

A high order discontinuous Galerkin method for fluid-structure interaction

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1. Introduction

Many important scientific and engineering problems require analysis of fluid-structure interaction (FSI). For example, aeroelastic flutter can produce large and potentially destructive vibrations in aircraft [7], turbines [1], and other structures [3] or biological applications such as study of fluid flow inside human vocal tract [5]. The presented study deals with a high order discontinuous Galerkin method for fluid-structure interaction.

2. Mathematical model

The model consists of the Navier-Stokes equations governing the motion of a compressible fluid flow coupled to a rigid body dynamics, i.e., a movement of a structure, described by a second order ordinary differential equation. Arbitrary Lagrangian Eulerian (ALE) method is used to treat the deformable domain. The viscous gas dynamics in computational domain $\Omega_t \subset \mathbb{R}^2$ for any $t \in (0, T), T > 0$ is described by the Navier-Stokes equations, see e.g. [2]. The Navier-Stokes equations written in the conservative form reads

$$\begin{aligned} \frac{\partial}{\partial t}(\varrho) + \sum_{i=1}^2 \frac{\partial}{\partial x_i}(\varrho v_i) &= 0, \\ \frac{\partial}{\partial t}(\varrho v_i) + \sum_{j=1}^2 \frac{\partial}{\partial x_j}(\varrho v_i v_j + p \delta_{ij}) &= \sum_{j=1}^2 \frac{\partial}{\partial x_j}(\tau_{ij}), \quad \text{for } i = 1, 2, \\ \frac{\partial}{\partial t}(\varrho E) + \sum_{j=1}^2 \frac{\partial}{\partial x_j}(\varrho v_j E + v_j p) &= \sum_{j=1}^2 \frac{\partial}{\partial x_j}(-q_j + v_i \tau_{ij}), \end{aligned} \quad (1)$$

where ϱ is the fluid density, p is the pressure, v_1, v_2 are the velocity components of the velocity vector \mathbf{v} , and E is the total energy. The components of the viscous stress tensor $\boldsymbol{\tau}$ and the heat flux \mathbf{q} are given by

$$\tau_{ij} = \mu \left(\frac{\partial v_i}{\partial x_j} + \frac{\partial v_j}{\partial x_i} - \sum_{k=1}^2 \frac{2}{3} \frac{\partial v_k}{\partial x_k} \delta_{ij} \right) \quad (2)$$

and

$$q_i = \frac{\mu}{\text{Pr}} \frac{\partial}{\partial x_i} \left(E + \frac{p}{\varrho} - \sum_{j=1}^2 \frac{1}{2} v_j v_j \right), \quad (3)$$

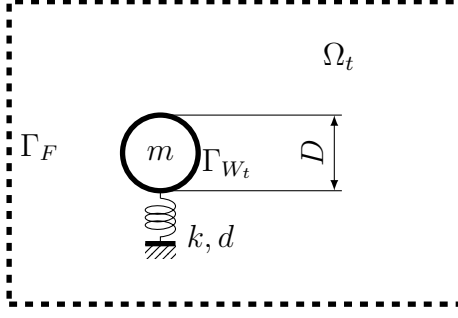


Fig. 1. Sketch of the computation domain Ω_t

where μ is dynamic viscosity and Pr , assumed to be constant $\text{Pr} = 0.72$, is the Prandtl number. For an ideal gas, the pressure p has the form

$$p = (\gamma - 1)\varrho \left(E - \frac{1}{2} \sum_{i=1}^2 v_i v_i \right), \quad (4)$$

where γ is the adiabatic gas constant, γ is set to 1.4 in presented study. Imposed boundary conditions are either free-stream at the far field, or adiabatic no-slip conditions at the boundaries of the structure, i.e., $\mathbf{v}|_{\Gamma_{W_t}}$ is equal to the velocity of the structure. System (1) is supplemented with suitable initial conditions.

The motion of the structure in the one-direction is modeled by the second-order differential linear equation, i.e.,

$$m\ddot{h} + d\dot{h} + kh = L, \quad (5)$$

where m is the oscillating mass of the system, d and k denote the mechanical damping and stiffness of the oscillator unit, respectively, h is the displacement of the oscillator and L is the force exerted by the fluid on the structure in the transverse direction, see Fig 1. Fluid flow model (1) is coupled with the rigid body model (5) via L in the following manner, i.e.,

$$L = -l \int_{\Gamma_{W_t}} \sum_{j=1}^2 (\tau_{2j} - p\delta_{2j}) n_j dS, \quad (6)$$

where l is the depth of the structure, $\mathbf{n} = (n_1, n_2)$ is the unit outer normal to $\partial\Omega_t$ on Γ_{W_t} (pointing into the structure). System (5) is supplemented with initial suitable initial conditions.

3. Numerical approximation

The fluid flow model is discretized using a high-order discontinuous Galerkin formulation with triangular grid elements and nodal basis functions and the domain movement is taken into account with aid of arbitrary Lagrangian-Eulerian method, see e.g. [7]. Following standard procedure for DG discretization of second-derivatives, first the auxiliary gradient variable \mathbf{g} is introduced, and then governing equations are rewritten as the system of first order equations, i.e.,

$$\frac{\partial}{\partial t}(\mathbf{u}) + \nabla \cdot \mathbf{F}^i(\mathbf{u}) - \nabla \cdot \mathbf{F}^v(\mathbf{u}, \mathbf{g}) = \mathbf{0}, \quad (7)$$

$$\nabla \mathbf{u} = \mathbf{g}, \quad (8)$$

where $\mathbf{u} = [\varrho, \varrho v_1, \varrho v_2, \varrho E]^T$ is the solution vector, \mathbf{F}^i is inviscid flux given as

$$\mathbf{F}^i(\mathbf{u}) = \begin{bmatrix} \varrho v_1 & \varrho v_2 \\ \varrho v_1 v_1 & \varrho v_2 v_1 \\ \varrho v_1 v_2 & \varrho v_2 v_2 \\ (\varrho E + p)v_1 & (\varrho E + p)v_2 \end{bmatrix} \quad (9)$$

and \mathbf{F}^v is viscous flux given as

$$\mathbf{F}^v(\mathbf{u}, \mathbf{g}) = \begin{bmatrix} 0 & 0 \\ \tau_{11}(\mathbf{g}) & +\tau_{12}(\mathbf{g}) \\ \tau_{21}(\mathbf{g}) & +\tau_{22}(\mathbf{g}) \\ -q_1(\mathbf{u}) + v_1\tau_{11}(\mathbf{g}) + v_2\tau_{21}(\mathbf{g}) & -q_2(\mathbf{u}) + v_1\tau_{12}(\mathbf{g}) + v_2\tau_{22}(\mathbf{g}) \end{bmatrix}. \quad (10)$$

The inviscid fluxes are computed using Roe's method [6], and the numerical fluxes for the viscous terms are chosen according to the compact discontinuous Galerkin (CDG) method [4]. The computational domain Ω is discretized by the computational mesh with elements $\mathcal{T}_h = \{K\}$. The solution (\mathbf{u}, \mathbf{g}) is sought in $[V_h^p]^4$ and $[V_h^p]^{4 \times 2}$, respectively, where $V_h^p = \{v \in L^2(\Omega), v|_K \in P^p(K), \forall K \in \mathcal{T}_h\}$ with P^p being the space of polynomial functions of degree at most $p \geq 1$ on K . The semi-discrete DG formulation is expressed as: find $\mathbf{u}_h \in [V_h^p]^4$ and $\mathbf{g}_h \in [V_h^p]^{4 \times 2}$ such that for all $K \in \mathcal{T}_h$

$$\begin{aligned} \int_K \frac{\partial \mathbf{u}_h}{\partial t} \cdot \boldsymbol{\varphi} + \int_K (\mathbf{F}^i(\mathbf{u}_h) - \mathbf{F}^v(\mathbf{u}_h, \mathbf{g}_h)) : \nabla \boldsymbol{\varphi} - \\ - \int_{\partial K} (\mathbf{F}^i(\mathbf{u}_h) - \mathbf{F}^v(\mathbf{u}_h, \mathbf{g}_h)) \cdot \boldsymbol{\varphi} = \mathbf{0}, \quad \forall \boldsymbol{\varphi} \in [P^p(K)]^4, \end{aligned} \quad (11)$$

$$\int_K \mathbf{g}_h : \boldsymbol{\psi} + \int_K \mathbf{u}_h \cdot (\nabla \cdot \boldsymbol{\psi}) - \int_{\partial K} (\mathbf{u}_h \otimes \mathbf{n}) : \boldsymbol{\psi} = \mathbf{0}, \quad \forall \boldsymbol{\psi} \in [P^p(K)]^{4 \times 2}. \quad (12)$$

Fluxes in Eqs. (11) and (12) are modified according to the ALE method, see e.g. [7]. All integrals in Eqs. (11) and (12) are integrated using high-order Gaussian quadrature rules. Time integration is done with aid of a high-order Runge-Kutta (RK) method.

4. Numerical results

To validate the high-order scheme, we considered a test problem consisting of flow-induced vibration of a circular cylinder, where the cylinder is allowed to move in vertical direction, see Fig. 1. The far field fluid has velocity $\mathbf{v} = (1, 0)$ m/s, density $\varrho = 1$ kg/m³, Mach is equal to 0.2, and a Reynolds number with respect to diameter D is equal to 100. The constants chosen for this problem were $D = 1$ m, $m = 1$ kg, $k = 0.64$ N/m, $d = 10^{-3}k$ Ns/m, and $l = 1$ m. Fig. 2 shows position h of the cylinder during the computation. Numerical solutions for two cases of our discontinuous Galerkin scheme (RK2-DG1 – second order RK method, first order of polynomials, RK4-DG3 – fourth order RK method, third order of polynomials) are compared to finite volume approximation on very fine grid these solutions indicate very good converge of our scheme.

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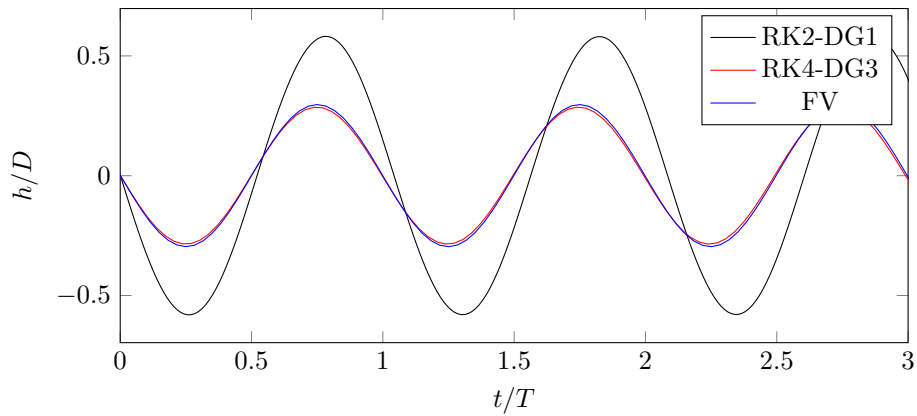


Fig. 2. Position of the cylinder h during selected time period

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