the sub-grid scales, the mass conservation equation is re-written as

$$\varepsilon^{-1} (\nabla \cdot (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}), q_h) = \varepsilon^{-1} (\nabla \cdot \boldsymbol{u}_h, q_h)$$
$$- \varepsilon^{-1} (\tilde{\boldsymbol{u}}, \nabla q_h) + \varepsilon^{-1} (\boldsymbol{n} \cdot \tilde{\boldsymbol{u}}, q_h)_{\partial \Omega},$$

after integration by parts. As before, the sub-grid scale boundary term is neglected.

Taking into account the previous considerations, the equations we have are:

$$(\partial_t \boldsymbol{u}_h, \boldsymbol{v}_h) + \langle \boldsymbol{a} \cdot \nabla \boldsymbol{u}_h, \boldsymbol{v}_h \rangle + \nu (\nabla \boldsymbol{u}_h, \nabla \boldsymbol{v}_h)$$
$$- (p_h, \nabla \cdot \boldsymbol{v}_h) - (\tilde{p}, \nabla \cdot \boldsymbol{v}_h) - \langle \tilde{\boldsymbol{u}}, \boldsymbol{a} \cdot \nabla \boldsymbol{v}_h \rangle = 0,$$
$$(\partial_t p_h, q_h) + \varepsilon^{-1} (\nabla \cdot \boldsymbol{u}_h, q_h) - \varepsilon^{-1} (\tilde{\boldsymbol{u}}, \nabla q_h) = 0,$$

where $\boldsymbol{a} \doteq \boldsymbol{u}_h + \tilde{\boldsymbol{u}}$. The problem will be closed once an approximation for the sub-grid scale velocity and pressure is proposed.

### 3.2 Local sub-grid scale problem

In this section, we discuss the sub-grid scale formulation and then give the analytical approximation of the solution. As a result, the orthogonal sub-grid scale in an explicit VMS method will be given. The sub-grid scale equation reads (in strong form):

$$\partial_t \tilde{\boldsymbol{u}} + (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla \tilde{\boldsymbol{u}} - \nu \Delta \tilde{\boldsymbol{u}} + \nabla \tilde{p} = \boldsymbol{R}_u,$$
$$\partial_t \tilde{p} + \varepsilon^{-1} \nabla \cdot \tilde{\boldsymbol{u}} = R_p,$$

where $\boldsymbol{R}_u$ and $R_p$ are appropriate residuals of the FE components defined as

$$\boldsymbol{R}_u = -\mathcal{P}_u (\partial_t \boldsymbol{u}_h + (\boldsymbol{u}_h + \tilde{\boldsymbol{u}}) \cdot \nabla \boldsymbol{u}_h - \nu \Delta \boldsymbol{u}_h + \nabla p_h - \boldsymbol{f}),$$
$$R_p = -\mathcal{P}_p (\partial_t p_h + \varepsilon^{-1} \nabla \cdot \boldsymbol{u}_h).$$

$\mathcal{P}_u$ and $\mathcal{P}_p$ are projection operators onto the sub-grid scale spaces of velocities and pressure. In the case in which we take the identity operator in both cases, we recover the original VMS method proposed [22]. Codina [11] proposed to enforce the subscales to be orthogonal to the FE velocity and pressure spaces. It leads to $\mathcal{P}_u \doteq \Pi_u^{\perp} \doteq 1 - \Pi_u$, where $\Pi_u$ is the $L^2$-projector onto $\boldsymbol{V}_h$ (using mass lumping), i.e., for a function $f$, we compute the nodal values of $\Pi_u(f)$ at any node $a$ of the FE mesh as $\Pi_u(f)^a = \frac{\int_\Omega f \varphi^a}{\int_\Omega \varphi^a}$, where $\varphi^a$ is the corresponding nodal shape function. We proceed analogously for $\mathcal{P}_p$ and $L_h$.

Next, we consider the following algebraic approximation of this nonlinear differential system. We refer to [10] for a detailed exposition of the sub-grid problem approximation, e.g., using a motivation based on Fourier transforms. After time discretisation, e.g., using an implicit–explicit first

order time integration, the approximated dynamic sub-grid model for the momentum equation reads as: compute $\tilde{\boldsymbol{u}}^{n+1}$ from

$$\frac{1}{\Delta t}\tilde{\boldsymbol{u}}^{n+1} + \frac{1}{\tau_u^n}\tilde{\boldsymbol{u}}^{n+1} = \frac{1}{\Delta t}\tilde{\boldsymbol{u}}^n + \boldsymbol{R}_u^n,$$

$$\text{where} \quad \tau_u^n = \left(\frac{c_1\nu}{h^2} + \frac{c_2|\boldsymbol{u}_h^n + \tilde{\boldsymbol{u}}^n|}{h}\right)^{-1},$$

or analogously,

$$\tilde{\boldsymbol{u}}^{n+1} = \frac{\tau_t^n}{\Delta t}\tilde{\boldsymbol{u}}^n + \tau_t^n \boldsymbol{R}_u^n, \quad \text{where} \quad \frac{1}{\tau_t^n} = \frac{1}{\Delta t} + \frac{1}{\tau_u}, \qquad (8)$$

$c_1$ and $c_2$ being algorithmic constants, that we set to $c_1 = 4$, $c_2 = 2$ in the numerical experiments. We can treat the sub-grid scale term implicitly because it does not imply an increase in CPU cost. Note that the previous expressions for the sub-grid scales are required at the integration points.

The pressure sub-grid scale component can be treated in a similar way. Again, following the motivation explained in [10], this component can be found from:

$$\frac{1}{\Delta t}\tilde{p}^{n+1} + \varepsilon^{-1}\frac{1}{\tau_p^n}\tilde{p}^{n+1} = \frac{1}{\Delta t}\tilde{p}^n + R_p^n, \quad \text{where} \quad \tau_p^n = \frac{h^2}{\tau_u^n}.$$

This term is not really required to get stability and convergence bounds [2]. Further, the exhaustive numerical analysis in [16] advocates to switch off the pressure sub-grid component for turbulent flow simulations. In this work, we follow this approach.

### 3.3 Time integration and final algorithm

The motivation for using an artificial compressibility method is the possibility to use an explicit time integration scheme of the flow equations. Any choice is open, but in this paper we have used a simple second order time integration scheme for the FE solution and a first order time integration for the velocity sub-grid scales, as described above. This is consistent since the sub-grid scales are multiplied by the stabilisation parameters, which are of the order of the time step size. Note that they essentially behave as the critical time step that would be found in an forward Euler time integration [15]. The solver has been implemented in our in-house code using Fortran90 and parallelised using OpenMP. A 2D matrix-vector implementation is illustrated in Appendix. The final algorithm is depicted below.

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**Algorithm**: A pseudo-compressible VMS method for incompressible flows

- Read $\boldsymbol{u}_h^0$, $p_h^0$ and set $\tilde{\boldsymbol{u}} = \boldsymbol{0}$, loop over time step $n$
- Loop over elements, loop over integration points

  1. Evaluate velocity prediction $\hat{\boldsymbol{u}}_h^{n+1}$ and pressure prediction $\hat{p}_h^{n+1}$ at time step $n+1$ using the second order approximation:

     $$\hat{\boldsymbol{u}}_h^{n+1} = \frac{3}{2}\boldsymbol{u}_h^n - \frac{1}{2}\boldsymbol{u}_h^{n-1}, \quad \hat{p}_h^{n+1} = \frac{3}{2}p_h^n - \frac{1}{2}p_h^{n-1}.$$

  2. Add the nonlinear sub-grid scale component for the convective velocity $\boldsymbol{a}$:

     $$\boldsymbol{a} = \hat{\boldsymbol{u}}_h^{n+1} + \tilde{\boldsymbol{u}}^n.$$

  3. Calculate the stabilisation parameters $\tau_t^n$ as in (8).

  4. Evaluate velocity sub-grid scales:

     $$\tilde{\boldsymbol{u}}^{n+1} = \frac{\tau_t}{\Delta t}\tilde{\boldsymbol{u}}^n - \Pi_u^\perp\left[\tau_t\left(\boldsymbol{a}\cdot\nabla\boldsymbol{u}_h^n + \nabla p_h^n\right)\right]$$

  5. Update the velocity and pressure for the next time step $n+1$ by solving:

     $$(\boldsymbol{u}_h^{n+1}, \boldsymbol{v}_h) = (\boldsymbol{u}_h^n, \boldsymbol{v}_h) - \Delta t\, \boldsymbol{R}_u^n,$$
     $$(q_h, p_h^{n+1}) = (q_h, p_h^n) - \Delta t\, \mathcal{R}_p^n$$

     for all test functions $\boldsymbol{v}_h$ and $q_h$, where

     $$\boldsymbol{R}_u^n = (\boldsymbol{a}\cdot\nabla\hat{\boldsymbol{u}}_h^{n+1}, \boldsymbol{v}_h) + \nu(\nabla\hat{\boldsymbol{u}}_h^{n+1}, \nabla\boldsymbol{v}_h)$$
     $$\quad - (\hat{p}_h^{n+1}, \nabla\cdot\boldsymbol{v}_h)$$
     $$\quad - (\tilde{\boldsymbol{u}}^n, \boldsymbol{a}\cdot\nabla\boldsymbol{v}_h) - <\boldsymbol{f}, \boldsymbol{v}_h>$$
     $$\mathcal{R}_p^n = \varepsilon^{-1}(\nabla\cdot\hat{\boldsymbol{u}}_h^{n+1}, q_h) - \varepsilon^{-1}(\tilde{\boldsymbol{u}}^n, \nabla q_h)$$

- Prescribe boundary conditions

  $$\boldsymbol{u}_h^{n+1} = \bar{\boldsymbol{u}} \quad \text{on } \partial\Omega$$

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## 4 Numerical examples

In this section, a variety of numerical examples will be presented in order to illustrate the performance of the pseudo-compressible VMS solver.
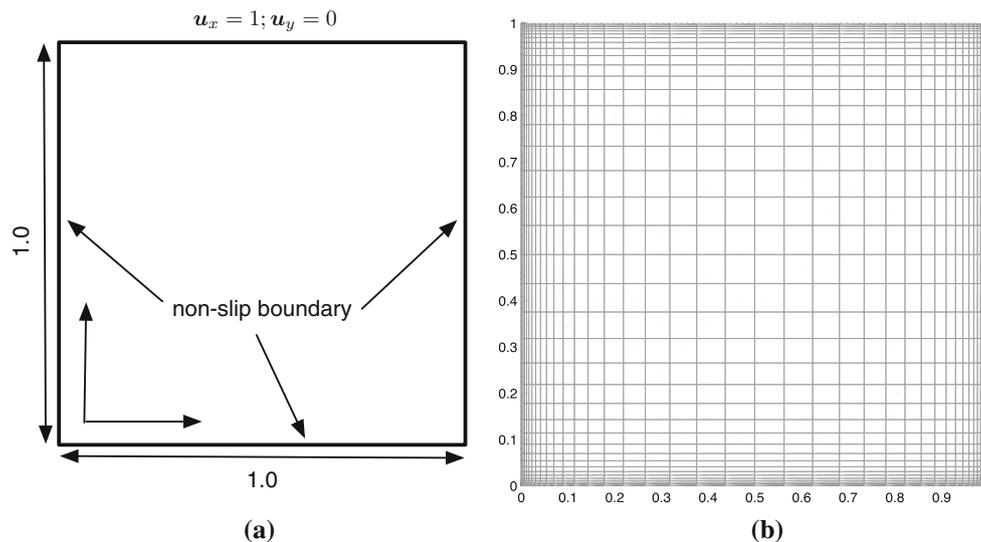
**Fig. 1** Two dimension lid driven flow. **a** Schematic diagram for the lid-driven flow problem. **b** Computational domain: a non-uniform structured meshes

### 4.1 Two-dimensional flow: lid-driven cavity

The first problem considered here is the well documented benchmark case of the lid-driven cavity flow to test the proposed method and the implementation in 2D. A scheme of the test problem configuration is shown in Fig. 1a with the top wall moving to the right at a velocity $u_x = 1$. The bottom and two vertical walls are non-slip boundaries.

A $100 \times 100$ non-uniform structured mesh for $Re = 100$, $Re = 1000$, $Re = 5000$ and $Re = 10,000$, containing finer grid points near the walls, is employed for the computation. Since the target is to compute the steady state, we do not study the effect of ASS in this case. The artificial sound speed is chosen to be $c = 2||\boldsymbol{u}||_\infty^2$. The mesh grid point distance in the $x$ and $y$ directions is set to be a hyperbolic tangent profile. For example, the $y$ coordinates in terms of the $x$ coordinates are given by:

$$y(x) = 0.5 + \frac{\tanh(5(x - 0.5))}{2\tanh(2.5)}, \quad 0 \le x \le 1,$$

as shown in Fig. 1b.

The plots of vertical velocity profiles at the horizontal centreline and of horizontal velocity profiles at the vertical centreline, at different Reynolds numbers, are compared with the second order accurate solution by Ghia et al. [19] in Fig. 2. A good agreement with the reference result is observed.

We also present here the numerical results at $Re = 100,000$ performed on the uniform grids of $512^2$, $1024^2$ and $2048^2$ for the transient case. The artificial sound speed is chosen to be $c = 3||\boldsymbol{u}||_\infty^2$. In Fig. 3, it can be noticed that the coarse grid is more diffusive than the fine grid and the bilinear element under meshsize $1024^2$ is not fine enough to

have a converged result for the lid-driven cavity flow under $Re = 100,000$. But the proposed method can capture the transient features of the cavity flow for very high Reynolds number.

### 4.2 Two-dimensional flow: Kelvin–Helmholtz instability

Kelvin–Helmholtz instability of a plane mixing layer is an important case in fluid dynamics. This example will not lead to what is usually perceived as turbulence, but it shows some features of the flow in common with turbulence in a certain sense, because the flow is extremely sensitive to the initial condition and a broadband energy spectrum of slope between $k^{-4}$ and $k^{-3}$ is developed [20]. The problem is set in the domain $\Omega = [0, 1] \times [0, 1]$. Free slip boundary conditions are applied at $y = 0$ and $y = 1$. At the left and right walls $x = 0$ and $x = 1$, periodic boundaries are imposed, as seen in Fig. 4.

According to [6,20], the initial velocity field is set to be a hyperbolic tangent profile as:

$$u_x = ||\boldsymbol{u}||_\infty \left[ \tanh\left(\frac{2y - 1}{\delta_0}\right) + c\frac{\partial \psi_{\text{per}}}{\partial y} \right]$$

$$u_y = -||\boldsymbol{u}||_\infty c\frac{\partial \psi_{\text{per}}}{\partial x}$$

where $u_x$ and $u_y$ are the velocity components, $\psi_{\text{per}} = ||\boldsymbol{u}||_\infty \exp\left(-\frac{y-0.5}{\delta_0}\right)^2 \cos(\alpha x)$ is a white noise perturbation expressed in a stream function formulation. The initial thickness is chosen as $\delta_0 = 1/28$ and the wave number as $\alpha = 8\pi$ to reach the most unstable wavelength [27]. $||\boldsymbol{u}||_\infty = 1$ and $c_{\text{noise}} = 10^{-3}$ are taken. With the above parameters, the
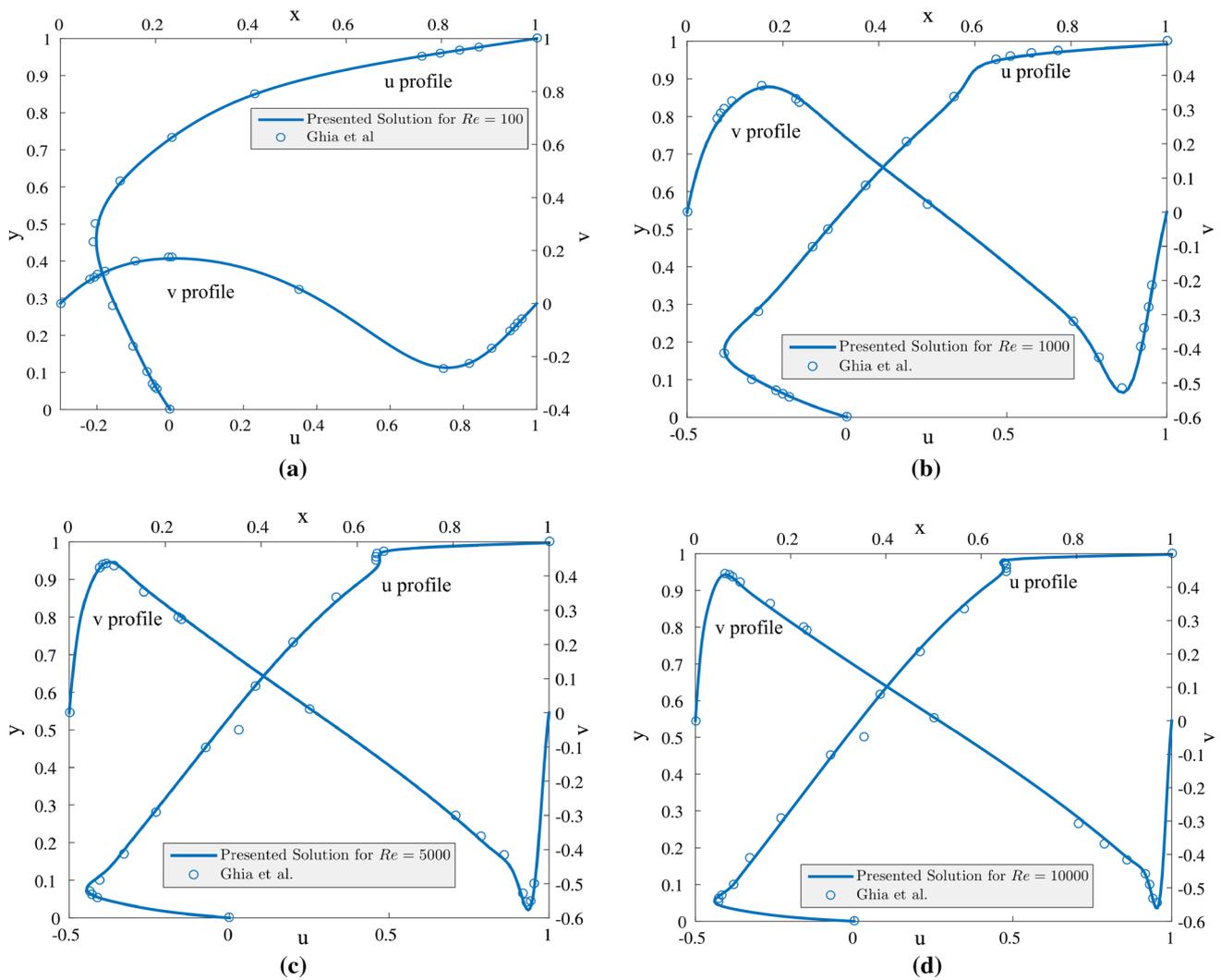
**Fig. 2** 2-D lid-driven flow. Comparison of velocity profiles with Ghia at. [19] at the mid-sections for Reynolds number: **a** $Re = 100$; **b** $Re = 1000$; **c** $Re = 5000$; **d** $Re = 10,000$

Reynolds number of the flow can be evaluated as

$$Re = \frac{||\boldsymbol{u}||_\infty \delta_0}{\nu} = 10,000.$$

The bilinear quadrilateral elements are employed for the simulation in a sequence of meshes $40^2$, $80^2$, $160^2$, $240^2$, $320^2$. The mesh refinement study is carried for the artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$.

In [20] it is reported that the Pressure Stabilised Petrov–Galerkin (PSPG) method fails to converge at the finest grid of $240 \times 240$ elements, whilst in the proposed method the solution is stable for all meshes. Four different stages can be distinguished, which can be seen in Fig. 5. Four Kelvin–Helmholtz vortices will appear along the $x$-axis, that merge into two vortices and then become one (Table 1).

1. *Development of four primary eddies* The four primary eddies can be seen at time unit 16. This development starts similarly for group $80^2$, $160^2$, $240^2$, $320^2$. For the coarsest mesh $40^2$, these stages cannot be observed.

2. *Pairing of four primary eddies* It can be seen that, after the mesh refinement, the pairing starts later. At time unit 43, the different developments of both pairings can be observed.

3. *Pairing of two secondary eddies* The second pairing is finished at about 80 time units, which is close to the results in [25] (75 time units), and earlier in comparison to [20] (115 time units) and [24] (140 time units)

4. *Rotation of the final eddy* At the end of the second pairing, the final eddy rotates at a rather fixed position. The value of the vorticity thickness oscillates during this stage due to the elliptic shape of this vortex.
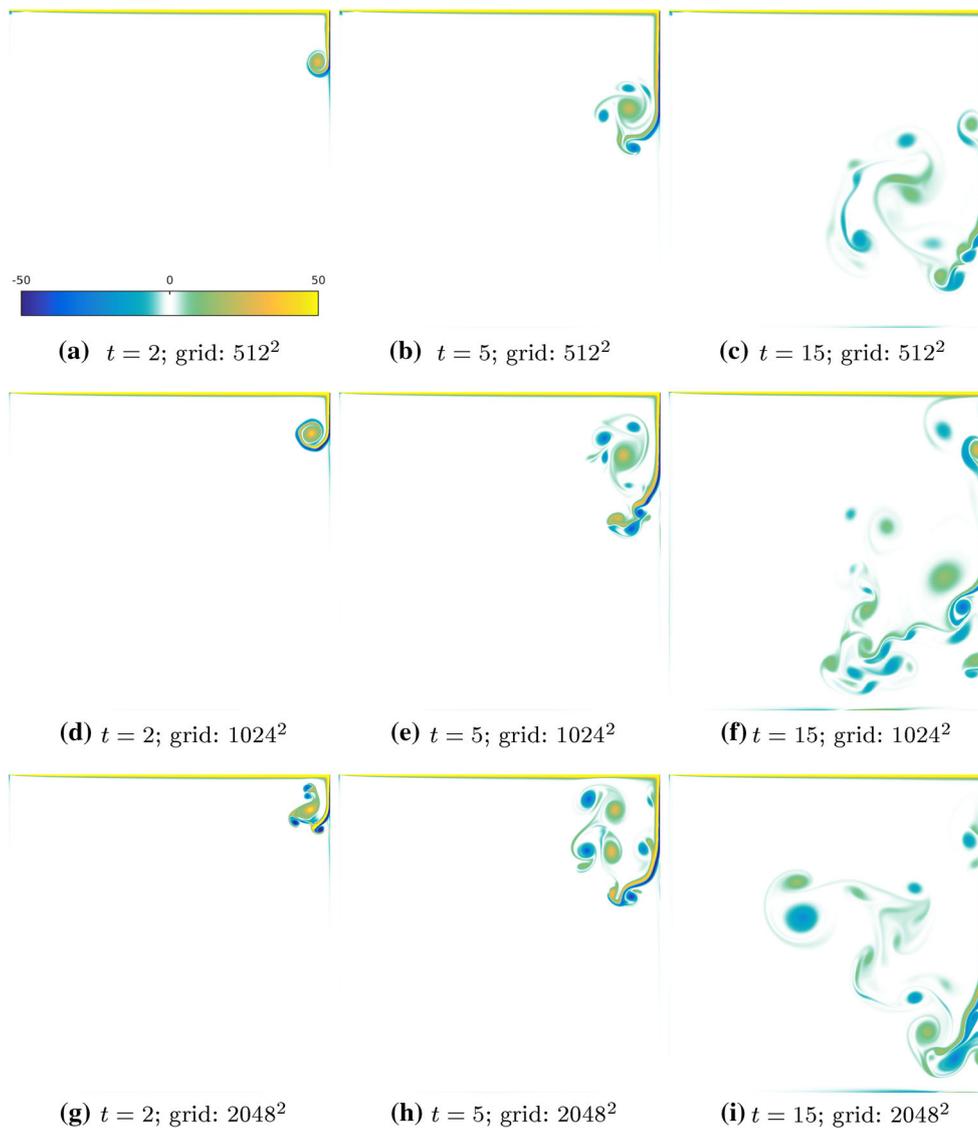
**(a)** $t = 2$; grid: $512^2$     **(b)** $t = 5$; grid: $512^2$     **(c)** $t = 15$; grid: $512^2$

**(d)** $t = 2$; grid: $1024^2$     **(e)** $t = 5$; grid: $1024^2$     **(f)** $t = 15$; grid: $1024^2$

**(g)** $t = 2$; grid: $2048^2$     **(h)** $t = 5$; grid: $2048^2$     **(i)** $t = 15$; grid: $2048^2$

**Fig. 3** 2-D lid-driven flow. Evolution of the vorticity field $-50 < \omega_z < -50$ for Reynolds number $Re = 100000$ at times $t = 2$, $t = 5$, $t = 15$ (from *left* to *right*). The artificial sound speed $c = 3||\boldsymbol{u}||_\infty^2$
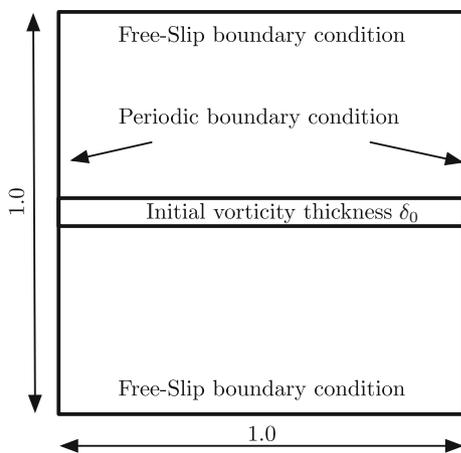


**Fig. 4** Schematic show of the plane mixing layer problem

The integrated kinetic energy is given by

$$E_k = \frac{1}{|\Omega|} \int_\Omega \frac{\boldsymbol{u} \cdot \boldsymbol{u}}{2} \, d\Omega$$

for comparison purposes. The temporal evolution of the kinetic energy dissipation rate is defined as

$$\epsilon = -\frac{dE_k}{dt}.$$

The vorticity thickness is also evaluated quantitatively. The scalar vorticity is defined as

$$\omega(x, y, t) = \frac{1}{2}\left(\frac{\partial u_y}{\partial x} - \frac{\partial u_x}{\partial y}\right).$$
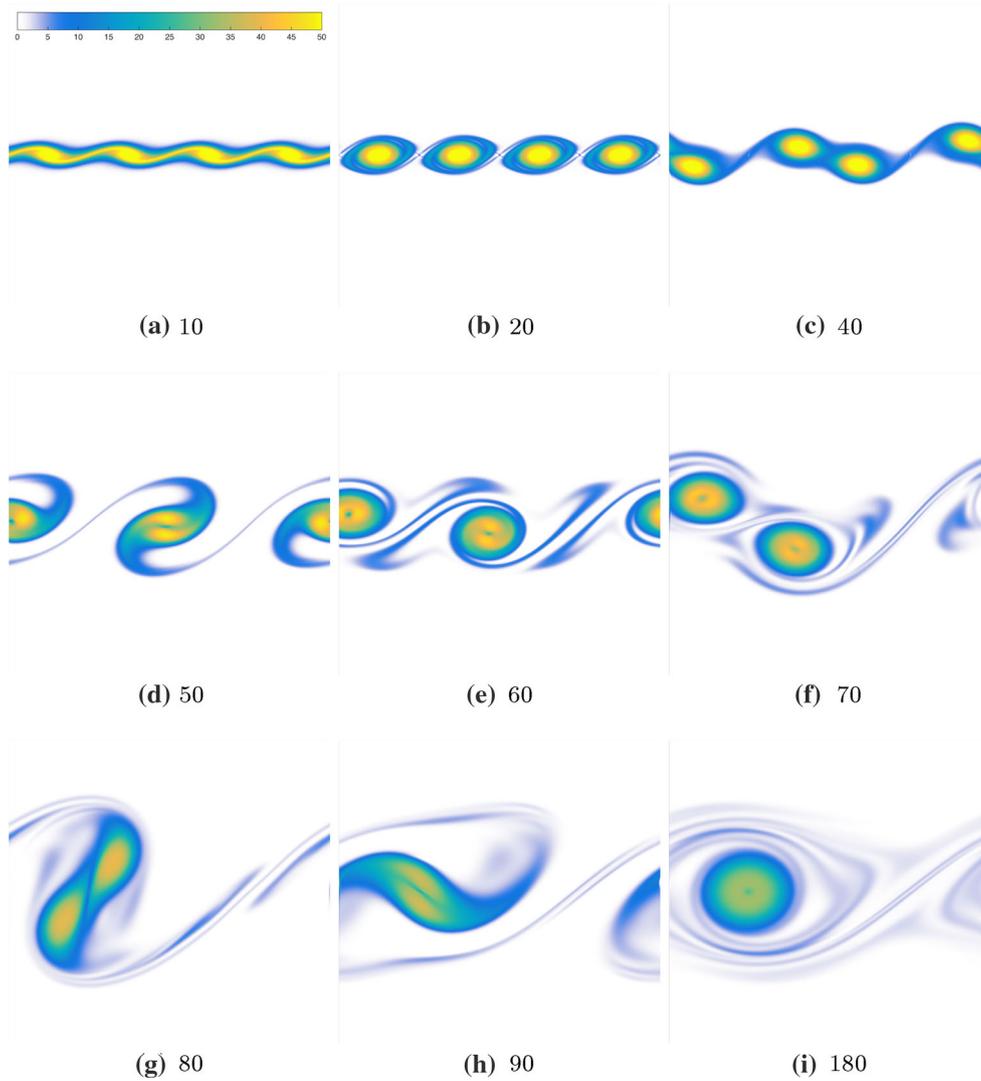
**Fig. 5** 2D plane mixing layer. Coloured vorticity $0 < \omega < 50$ of the velocity field for the finest mesh grid at time unit: 10, 20, 40, 50, 60, 70, 80, 90, 180 in the mesh $320^2$. The artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$

**Table 1** Computational settings for 2D plane mixing layer, running on a Desktop machine

| Elements | Time step $\Delta t$ | ASS | CPU time (hrs) | Peak memory usage (MB) |
|----------|---------------------|-----|----------------|------------------------|
| $40^2$ | $3.2724 \times 10^{-3}$ | $5||\boldsymbol{u}||_\infty$ | 0.0066 | 0.5 |
| $80^2$ | $1.6362 \times 10^{-3}$ | $5||\boldsymbol{u}||_\infty$ | 0.053 | 1.9 |
| $160^2$ | $4.9087 \times 10^{-4}$ | $5||\boldsymbol{u}||_\infty$ | 0.43 | 7.8 |
| $240^2$ | $1.6362 \times 10^{-4}$ | $5||\boldsymbol{u}||_\infty$ | 1.44 | 17.5 |
| $320^2$ | $9.66 \times 10^{-4}$ | $5||\boldsymbol{u}||_\infty$ | 3.45 | 31.1 |
| $320^2$ | $3.40 \times 10^{-3}$ | $2.5||\boldsymbol{u}||_\infty$ | 1.09 | 31.1 |
| $320^2$ | $2.77 \times 10^{-4}$ | $10||\boldsymbol{u}||_\infty$ | 13.4 | 31.1 |
| $320^2$ | $7.97 \times 10^{-5}$ | $20||\boldsymbol{u}||_\infty$ | 46.5 | 31.1 |

Then, define the maximum vorticity $\omega_{max}$ as

$$\omega_{max}(t) = \sup_{y \in [0,1]} \int_0^1 \omega(x, y, t) \, dx,$$

where $\int_0^1 \omega(x, y, t) \, dx$ denotes the integral mean in the periodic $x-$ direction. The vorticity thickness $\delta(t)$ is defined as

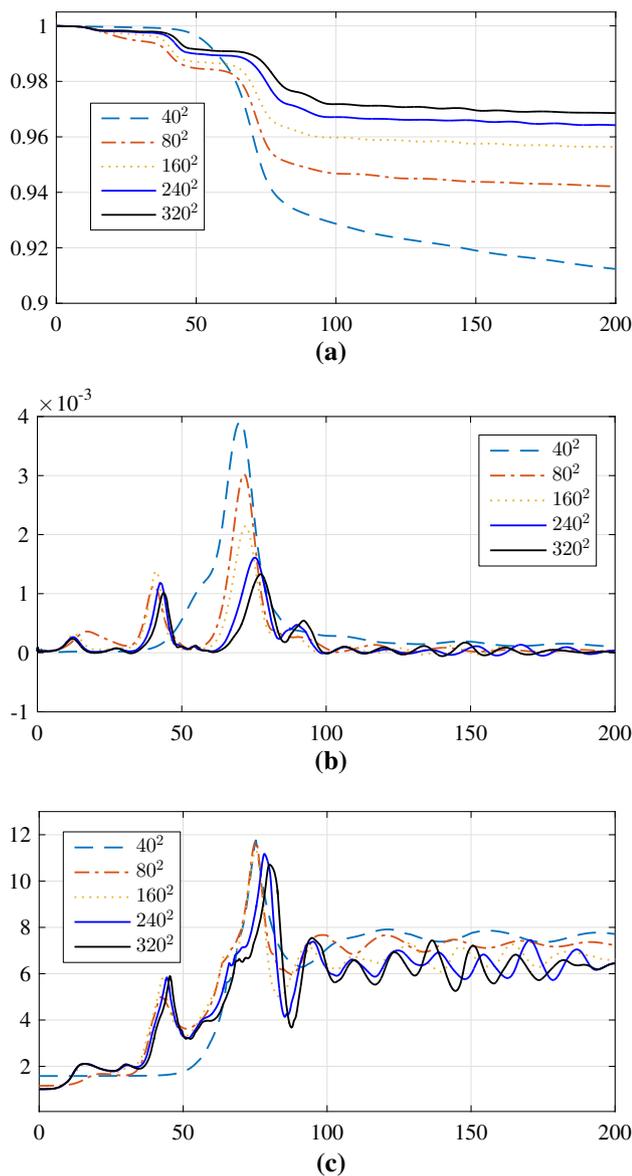$$\delta(t) = \frac{2U_{max}}{\omega_{max}(t)}.$$

**Fig. 6** 2D plane mixing layer. Time history of the kinetic energy, the kinetic energy dissipation rate and the vorticity thickness for a set of Q1Q1 elements $40^2$, $80^2$, $160^2$, $240^2$, $320^2$, the artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$. **a** Evolution of the total kinetic energy. **b** Evolution of the kinetic energy dissipation rate. **c** Evolution of the vorticity thickness

**Fig. 7** 2D plane mixing layer. The effect of different artificial speed $c = 2.5||\boldsymbol{u}||_\infty^2$, $c = 5||\boldsymbol{u}||_\infty^2$, $c = 10||\boldsymbol{u}||_\infty^2$ $c = 20||\boldsymbol{u}||_\infty^2$ for meshsize $320^2$. **a** Evolution of the total kinetic energy. **b** Evolution of the kinetic energy dissipation rate. **c** Evolution of the vorticity thickness

### 4.2.1 Mesh refinement

Figure 6 shows the time history of the kinetic energy, the kinetic energy dissipation rate and the vorticity thickness under mesh refinement. The eddies pairing causes the increases of the vorticity thickness, as well as the energy dissipation. For the coarsest mesh size $40 \times 40$, the simulation cannot capture the four and two eddies periods and it can only capture the final large eddy. The mesh of $80 \times 80$ elements is sufficient to clearly observe the development of four stages. The pairing happens relatively slower after refining the mesh.

### 4.2.2 Effect of artificial sound speed

The effect of the different artificial sound speed under mesh-size $320 \times 320$ is shown in Fig. 7. It can be seen that ASS = 5 is enough to capture the key features of the problem. Using the lower ASS number, higher kinetic energy dissipation rate and oscillation can be observed.

### 4.3 Three-dimensional laminar flow: lid-driven cavity

The 3D lid-driven cavity flow illustrated in Fig. 8 is studied to examine the performance of the proposed method
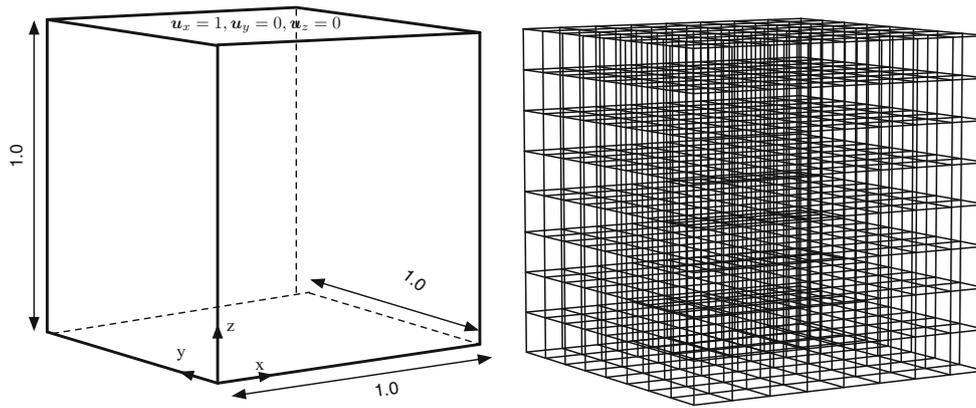
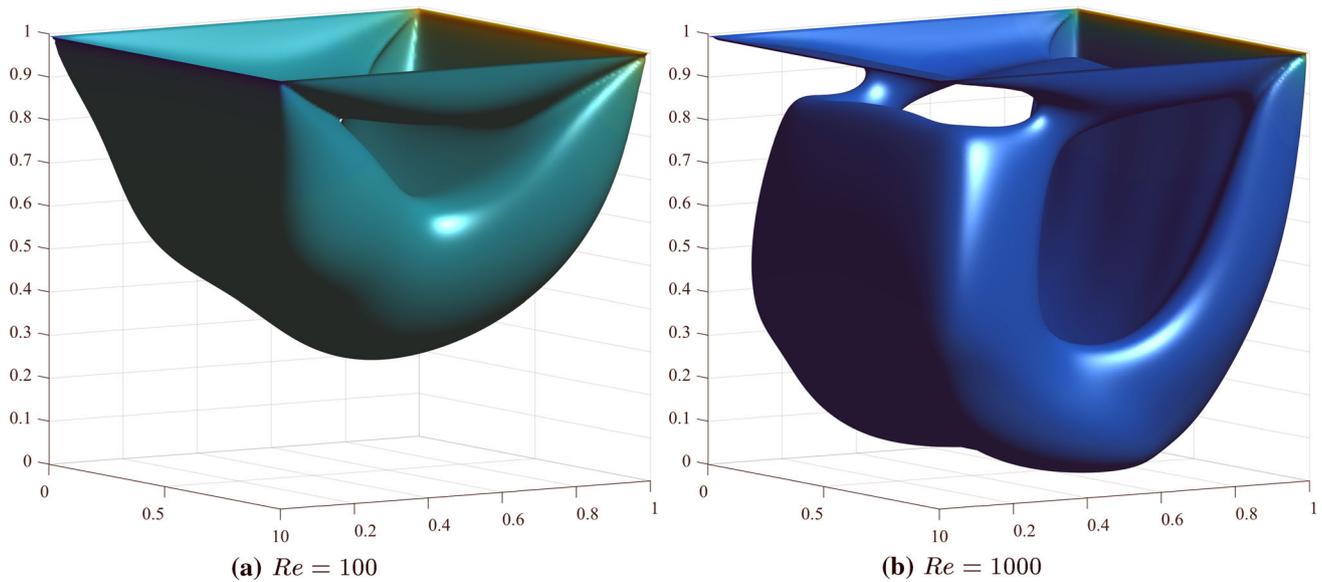**Fig. 8** Schematic show of the 3D lid-driven cavity flow



**(a)** $Re = 100$ **(b)** $Re = 1000$

**Fig. 9** 3D lid-driven cavity. Isosurfaces of magnitude of velocity $|\boldsymbol{u}| = 0.15$, with $100^3$ trilinear Q1Q1 elements for Reynolds numbers $Re = 100$ and $Re = 1000$
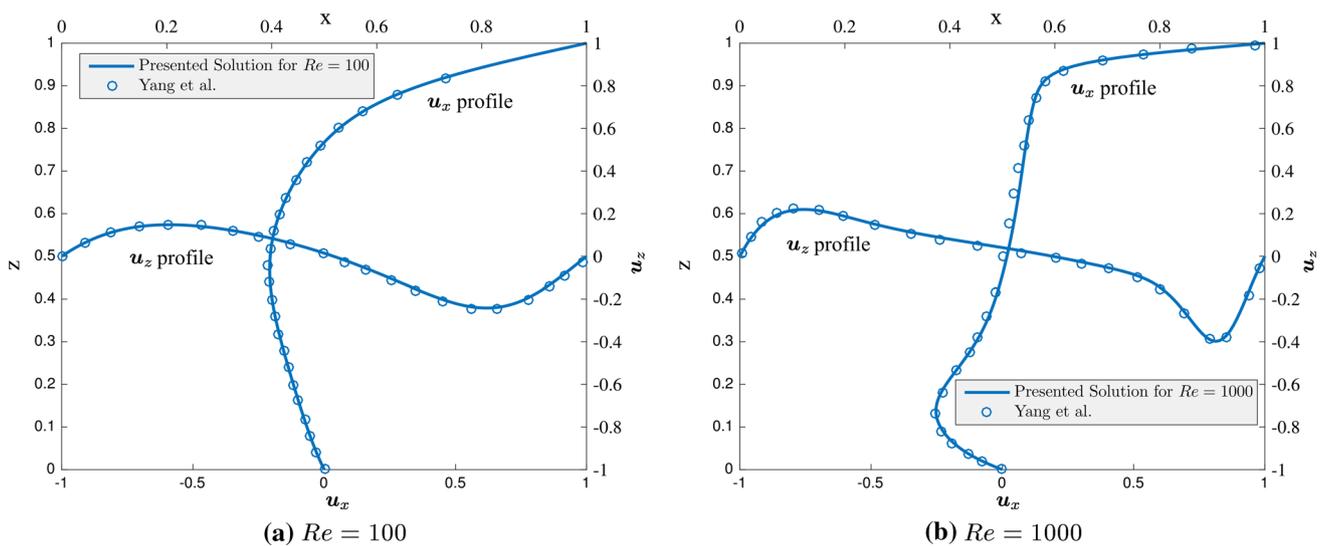


**(a)** $Re = 100$ **(b)** $Re = 1000$

**Fig. 10** 3D lid-driven cavity. Comparison of velocity profiles on the mid-plane $y = 0.5$ with reference data [32] for Reynolds number of 100 and 1000. The artificial sound speed $c = 2||\boldsymbol{u}||_\infty^2$
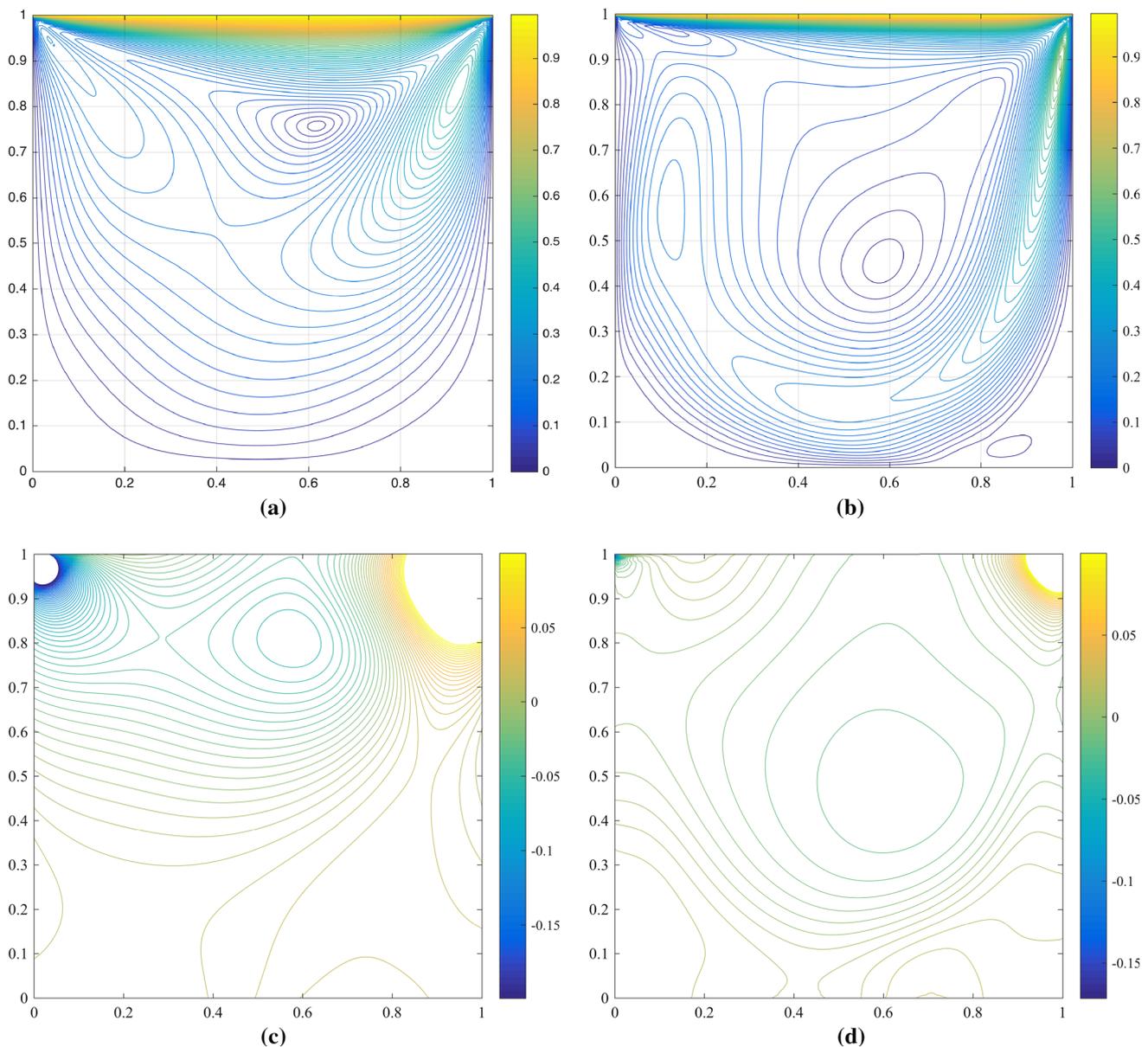
**Fig. 11** 3D lid-driven cavity. Isolines of velocity magnitude and pressure field on the mid-plane $y = 0.5$ of cubic cavity at $Re = 100$ and $Re = 1000$. The artificial sound speed $c = 2||\boldsymbol{u}||_\infty^2$. **a** Velocity magnitude, $Re = 100$. **b** Velocity magnitude, $Re = 1000$. **c** Pressure field, $Re = 100$. **d** Pressure field, $Re = 1000$

and the implementation. The domain in physical space is $\Omega = [0, 1] \times [0, 1] \times [0, 1]$. The top wall moves at a velocity $u_x = 1, u_y = 0, u_z = 0$. The bottom and four vertical walls are non-slip boundaries. A $100^3$ uniform structured mesh for $Re = 100$ and $Re = 1000$ is employed for the computation of the laminar flow. Figure 9 shows the isosurfaces of magnitude of the velocity $|\boldsymbol{u}| = 0.15$ when the flow is reached to the steady state. The artificial sound speed is chosen to be $c = 2||\boldsymbol{u}||_\infty^2$, as in the 2D. Quantitive velocity profiles were compared with the reference by Yang [32].

*4.3.1 One-dimensional profile*

The velocity profiles on the mid-plane $y = 0.5$ centrelines is illustrated in Fig. 10 for the quantitive comparisons. A good agreement has been reached with the results obtained by Yang [32].

*4.3.2 Two-dimensional profile*

The velocity magnitude in the mid-plane $y = 0.5$ and its corresponding pressure field is shown in Fig. 11. The pressure

**Fig. 12** 3D lid-driven cavity. Streamlines on the mid-plane $x = 0.5$, $y = 0.5$ and $z = 0.5$ of the cubic cavity at Reynolds number of 100 and 1000. The artificial sound speed $c = 2||\boldsymbol{u}||_\infty^2$. **a** $Re = 100$ on the mid-plane $x = 0.5$. **b** $Re = 1000$ on the mid-plane $x = 0.5$. **c** $Re = 100$ on the mid-plane $y = 0.5$. **d** $Re = 1000$ on the mid-plane $y = 0.5$. **e** $Re = 100$ on the mid-plane $z = 0.5$. **f** $Re = 1000$ on the mid-plane $z = 0.5$



(a)

(b)

(c)

(d)

(e)

(f)
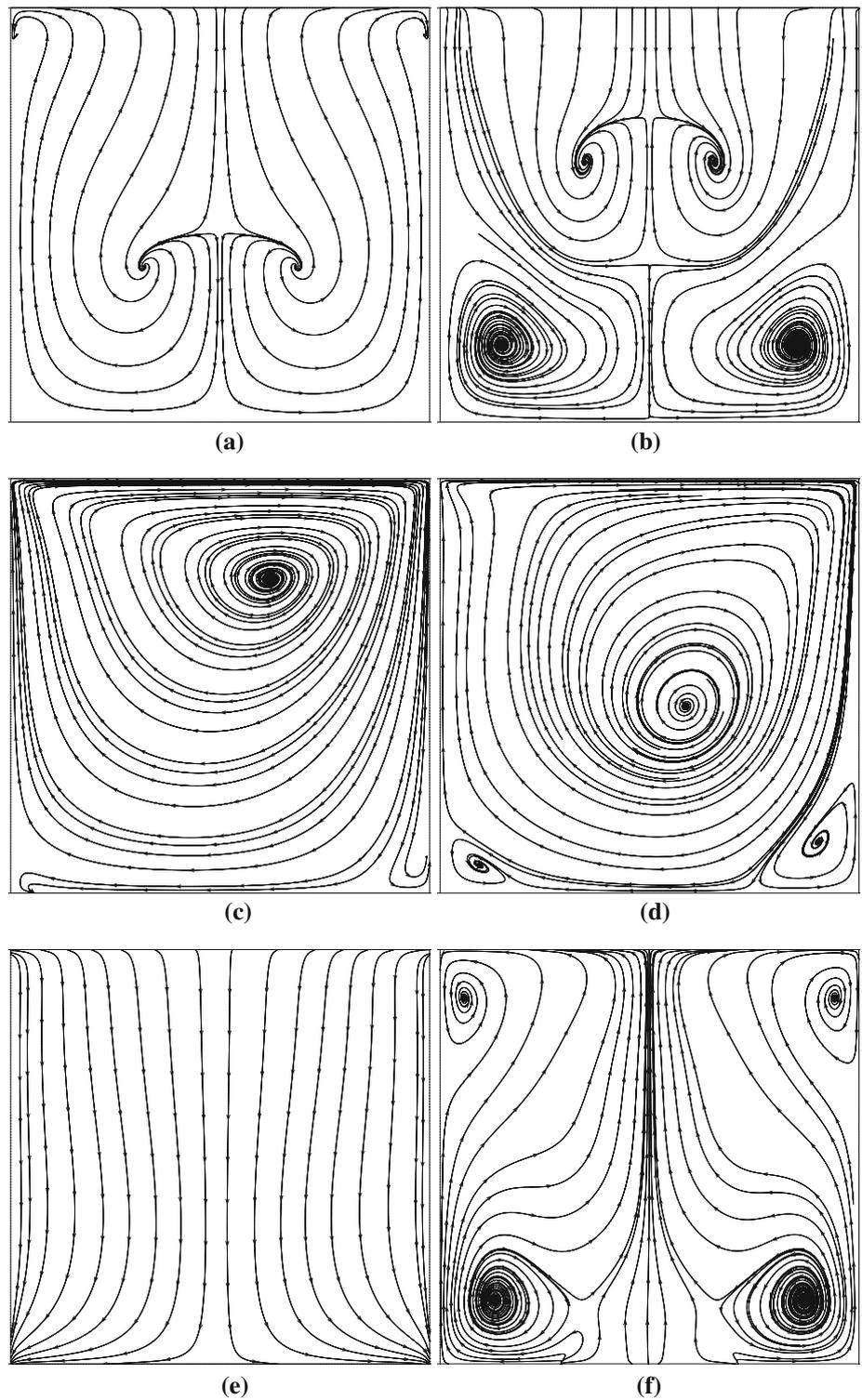
field is stable using the equal velocity and pressure interpolation, which qualitatively agrees well with the results of [18].

In order to show the flow patterns of the cubic cavity flow, streamlines on the plane of $x = 0.5$, $y = 0.5$ and $z = 0.5$ are presented in Fig. 12. The effects of Reynolds numbers on the flow patterns can be observed. As the Reynolds number increases, the strength of the secondary vortices is enhanced. A similar flow patten was reported in [18].

**Table 2** Computational settings for 3D Taylor Green vortex, run on a 12 cores machines, speed up by OpenMP

| Elements | Time step $\Delta t$ | ASS | CPU time (h) | Peak memory usage (GB) |
|---|---|---|---|---|
| $128^3$ | $1.63 \times 10^{-3}$ | $3\|\boldsymbol{u}\|_\infty$ | 1 | 0.68 |
| $128^3$ | $4.90 \times 10^{-4}$ | $5\|\boldsymbol{u}\|_\infty$ | 3 | 0.68 |
| $128^3$ | $1.64 \times 10^{-4}$ | $10\|\boldsymbol{u}\|_\infty$ | 11 | 0.68 |
| $128^3$ | $3.08 \times 10^{-5}$ | $20\|\boldsymbol{u}\|_\infty$ | 55 | 0.68 |
| $256^3$ | $2.46 \times 10^{-4}$ | $5\|\boldsymbol{u}\|_\infty$ | 16 | 5.44 |
| $512^3$ | $1.23 \times 10^{-4}$ | $5\|\boldsymbol{u}\|_\infty$ | 253 | 43.51 |

### 4.4 Three-dimensional turbulent flow: Taylor–Green vortex

The Taylor–Green vortex problem is widely used in the numerical simulation of turbulent flow. This flow transits to turbulence and represents a basic turbulence decay mode similar to decaying homogeneous turbulence [7,8]. The domain in physical space is $\Omega = [0, 2\pi] \times [0, 2\pi] \times [0, 2\pi]$ with periodic boundary conditions in all directions. Within the domain, initial velocity and pressure is defined as

$$u_x = u_0 \cos(x) \sin(y) \sin(z)$$
$$u_y = -u_0 \sin(x) \cos(y) \sin(z)$$
$$u_z = 0$$
$$p = p_0 + \frac{1}{16}(\sin(\cos(2x) + \cos(2y))(\cos(2z) + 2)$$

where $u_0 = 1$. The initial velocity field on the Fourier space has eight modes located at the wave number $\boldsymbol{k} = (\pm 1, \pm 1, \pm 1)$. The Reynolds number of flow is defined as $Re = \frac{u_0 L}{\nu}$ and is equal to 1600.

The primary method for evaluating the TGV solution is examining the energy dissipation rate. The integrated kinetic energy is given by

$$E_k = \frac{1}{|\Omega|} \int_\Omega \frac{\boldsymbol{u} \cdot \boldsymbol{u}}{2} \, d\Omega.$$

The temporal evolution of the kinetic energy dissipation rate is defined as

$$\epsilon_1 = -\frac{dE_k}{dt}.$$

A second kinetic energy dissipation rate can be defined from the integrated enstrophy of the problem:

$$\epsilon_2 = \frac{2\nu}{|\Omega|} \int_\Omega \frac{\boldsymbol{\omega} \cdot \boldsymbol{\omega}}{2} \, d\Omega,$$

where $\boldsymbol{\omega}$ is the vorticity of the velocity field.

Trilinear elements are employed for the simulation in a sequence of meshes $100 \cdot 2^i \times 128 \cdot 2^i$ for $i = 1, 2, 3$. All the simulation were run on a single processor desktop, including the case for $512^3$. The details of each cases setting were given in Table 2. We also report the peak memory consumption for every run. In Fig. 13 we present some vorticity iso-surface images. The initial condition Fig. 13a has eight vortices with the same scale corresponding to the eight Fourier modes. The vorticity iso-surfaces are very similar to the results obtained by the implicit VMS [16].

#### 4.4.1 Effect of artificial sound speed

The effect of the artificial sound speed (ASS) was tested under a mesh size of $100^3$. The ASS is chosen to be 3, 5, 10, 20 times the maximum of the velocity $\|\boldsymbol{u}\|_\infty$. Figure 14 shows the effect of ASS on the resolved energy, the total kinetic energy evolution and dissipation rate. It can be observed from that the resolved dissipation is very similar for all the cases. However, the energy oscillation can be seen from time $t = 0$ to $t = 3$ due to the artificial compressibility scheme. As the ASS number increases, the oscillation decreases, as the relaxation of the incompressibility becomes strict. As expected, the large ASS number brings the time stepping restrictions and causes more computational time. Table 2 also illustrates the computation times required for the different ASS values. For the Taylor–Green vortex problem, the interest is to study the transition period from the laminar to turbulent flow and the turbulence decay. The energy oscillation caused by the artificial compressibility scheme is very small before the turbulent transition. As a result, ASS equal to $5\|\boldsymbol{u}\|_\infty$ is chosen for the mesh refinement simulations.

#### 4.4.2 Mesh refinement

Figure 15 shows the total kinetic energy and the total energy dissipation rate. The directly computing energy dissipation rate is well predicated compared with the DNS result by [7,17]. This indicates that the turbulent structures are under-resolved and the sub-grid scales plays a significant role in the energy dissipation. And the proposed explicit variational
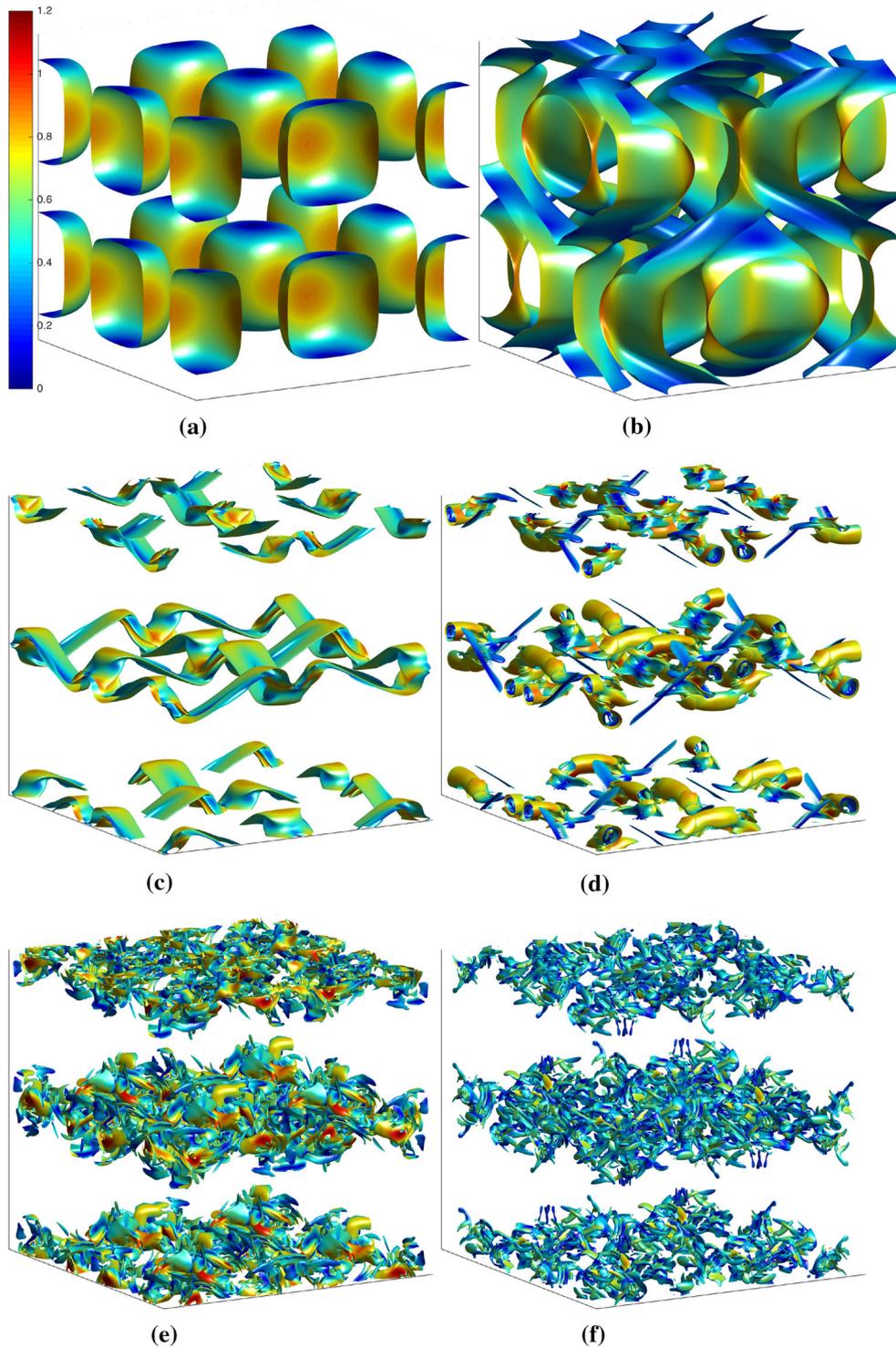
**Fig. 13** Taylor-Green vortex. Vorticity isosurfaces with velocity colored at different time step, mesh of $512^3$ trilinear elements, the artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$. **a** Isosurface for $|\boldsymbol{\omega}| = 1$ at $t = 0.0$. **b** Isosurface for $|\boldsymbol{\omega}| = 1$ at $t = 2.0$. **c** Isosurface for $|\boldsymbol{\omega}| = 4$ at $t = 4.0$. **d** Isosurface for $|\boldsymbol{\omega}| = 7$ at $t = 6.0$. **e** Isosurface for $|\boldsymbol{\omega}| = 8$ at $t = 9.0$. **f** Isosurface for $|\boldsymbol{\omega}| = 9$ at $t = 11.0$
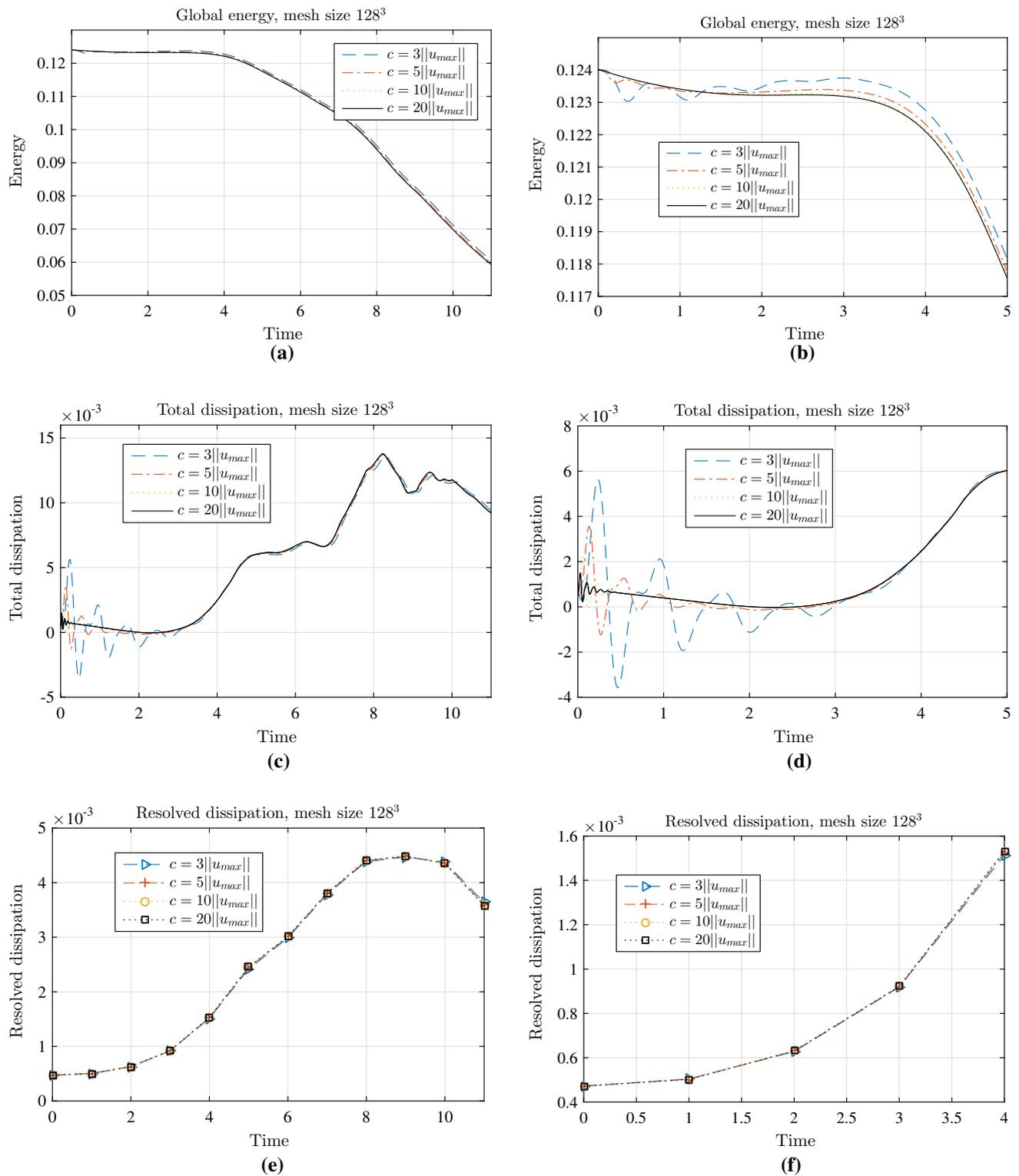
**Fig. 14** Taylor–Green vortex. Study the effect of different artificial sound speed for resolved energy, total kinetic energy evolution and dissipation rate under grid $128^3$. **a** Total kinetic energy. **b** Close view. **c** Total kinetic dissipation rate. **d** Close view. **e** Resolved dissipation rate. **f** Close view
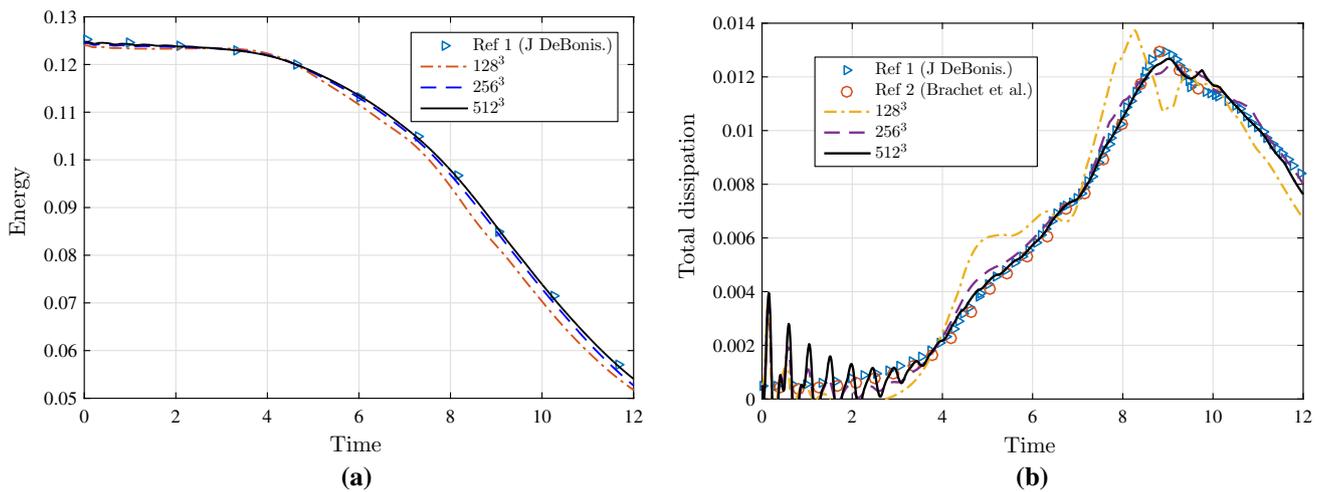
**Fig. 15** Taylor–Green vortex. The evolution of the total kinetic energy and the kinetic dissipation rate under mesh refinement, the artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$. **a** Total kinetic energy. **b** Total kinetic dissipation rate
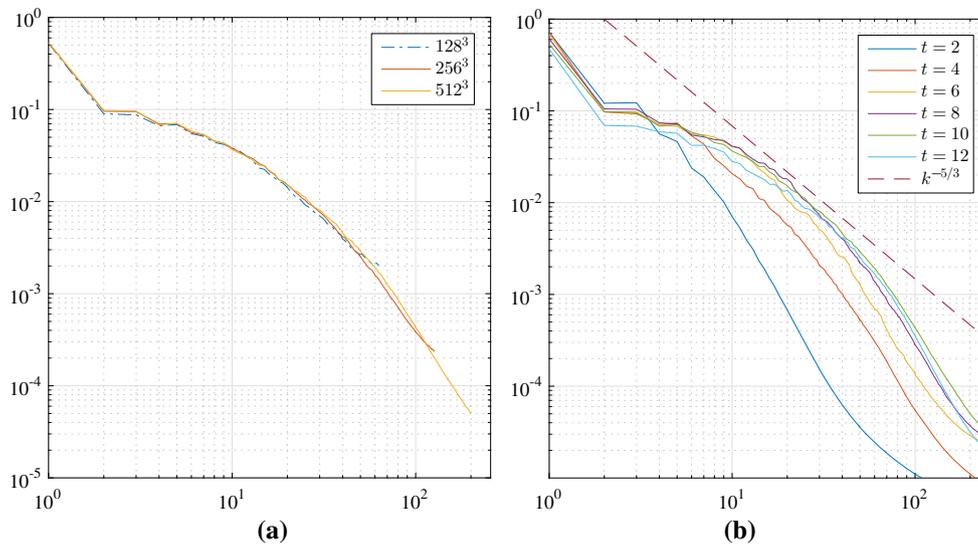


**Fig. 16** Taylor–Green vortex. Energy spectrum for the Taylor–Green Vortex problem, the artificial sound speed $c = 5||\boldsymbol{u}||_\infty^2$. **a** Energy spectrum under different mesh size at $t = 10$. **b** Evolution of the energy spectra under different time on the $512^3$ grid

multiscale method is able to approximate the un-resolved scales and simulate the turbulent flow.

### 4.4.3 Energy spectra

The progression of the energy spectra with time on $512^3$ is shown in Fig. 16a. At $t = 2$ all the energy is confined to small wave numbers. As the time progresses, the energy begins to cascade down to smaller scales. The energy in the smallest scales peaks around $t = 10$. This corresponds to the maximum energy dissipation rate. The spectra for $t = 10$ for a set of meshes are shown in Fig. 16b. The finer grid is able to capture smaller energy scales.

## 5 Conclusions

In this work, we have proposed a pseudo-compressible VMS solver for turbulent incompressible flows. As a result, VMS-type LES models can now be combined with explicit time stepping techniques. We have analyzed the resulting scheme for a set of benchmark tests. Out of these results, we can conclude that one can obtain accurate under-resolved LES simulations of turbulent flows by using the proposed explicit VMS scheme with a very limited memory consumption. In particular, the transient numerical examples, $2D$ Kelvin-Helmholtz instability and $3D$ Taylor-Green Vortex, show a good convergence with respect to the artificial compressibility sound speed and the mesh size.

## Appendix

Here we illustrate the matrix-vector implementation in 2D for linear FEs for simplicity. The nodal shape function associated to node $a$ is denoted by $N^a(x)$ and its value at gauss point $b$ is represent as $N^a_b$. Every node $a$ is represented by the lexicographical index $(i_a, j_a)$, with $i_a, j_a = 0, 1$. Further, given the $\alpha$-th axis direction, with $\alpha = 1, 2$, we define the function $tw(a, \alpha) = tw(i_a, j_a, \alpha)$ as $(1 - i_a, j_a)$ for $\alpha = 1$ and $(i_a, 1 - j_a)$ for $\alpha = 2$. We use a nodal quadrature for the numerical integration, which leads to a lumped mass matrix. As a result, using the expression of the nodal shape functions and their derivatives on nodes, we get:

$$
\begin{aligned}
N^a(x_b) &= \delta_{ab}, \quad N^a(x_b)\delta_{ab} \quad \partial_\alpha N^a(x_b) \\
&= -\partial_\alpha N^{tw(a,\alpha)}(x_b), \partial_\alpha N^a(x_{tw(a,\beta)}) \\
&= \delta_{\alpha\beta}\partial_\alpha N^a(x_a),
\end{aligned}
$$

where $x_b$ denotes the coordinates of the node $b$ and $\delta_{\alpha\beta}$ the Kronecker delta. Using these expressions, we can implement efficiently all the terms in our matrix-free formulation using the following expressions. The viscous term is implemented as follows:

$$
\begin{aligned}
\int_K \partial_\beta u^\alpha(x)\partial_\beta N^a(x)\mathrm{d}\Omega &= \sum_{b=1}^4 \partial_\beta u^\alpha(x_b)\partial_\beta N^a_b\, |J|\, w_b \\
&= (\partial_\beta u^\alpha(x_a) + \partial_\beta u^\alpha(x_{tw(a,\beta)}))\partial_\beta N^a_a\, |J|\, w_a.
\end{aligned}
$$

The nonlinear convective term is implemented as:

$$
\begin{aligned}
\int_K a_\beta \partial_\beta u^\alpha(x) a_\gamma \partial_\gamma N^a(x)\mathrm{d}\Omega \\
= \sum_{b=1}^4 (a_\beta \partial_\beta u^\alpha)(x_b) a_\gamma(x_b)\partial_\gamma N^a_b\, |J|\, w_b \\
= ((a_\beta \partial_\beta u^\alpha a_\gamma)(x_a) \\
+ (a_\beta \partial_\beta u^\alpha a_\gamma)(x_{tw(a,\gamma)}))\partial_\gamma N^a_a\, |J|\, w_a.
\end{aligned}
$$

We proceed analogously for the rest of terms.

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