

A bound-preserving multinumetrics scheme for steady-state convection–diffusion equations

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Abstract

We solve the convection–diffusion equation using a coupling of cell-centered finite volume (FV) and discontinuous Galerkin (DG) methods. The domain is divided into disjoint regions assigned to FV or DG, and the two methods are coupled through an interface term. DG is stable and resolves sharp layers in convection-dominated regimes, but it can produce sizable spurious oscillations and is computationally expensive; FV (two-point flux) is low-order and monotone, but inexpensive. We propose a novel adaptive partitioning strategy that automatically selects FV and DG subdomains: whenever the solution’s cell average violates the bounds, we switch to FV on a small neighborhood of that element. Viewed as a natural analog of p -adaptivity, this process is repeated until all cell averages are bound-preserving (up to some specified tolerance). Thereafter, standard conservative limiters may be applied to ensure the full solution is bound-preserving. Standard benchmarks confirm the effectiveness of the adaptive technique.

Keywords: finite elements, discontinuous Galerkin methods, bound-preserving, slope limiters, convection-dominated flow

1. Introduction

As a core transport model, the convection-diffusion equation captures the interplay of flow and diffusion. Transport of scalar fields (e.g., temperature in energy models, concentration in mass transport) is commonly represented by partial differential equations that combine convection by a velocity field with diffusion. When convection outweighs diffusion, the problem enters an convection-dominated regime and solutions develop narrow boundary or internal layers with very steep spatial variation. Such layers occur in both steady and time-dependent cases; however, the latter typically do not introduce similarly small time scales [1]. Our focus concentrates on the steady-state problem.

Mesh-based discretizations struggle when convection overwhelms diffusion because the characteristic layers are much narrower than a feasible mesh spacing. Stabilization is essential for many classical methods; otherwise oscillations contaminate the entire solution [2]. A thorough review of stabilization techniques in the finite element setting is given in [3]. This work studies the coupling of classical cell-centered finite volume (CCFV or FV) [4] and discontinuous Galerkin (DG) methods [5, 6]. Among methods for convection-diffusion, CCFV

and DG are two of the most commonly employed discretizations [7]. Cell-centered finite volume schemes are locally conservative, inexpensive, and under standard two-point flux approximations are monotone, which makes them robust for convection-dominated problems. Their drawbacks are low formal accuracy (typically first order), layer smearing, possible grid-orientation effects, and loss of consistency on non-orthogonal meshes or with strong anisotropy unless more elaborate fluxes or reconstructions are used [8].

Discontinuous Galerkin methods, by contrast, offer high-order accuracy, natural *hp*-adaptivity, compact element-local couplings that parallelize well, and sharp resolution of layers via upwind and penalty fluxes; however, they introduce many more unknowns, in many cases require parameter tuning, are costlier in memory and time, and are not inherently monotone; overshoots/undershoots can appear without limiting. Coupling CCFV and DG capitalizes on their strengths, and results in a locally conservative scheme. The analyses in [9, 10] establish theoretical foundations for CCFV–DG coupling for linear convection-diffusion problems, and [11] demonstrates its effectiveness for nonlinear multiphase flow. Nevertheless, a key unresolved issue is how to decide in an automated way where within the computational domain to deploy FV versus DG, i.e., when and where each scheme should be activated.

We introduce a novel bound-preserving technique for coupling cell-centered finite volume (CCFV) with DG in steady convection-diffusion problems. It targets the spurious overshoots and undershoots that can arise in the high-order DG region. One way to deal with these spurious oscillations is to leverage slope limiters [12]. Most popular slope limiters assume the cell averages are already correct and therefore preserve them during limiting. To repurpose conservative limiters, we focus on generating bound-preserving cell averages. Rather than relying on fix-ups or postprocessing of the DG cell averages, we perform a procedure similar to *p*-adaptivity. If the DG cell average bound-violating (up to some tolerance, e.g., a scaling of machine precision), we revert the approximation to CCFV in a small neighborhood of the troubled cell. Our working ansatz is that a suitable threshold between high-order DG and CCFV yields cell averages that are bound-preserving up to a prescribed tolerance. The intuition is that if we use CCFV over the entire mesh, then the method will be bound-preserving. Once satisfactory cell averages are obtained, any standard conservative slope limiters can be used to ensure the entire solution is bound-preserving.

Instead of manipulating cell averages via postprocessing, we leave them unchanged (we force the cell averages from the approximation to be bound-preserving, without manipulation [13, 14]). The adaptivity yields bound-preserving cell averages up to a prescribed tolerance (and, irrespective of adaptivity, cell averages are generally only accurate up to floating-point round-off). Moreover, *hp*-adaptivity is widely recognized as a state-of-the-art strategy, often attaining a specified error tolerance with markedly fewer degrees of freedom [15]. Our adaptive technique can easily be incorporated into standard *hp*-adaptivity routines, and does not modify the CCFV scheme, DG scheme, or postprocess cell averages.

As an alternative, full DG frameworks may be employed where applicable (e.g., the hybridized LDG method [16] permits piecewise-constant approximations, and leads to a finite volume analog [17]). In addition, the proposed adaptivity can be modified to ensure nodal bound-preservation.

We next survey some related works. Due to the large body of literature surrounding the subject, this survey is non-exhaustive. The work [18] develops nonintrusive postprocessings for DG solutions to suppress spurious oscillations. While the procedure is computationally cheap, it does not enforce bounds preservation. A style of FV-DG limiter computation is introduced in [19, 20]. Here, the authors modify the convective form by using the L^2 projection of the high-order DG solution onto piecewise constants. Their error indicator is based on the jump of the DG approximation across element interfaces, and this is used to determine which cells activate the modified convective form.

Several nonlinear approaches are known to perform well in mitigating spurious oscillations. Artificial viscosity approaches introduce extra diffusion into the discretization, in the hope that it will stabilize as well as prevent unphysical oscillations [21, 22, 23]. In flux-corrected transport, one uses a diffusive, positivity-preserving low-order scheme, adds a high-order correction to improve accuracy, and limits the resulting antidiffusive fluxes to avoid new oscillations [24, 25]. Finally, we mention a review paper on nonlinear bound-preserving modifications (such as variational-inequality formulations) [26, 27]. Each of the aforementioned approaches has trade-offs: some are inexpensive but not bound-preserving; others are bound-preserving but costly; some depart from the original DG discretization or lose conservation of mass; and most lack a formal error analysis.

2. Model problems

The model PDE considered in this work are detailed here. Let $\Omega \subset \mathbb{R}^d$ ($d = 2, 3$) be a bounded polygonal domain.

$$\begin{aligned} -\nabla \cdot (K \nabla u) + \nabla \cdot (\vec{\beta} u) + cu &= f, & \text{in } \Omega, \\ u &= g_D, & \text{on } \Gamma_D, \\ -K \nabla u \cdot \mathbf{n} &= 0, & \text{on } \Gamma_N, \end{aligned} \tag{1}$$

with $\Omega \subset \mathbb{R}^n$ is a bounded domain with polyhedral Lipschitz boundary $\partial\Omega = \Gamma_D \cup \Gamma_N$ (and $\Gamma_D \cap \Gamma_N = \emptyset$). The remaining parameters are as follows: K is the diffusion coefficient (bounded above/below by positive constants), $\vec{\beta}$ is the convection field, c is the reaction coefficient, and f is a source term. In addition, we have nonhomogeneous Dirichlet boundary conditions (g_D) imposed on Γ_D , and homogeneous Neumann conditions on Γ_N , where $\mathbf{n}$ is the unit outward normal to $\partial\Omega$. The inflow boundary is defined as $\Gamma_{\text{inflow}} = \{x \in \Gamma : \vec{\beta}(x) \cdot \mathbf{n} < 0\} \subset \Gamma_D$.

3. Preliminaries

The bounded polygonal domain $\Omega \subset \mathbb{R}^d$ is decomposed into a finite volume region Ω_F and a discontinuous Galerkin region Ω_D with matching interface $\Gamma_{DF} := \partial\Omega_F \cap \partial\Omega_D$, so that $\Omega = \Omega_F \cup \Omega_D$ and $\Omega_F \cap \Omega_D = \emptyset$. Let $\mathcal{T}_h^{FV}$ and $\mathcal{T}_h^{DG}$ be conforming partitions of Ω_F and Ω_D , respectively. FV cells are denoted by $V \in \mathcal{T}_h^{FV}$ (Voronoi in the paper), and DG cells by $W \in \mathcal{T}_h^{DG}$ (simplices/hexahedra or Voronoi). Write $h_V := \text{diam}(V)$, $h_W := \text{diam}(W)$,

$h_{FV} := \max_{V \in \mathcal{T}_h^{FV}} h_V$, $h_{DG} := \max_{W \in \mathcal{T}_h^{DG}} h_W$, and $h := \max(h_{FV}, h_{DG})$. For an edge/face γ , let $|\gamma|$ denote its $(d-1)$ -dimensional measure. We split facets into interior and boundary parts:

$$\Gamma_h^{FV} = \Gamma_{h,I}^{FV} \cup \Gamma_{h,\partial}^{FV}, \quad \Gamma_h^{DG} = \Gamma_{h,I}^{DG} \cup \Gamma_{h,\partial}^{DG}, \quad \Gamma_h^{FV-DG} := \{\gamma \subset \Gamma_h^{FV} \cap \Gamma_h^{DG}\}.$$

Here $\Gamma_{h,I}^{FV}$ are edges interior to Ω_{FV} , $\Gamma_{h,\partial}^{FV}$ lie on $\partial\Omega_{FV} \cap \partial\Omega$, and analogously for DG . We will also use the following notations

$$\begin{aligned} \Gamma_{h,\text{dir}}^{DG} &= \Gamma_h^{DG} \cap \Gamma_D, & \Gamma_{h,\text{dir}}^{FV} &= \Gamma_h^{FV} \cap \Gamma_D, \\ \Gamma_{h,ID}^{DG} &= \Gamma_{h,I}^{DG} \cup \Gamma_{h,\text{dir}}^{DG}, & \Gamma_{h,ID}^{FV} &= \Gamma_{h,I}^{FV} \cup \Gamma_{h,\text{dir}}^{FV}, \\ \Gamma_{h,\partial-}^{DG} &= \Gamma_h^{DG} \cap \Gamma_{\text{inflow}}, & \Gamma_{h,\partial-}^{FV} &= \Gamma_h^{FV} \cap \Gamma_{\text{inflow}}. \end{aligned}$$

For the FV region, we require a so-called two-point flux (TPFA) admissible mesh [28].

Assumption 1 (TPFA admissibility). *Let $\mathcal{T}_h^{FV}$ be a mesh for the finite volume region, and $V, W \in \mathcal{T}_h^{FV}$. We say that $\mathcal{T}_h^{FV}$ is an admissible mesh, in the classical two-point flux finite volume sense if:*

1. *(Cell centers and orthogonality) There exist points $\{x_V\}_{V \in \mathcal{T}_h^{FV}}$ with $x_V \in V$ such that, for every interior edge $\gamma = \partial V \cap \partial W$ with $V \neq W$, the line segment between x_V and x_W is orthogonal to γ .*
2. *(Boundary edges) For each boundary edge $\gamma = \partial V \cap \partial\Omega$, one has $x_V \notin \gamma$. Define y_γ as the orthogonal projection of x_V onto γ .*

Although Assumption 1 may appear to be restrictive, robust unstructured Voronoi generators are well established [29, 30]. In addition, enforcing Assumption 1 ensures optimal convergence of the CCFV method. Using meshes that do not satisfy the assumption forfeits any convergence guarantee. Fig. 1 has an examples Voronoi meshes. An unstructured Voronoi mesh is given in Fig. 1a, and Fig. 1b has an example FV/DG partitioning.

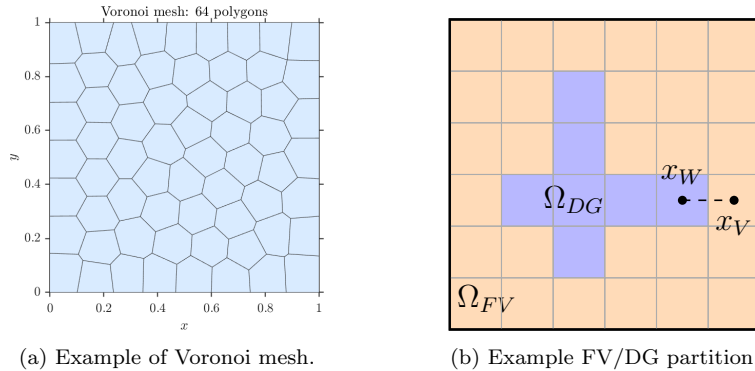


Figure 1: Example Voronoi meshes and illustration of FV (orange) and DG regions (blue), as well as the K -orthogonality property.

For any facet γ fix a unit normal $\mathbf{n}_\gamma$. On $\partial\Omega$ take $\mathbf{n}_\gamma$ outward; on Γ_h^{FV-DG} orient $\mathbf{n}_\gamma$ from the DG side into the FV side; on interior edges choose either orientation consistently. When $\gamma = \partial V \cap \partial W$ lies in Γ_h^{DG} , set $h_\gamma := \max\{\text{diam}(V), \text{diam}(W)\}$ (used in DG penalties).

Define the scalar d_γ for each facet:

$$d_\gamma := \begin{cases} \|x_V - x_W\|_2, & \gamma = \partial V \cap \partial W \in \Gamma_{h,I}^{FV}, \\ \text{dist}(x_V, \gamma) = \|x_V - y_\gamma\|_2, & \gamma = \partial V \cap \partial\Omega \in \Gamma_{h,\partial}^{FV}, \\ \|x_V - y_\gamma\|_2, & \gamma = \partial V \cap \partial W \in \Gamma_h^{FV-DG}, \end{cases}$$

and assume a uniform shape-regularity: there exists $\theta > 0$ such that

$$\begin{aligned} \gamma = \partial V \cap \partial W \in \Gamma_{h,I}^{FV} : \quad d_\gamma &\geq \theta \max(h_V, h_W), \\ \gamma = \partial V \cap \partial\Omega \in \Gamma_{h,\partial}^{FV} : \quad d_\gamma &\geq \theta h_V, \\ \gamma = \partial V \cap \partial W \in \Gamma_h^{FV-DG} : \quad d_\gamma &\geq \theta h_V. \end{aligned}$$

For each FV facet γ (including FV–DG interface facets), let L_γ denote the normal line used by the two-point flux (the straight line orthogonal to γ):

$$L_\gamma = \begin{cases} [x_V, x_W], & \gamma = \partial V \cap \partial W \in \Gamma_{h,I}^F, \\ [x_V, y_\gamma], \quad y_\gamma := \text{proj}_\gamma(x_V), & \gamma \in \Gamma_{h,\partial}^F \text{ or } \gamma \in \Gamma_h^{DF}, \end{cases}$$

and set $d_\gamma := |L_\gamma|$. The harmonic normal average along L_γ is

$$K_\gamma := d_\gamma \left(\int_{L_\gamma} \frac{ds}{K(s)} \right)^{-1}.$$

These K_γ inherit the same upper/lower bounds as K . The finite volume literature often refers to $(|\gamma|/d_\gamma)K_\gamma$ as the TPFA face transmissibility [4].

4. CCFV, DG, and coupled discretization

For a fixed polynomial degree $k \geq 1$ in the DG region, define

$$X_h := \{ v \in L^2(\Omega) : v|_W \in \mathcal{P}_k(W) \quad \forall W \in \mathcal{T}_h^{DG}, v|_V \in \mathcal{P}_0(V) \quad \forall V \in \mathcal{T}_h^{FV} \},$$

where $\mathcal{P}_k(W)$ denotes the space of polynomials on W of total degree at most k .

Functions in X_h can be discontinuous across element interfaces, so we require some jump and average notations to provide meaning to trace values. To make the exposition more clear, given any edge γ , we fix a unit normal vector $\mathbf{n}_\gamma$ to γ . More specifically, we assume that if γ is a boundary facet (belongs to $\partial\Omega$), then $\mathbf{n}_\gamma$ points outward of $\partial\Omega$. If γ belongs to the interface Γ_h^{FV-DG} , then we assume that $\mathbf{n}_\gamma$ points from the DG region into the FV

region. In the following, put $W, V \in \mathcal{T}_h$ so that the vector $\mathbf{n}_\gamma$ points from ∂V into ∂W . Given $u \in X_h$, the jump of u is defined to be:

$$\begin{aligned}\gamma \in \Gamma_{h,I}^{DG} : \quad [u]_\gamma &:= u|_V - u|_W, \\ \gamma \in \Gamma_{h,\partial}^{DG} : \quad [u]_\gamma &:= u|_V, \\ \gamma \in \Gamma_{h,I}^{FV} : \quad [u]_\gamma &:= u(x_V) - u(x_W), \\ \gamma \in \Gamma_{h,\partial}^{FV} : \quad [u]_\gamma &:= u(x_V), \\ \gamma \in \Gamma_h^{FV-DG} : \quad y_\gamma &:= \text{proj}_\gamma(x_V), \quad [u]_\gamma := u|_{\Omega_{DG}}(y_\gamma) - u|_{\Omega_{FV}}(x_V).\end{aligned}$$

The IPDG method requires the notion of the average of $u \in X_h$. Suppose γ is an interior facet in the DG region, with $\gamma = E_1 \cap E_2$, and $E_1, E_2 \in \mathcal{T}_h^{DG}$. We then define

$$\{u\}_\gamma := \frac{1}{2}(u|_{E_1} + u|_{E_2}).$$

Similarly, if γ is a boundary facet in the DG region, with $\gamma \subset \partial E_1$, and $E_1 \in \mathcal{T}_h^{DG}$, we define

$$\{u\}_\gamma := u|_{E_1}.$$

Given an interior facet γ , shared by elements V and W (with $\mathbf{n}_\gamma$ oriented from V to W), we define the upwind value of u to be

$$u^\uparrow = \begin{cases} u|_V & \text{if } \vec{\beta} \cdot \mathbf{n}_\gamma \geq 0, \\ u|_W & \text{if } \vec{\beta} \cdot \mathbf{n}_\gamma < 0. \end{cases}$$

For $E_1, E_2 \in \mathcal{T}_h^{DG}$, let $\gamma = \partial E_1 \cap \partial E_2$, and $h_\gamma = \max(\text{diam}(V), \text{diam}(W))$. With parameters $\sigma > 0$ and $\epsilon \in \{-1, 0, 1\}$, define the following bilinear forms for $u, v \in X_h$:

$$\begin{aligned}a_{DG}(u, v) &= \sum_{W \in \mathcal{T}_h^{DG}} \int_W K \nabla u \cdot \nabla v - \sum_{\gamma \in \Gamma_{h,ID}^{DG}} \int_\gamma \{K \nabla u \cdot \mathbf{n}_\gamma\} [v] + \epsilon \sum_{\gamma \in \Gamma_{h,ID}^{DG}} \int_\gamma \{K \nabla v \cdot \mathbf{n}_\gamma\} [u] \\ &+ \sum_{\gamma \in \Gamma_{h,ID}^{DG}} \frac{\sigma}{h_\gamma} \int_\gamma [u][v] - \sum_{W \in \mathcal{T}_h^{DG}} \int_W (\vec{\beta} \cdot \nabla v) u + \sum_{\gamma \in \Gamma_{h,I}^{DG}} \int_\gamma (\vec{\beta} \cdot \mathbf{n}_\gamma) u^\uparrow [v]_\gamma \\ &+ \sum_{\gamma \in \Gamma_{h,\partial-}^{DG}} \int_\gamma (\vec{\beta} \cdot \mathbf{n}_\gamma) uv + \sum_{W \in \mathcal{T}_h^{DG}} \int_W cuv,\end{aligned}$$

$$\begin{aligned}a_{FV}(u, v) &= \sum_{\gamma \in \Gamma_{h,ID}^{FV}} \frac{|\gamma|}{d_\gamma} K_\gamma [u]_\gamma [v]_\gamma + \sum_{\gamma \in \Gamma_{h,I}^{FV}} \int_\gamma (\vec{\beta} \cdot \mathbf{n}_\gamma) u^\uparrow [v]_\gamma + \sum_{\gamma \in \Gamma_{h,\partial-}^{FV}} \int_\gamma (\vec{\beta} \cdot \mathbf{n}_\gamma) uv \\ &+ \sum_{V \in \mathcal{T}_h^{FV}} \int_V cuv,\end{aligned}$$

$$a_{FG}(u, v) = \sum_{\gamma \in \Gamma_h^{FV-DG}} \frac{|\gamma|}{d_\gamma} K_\gamma [u]_\gamma [v]_\gamma + \sum_{\gamma \in \Gamma_h^{FV-DG}} \int_\gamma (\vec{\beta} \cdot \mathbf{n}_\gamma) u^\uparrow [v]_\gamma,$$

$$a(u, v) = a_{DG}(u, v) + a_{FV}(u, v) + a_{FG}(u, v).$$

For the linear form, we have

$$\begin{aligned} \ell(v) = & \epsilon \sum_{\gamma \in \Gamma_{h,\text{dir}}^{DG}} \int_{\gamma} \frac{\sigma}{h_{\gamma}} (\nabla v \cdot \mathbf{n}_{\gamma}) g_D + \sum_{\gamma \in \Gamma_{h,\text{dir}}^{DG}} \int_{\gamma} \frac{\sigma}{h_{\gamma}} g_D v - \sum_{\gamma \in \Gamma_{h,\partial-}^{DG}} \int_{\gamma} (\vec{\beta} \cdot \mathbf{n}_{\gamma}) g_D v \\ & + \sum_{W \in \mathcal{T}_h^{DG}} \int_W f v + \sum_{\gamma \in \Gamma_{h,\text{dir}}^{FV}} \frac{|\gamma|}{d_{\gamma}} K_{\gamma} g_D v|_{\gamma} - \sum_{\gamma \in \Gamma_{h,\partial-}^{FV}} \int_{\gamma} (\vec{\beta} \cdot \mathbf{n}_{\gamma}) g_D v. \end{aligned}$$

The discrete coupled FV-DG scheme for (1) is: find $u_h \in X_h$ so that

$$a(u_h, v_h) = \ell(v_h), \quad \forall v_h \in X_h. \quad (2)$$

If $\mathcal{T}_h^{DG} = \emptyset$, then (2) reduces to the classical cell-centered finite-volume method (CCFV). Similarly, if $\mathcal{T}_h^{FV} = \emptyset$, then (2) reduces to an interior-penalty DG (IPDG) method with upwinding for the convection. In Section 6, we compare and contrast the three schemes: full CCFV ($\mathcal{T}_h^{DG} = \emptyset$), full DG ($\mathcal{T}_h^{FV} = \emptyset$), and FV-DG ($\mathcal{T}_h^{FV} \neq \emptyset$, $\mathcal{T}_h^{DG} \neq \emptyset$).

5. Description of adaptive technique

Here we describe the bound-preserving mechanism used in this paper. The method runs in two phases:

- (i) perform standard h -adaptivity until the accuracy estimator meets tolerance;
- (ii) repeatedly scan cell averages against the admissible interval $I = [u_* - \delta_{\text{under}}, u^* + \delta_{\text{over}}]$.

Any violating cells (and their facet/vertex neighbors, see Fig. 2) are reassigned to FV, and update the partition: $\mathcal{T}_h^{FV} \cup \mathcal{T}_h^{DG}$. The FV and DG regions are updated until the bound-violation metric falls below tolerance, then apply a conservative slope limiter on the DG region. We also consider a slight modification that scans nodal values instead of cell averages (see Variant 2 in subsection 5.2).

5.1. Variant 1: Bound-driven FV/DG partitioning (fixed slack)

Pick slack parameters $\delta_{\text{under}}, \delta_{\text{over}} > 0$ and tolerances $\varepsilon_h, \varepsilon_{\text{bp}} > 0$. Define the admissible interval $I = [u_* - \delta_{\text{under}}, u^* + \delta_{\text{over}}]$ and let $\mathcal{N}(E)$ denote the set that contains the facet and vertex neighbors of E .

1. **Accuracy phase (h -adapt).** While $\eta_h(\mathcal{T}_h) > \varepsilon_h$, refine/coarsen $\mathcal{T}_h$ and recompute u_h .
2. **Bound-screened FV/DG partitioning.** Repeat *until* $\eta_{\text{bp}} \leq \varepsilon_{\text{bp}}$:
 - (a) Compute cell averages $\bar{u}_h|_E$ for all $E \in \mathcal{T}_h$.
 - (b) Form the violation set $V := \{E : \bar{u}_h|_E \notin I\}$ and expand $V \leftarrow V \cup \mathcal{N}(V)$.
 - (c) Set $\mathcal{T}_h^{FV} \leftarrow V$ and $\mathcal{T}_h^{DG} \leftarrow \mathcal{T}_h \setminus V$; recompute u_h and evaluate

$$\eta_{\text{bp}} := \max_{E \in \mathcal{T}_h} \text{dist}(\bar{u}_h|_E, I), \quad \text{dist}(x, I) := \max\{u_* - x, 0, x - u^*\}.$$

3. **DG limiting.** Apply a conservative slope limiter on $\mathcal{T}_h^D$.

5.2. Variant 2: Node-screened FV/DG partitioning.

Let $I = [u_* - \delta_{\text{under}}, u^* + \delta_{\text{over}}]$ and, for each cell $E \in \mathcal{T}_h$, let $\mathcal{X}(E)$ denote the set of nodal points to be screened (e.g., the vertices of E , or quadrature points). Define

$$\eta_{\text{bp}}^{\text{node}} := \max_{E \in \mathcal{T}_h} \max_{x \in \mathcal{X}(E)} \text{dist}(u_h|_E(x), I), \quad \text{dist}(z, I) := \max\{u_* - z, 0, z - u^*\}.$$

1. **Accuracy phase (h -adapt).** While $\eta_h(\mathcal{T}_h) > \varepsilon_h$, refine/coarsen $\mathcal{T}_h$ and recompute u_h .
2. **Bound-screened FV/DG partitioning (node-based).** Repeat *until* $\eta_{\text{bp}}^{\text{node}} \leq \varepsilon_{\text{bp}}$:
 - (a) For every $E \in \mathcal{T}_h$, evaluate $u_h|_E(x)$ at all $x \in \mathcal{X}(E)$.
 - (b) Form the violation set

$$V := \left\{ E \in \mathcal{T}_h : \exists x \in \mathcal{X}(E) \text{ s.t. } u_h|_E(x) \notin I \right\}.$$

- (c) Expand $V \leftarrow V \cup \mathcal{N}(V)$ and set $\mathcal{T}_h^{FV} \leftarrow V$, $\mathcal{T}_h^{DG} \leftarrow \mathcal{T}_h \setminus V$.
 - (d) Recompute u_h on the new partition and update $\eta_{\text{bp}}^{\text{node}}$.
3. **DG limiting.** Apply a conservative slope limiter on $\mathcal{T}_h^{DG}$.

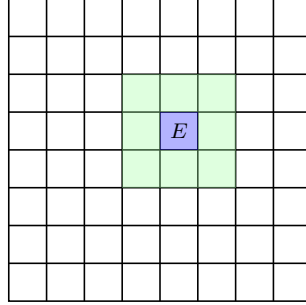


Figure 2: Element E marked for adaptivity (blue), and its vertex neighborhood $\mathcal{N}(E)$ (green).

Remark 1. Step 3 may be skipped if it is enough that the nodal values on $\mathcal{X}(E)$ are bound-preserving. For $k > 1$ DG, nodal bounds do not preclude between node overshoots; $u_h|_E(x)$ can violate the bounds for $x \in E \setminus \mathcal{X}(E)$.

Remark 2. Through various numerical tests, we find that taking $\mathcal{N}(E)$ to be the vertex neighbors of E provides good results.

5.3. Slope limiter

To make the exposition more clear, let $\Omega \subset \mathbb{R}^2$. We use a vertex limiter from [31]. For an element $E \in \mathcal{T}_h$ with centroid $c_0 = (x_0, y_0)$ and vertices $\{v_i = (x_i, y_i)\}_{i=1}^M$, the piecewise linear DG solution is

$$u_h|_E(x, y) = \bar{u}_E + u_x(x - x_0) + u_y(y - y_0),$$

where $(u_x, u_y) = \nabla u_h|_E$ is constant on E . For each vertex v_i , let $\omega(v_i)$ be the patch of elements that share v_i . Define vertexwise bounds

$$u_i^{\min} = \min_{L \in \omega(v_i)} \bar{u}_L, \quad u_i^{\max} = \max_{L \in \omega(v_i)} \bar{u}_L.$$

If v_i lies on a Dirichlet boundary, also include the boundary value $g_D(v_i)$ when forming $u_i^{\min}, u_i^{\max}$. The objective of the limiter is to construct a new polynomial

$$\tilde{u}_h|_E(x, y) = \bar{u}_E + \alpha_x u_x (x - x_0) + \alpha_y u_y (y - y_0), \quad \alpha_x, \alpha_y \in [0, 1],$$

such that the vertex values satisfy the discrete maximum principle

$$u_i^{\min} \leq \tilde{u}_h(v_i) \leq u_i^{\max} \quad \text{for all } i = 1, \dots, M.$$

The scalars α_x and α_y are computed as:

1. Choose $\alpha_x \in [0, 1]$ so that $u_i^{\min} \leq u_0 + \alpha_x u_x (x_i - x_0) \leq u_i^{\max}$ for all i . Define nodal factors

$$\alpha_{x,i} = \begin{cases} \min\left(1, \frac{u_i^{\max} - u_0}{u_x (x_i - x_0)}\right), & u_x (x_i - x_0) > 0, \\ 1, & u_x (x_i - x_0) = 0, \\ \min\left(1, \frac{u_i^{\min} - u_0}{u_x (x_i - x_0)}\right), & u_x (x_i - x_0) < 0, \end{cases} \quad \alpha_x = \min_{1 \leq i \leq M} \alpha_{x,i}.$$

Set the prelimiting polynomial $\hat{u}_h(x, y) = u_0 + \alpha_x u_x (x - x_0)$ and its vertex values $\hat{u}_i = \hat{u}_h(v_i)$, which are used in step 2.

2. Choose $\alpha_y \in [0, 1]$ so that $u_i^{\min} \leq \hat{u}_i + \alpha_y u_y (y_i - y_0) \leq u_i^{\max}$ for all i . Define

$$\alpha_{y,i} = \begin{cases} \min\left(1, \frac{u_i^{\max} - \hat{u}_i}{u_y (y_i - y_0)}\right), & u_y (y_i - y_0) > 0, \\ 1, & u_y (y_i - y_0) = 0, \\ \min\left(1, \frac{u_i^{\min} - \hat{u}_i}{u_y (y_i - y_0)}\right), & u_y (y_i - y_0) < 0, \end{cases} \quad \alpha_y = \min_{1 \leq i \leq M} \alpha_{y,i}.$$

3. Form the limited polynomial:

$$\bar{u}_h|_E(x, y) = \bar{u}_E + \alpha_x u_x (x - x_0) + \alpha_y u_y (y - y_0).$$

Theorem 5.1. *Let u_h be the solution to the FV-DG problem in (2) after application of the adaption technique Variant 1 from Section 5). Then, the cell-averages of u_h are bound-preserving up to some tolerance δ .*

Proof. Variant 1 of the adaption technique is guaranteed to terminate in a finite number of steps because the pessimistic choice $\mathcal{T}_h^{DG} = \emptyset$ (expand the neighborhoods until $\mathcal{T}_h = \mathcal{T}_h^{DG} \cup \mathcal{T}_h^{FV} = \mathcal{T}_h^{FV}$) ensures the scheme reverts to CCFV on the entire mesh; and the CCFV discretization is monotone [32, 4]. $\square$

Remark 3. Empirically, we find that the DG region is typically sizable (30%-90% of mesh elements), though this varies with the tolerance δ , the mesh size, and the underlying problem.

6. Numerical experiments

Unless otherwise stated, all experiments for DG have $k = 1$ (piecewise linear approximation), $\epsilon = -1$ (symmetric IPDG) and $\sigma = 14$. All computational meshes are unstructured (Voronoi) generated with CGAL [33] and PolyMesher [34]. The meshes are matching at the interface between the FV and DG regions.

Variant 1 from Section 5 governs the cell averages, while the limiter in Section 5.3 constrains the higher-order DG modes/slopes. While distinct slack parameters for undershoot and overshoot are admissible, in the experiments we take a single value, $\delta := \delta_{\text{under}} = \delta_{\text{over}}$. We take $f = c = 0$ in the model PDE (1). A suite of standard benchmark problems is presented in Sections 7, 8, 9, and 10. Section 11 summarizes the results and presents supplementary metrics: iteration counts, and the size of spurious oscillations.

7. Triple layer

We consider a triple layer problem, which is a common benchmark for convection-dominated flows. No exact solution can be easily written down for this problem. The domain is $\Omega = [0, 2]^2$, and $K = 10^{-6}$, with the advective field $\vec{\beta} = [y, (1 - x)^2]^T$. On the domain $\Gamma_N := \{2\} \times (0, 1) \cup (0, 2) \times \{1\}$ homogeneous Neumann conditions are imposed. For the remaining boundaries, Dirichlet conditions are imposed:

$$g_D(x, y) = \begin{cases} 1, & \text{if } x \in (1/8, 1/2), y = 0, \\ 2, & \text{if } x \in (1/2, 3/4), y = 0, \\ 0, & \text{elsewhere on } \Gamma_D \setminus \Gamma_N. \end{cases}$$

Due to the small amount of diffusion the discontinuous profile is basically transported along the characteristic curves leading to sharp characteristic interior layers.

Baseline tests for this benchmark are conducted in Fig. 3. The CCFV method is monotone and bound-preserving, at the cost of reduced feature sharpness. Conversely, the DG solution offers greater detail but is not bound-preserving. Figs. 3c and 3f visualize the spatial distribution of DG cell average overshoots and undershoots. Mesh refinement alone does not eliminate violations in the cell average.

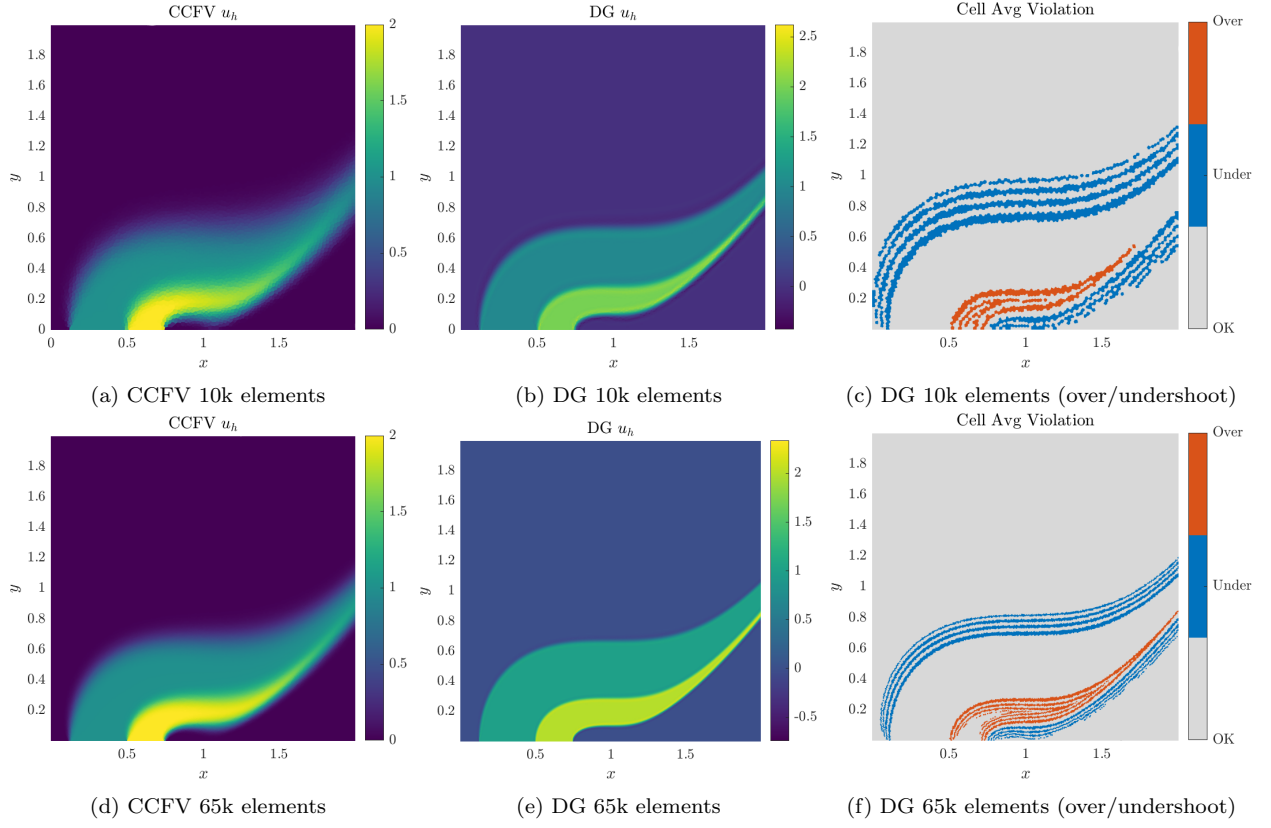


Figure 3: Baseline tests from Section 7, full CCFV ($\mathcal{T}_h^{DG} = \emptyset$) or full DG ($\mathcal{T}_h^{FV} = \emptyset$) with various mesh sizes.

We next examine the FV-DG adaptive results in Fig. 4. The FV-DG solution is markedly better than CCFV and improves further with mesh refinement (Figs. 4a–4c). In addition, the FV-DG solution bound-preserving up to the specified tolerance. Compared across refinements, the DG region becomes progressively larger (Figs. 4d–4f).

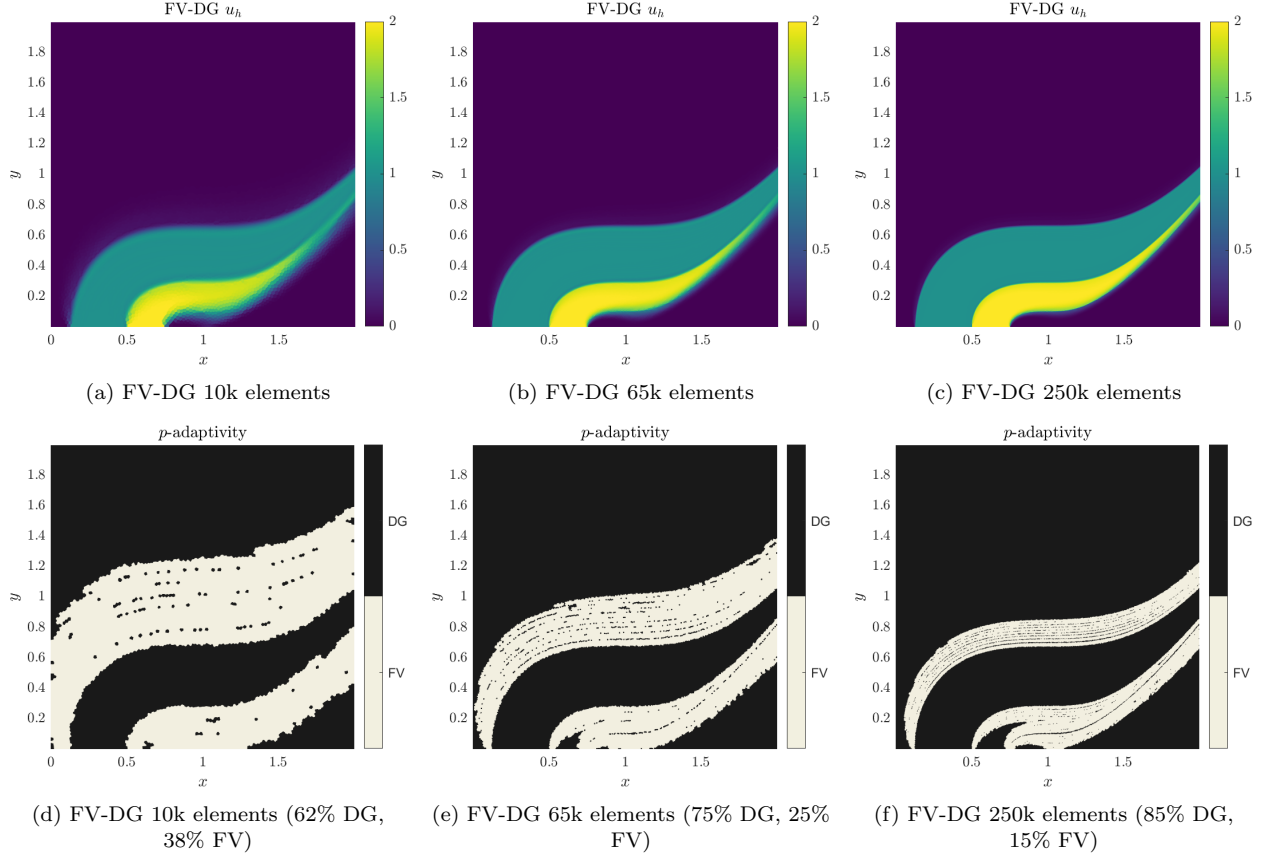


Figure 4: Section 7 adaptive FV-DG with $\delta = 10^{-7}$ and various mesh sizes.

7.1. Sub-extremal oscillations

In practice, the adaptive procedure functioned robustly across all test cases. We note that the adaptive technique specifically enforces bound-preservation, so it is possible the approximations have spurious oscillations within the bounds. For this benchmark problem, the exact solution to (1) obeys $0 \leq u \leq 2$. When the unmodified approximation (Figs. 3b and 3e) yields $\bar{u}_h \approx 1$, the indicator based on cell-average bounds will not flag that region (no overshoot/undershoot). To better explain this, Fig. 5 shows solution profiles along $x = 1$. The FV-DG scheme is bound-preserving and resolves the profile more sharply than CCFV, though small local oscillations remain (e.g., near $x \approx 0.3$).

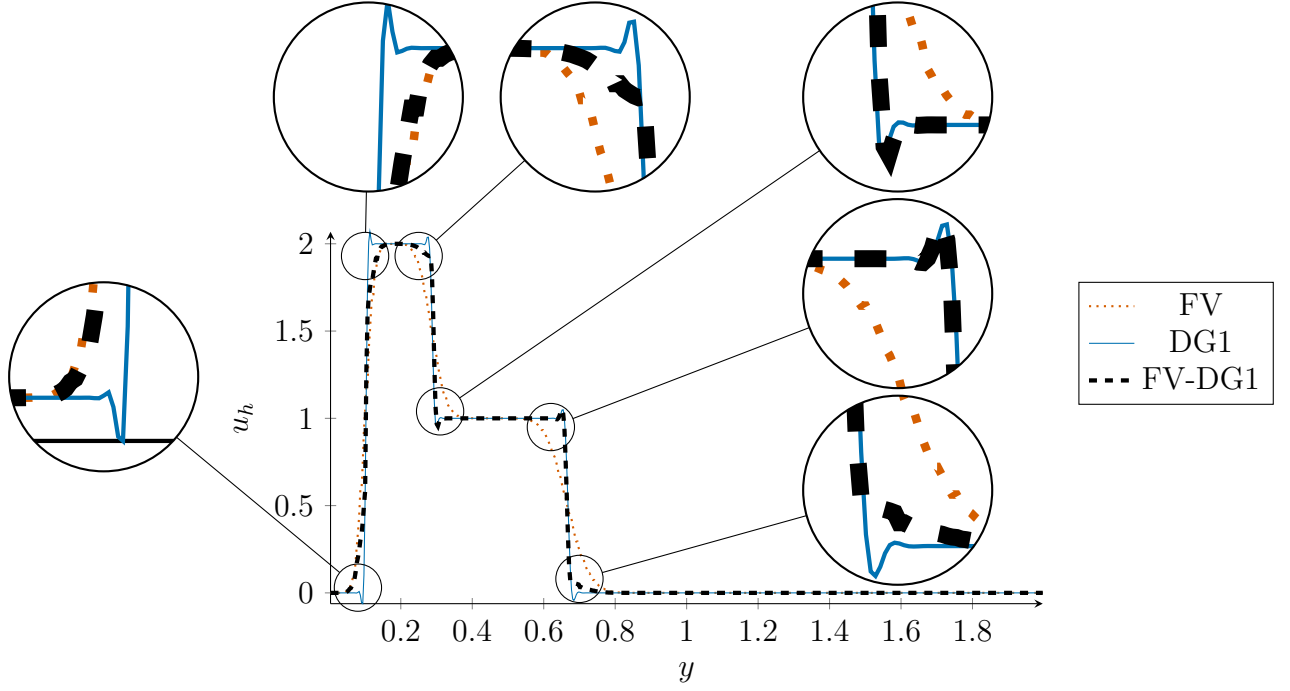


Figure 5: Profile along $x = 1$ for the triple layer problem, with magnifications. Fixed mesh with 262k elements.

Figure 6 compares the adaptive FV-DG scheme for varying polynomial degree k . The resulting FV and DG partitions are largely consistent across k , with only small differences.

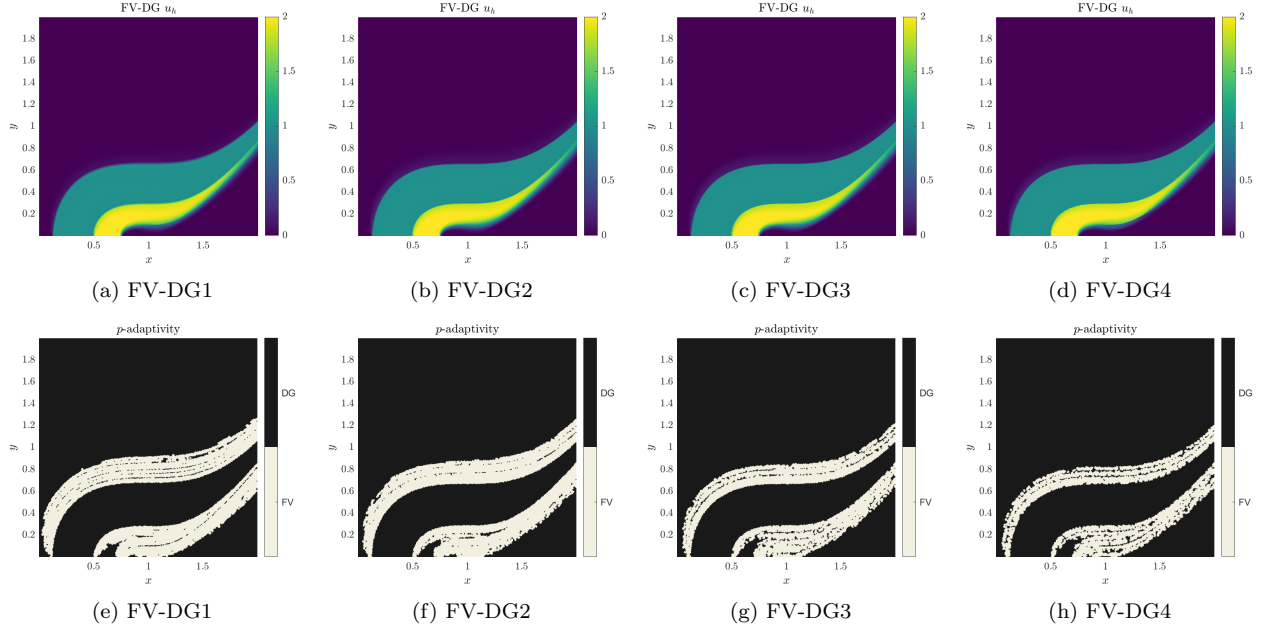


Figure 6: Comparison of the adaptive FV-DG scheme for different polynomial degrees.

8. Convection-dominated problem (L-shaped domain)

The next benchmark was taken from [35]. No exact solution can be easily written down for this problem. The domain is L-shaped with $\Omega = [0, 4]^2 \setminus [0, 2]^2$. The convective vector field is given as $\vec{\beta} = [y, -x]^T$ and $K = 10^{-6}$. The PDE boundary conditions are as follows:

$$\begin{aligned} u &= 1, & \text{on } \{(x, y) : x = 0, y \in (2, 4)\}, \\ K \nabla u \cdot \mathbf{n} &= 0, & \text{on } \{(x, y) : x = 4, y \in (0, 4)\} \cup \{(x, y) : x \in (2, 4), y = 0\}, \\ u &= 0, & \text{elsewhere.} \end{aligned}$$

The solution of this problem exhibits two boundary layers along $\{(x, y) : x \in (0, 4), y = 4\}$ and $\{(x, y) : x \in (0, 2), y = 2\}$. Moreover, the solution obeys $0 \leq u \leq 1$ and contains two circular-shaped interior layers [36].

A series of tests are conducted. We have some baseline experiments, in particular, high-order DG on all mesh elements, and CCFV on all mesh elements. Fig. 7 contrasts the CCFV and DG schemes. As seen in panels 7a and 7d, the CCFV solution is smeared, indicating stronger numerical diffusion. Conversely, DG is less diffusive and captures the sharp features (Figs. 7b, 7e). However, the DG cell averages exhibit overshoots/undershoots, as evidenced in Figs. 7c and 7f. Further h -refinement yields only marginal improvement.

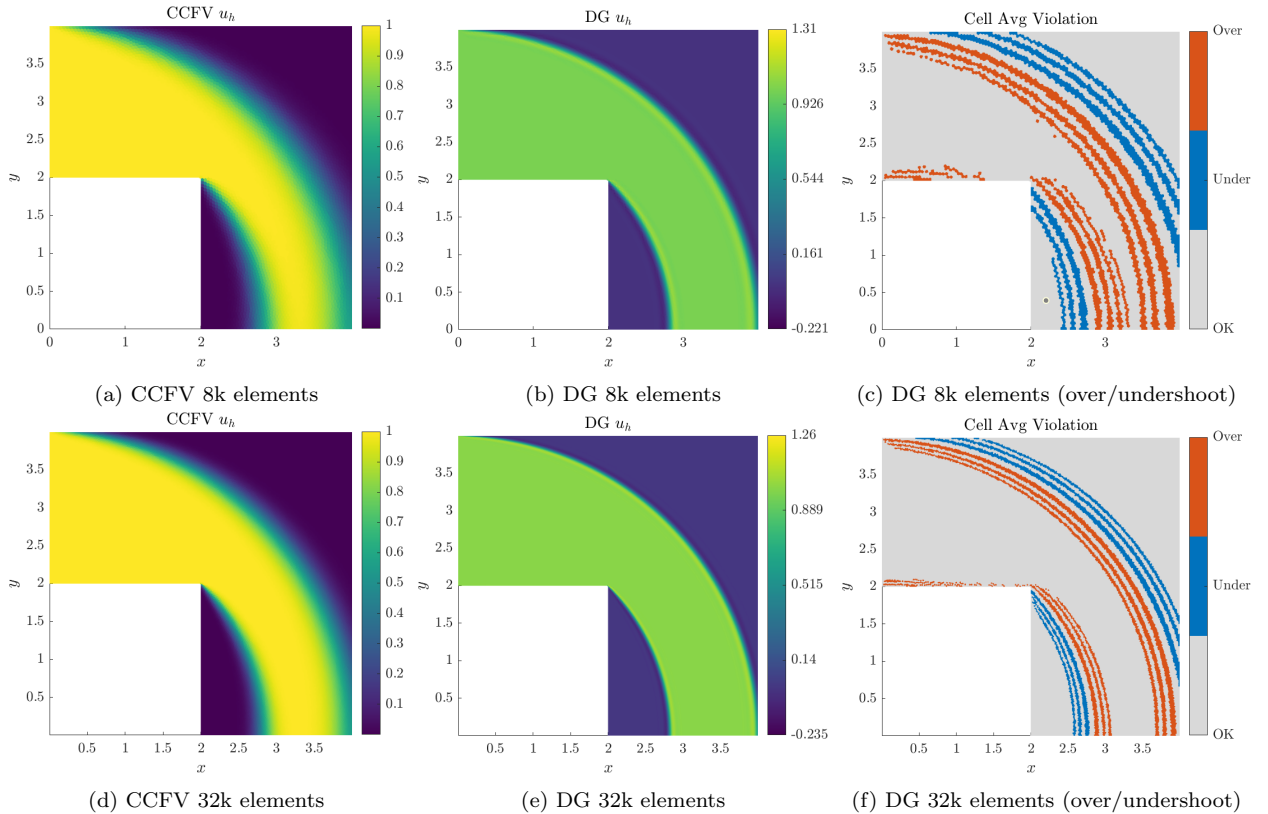


Figure 7: Baseline tests - use CCFV everywhere or DG everywhere with various mesh sizes.

8.1. FV-DG adapted scheme with relaxed tolerance

Next, we compare the adaptive FV-DG coupled scheme with $\delta = 10^{-6}$. Figs. 8a, 8b, and 8c show that the FV-DG solution is considerably less diffusive than the CCFV solution. The second row of Fig. 8 delineates the FV and DG regions. For this benchmark, the DG region expands with mesh refinement, occupying an increasing fraction of the domain.

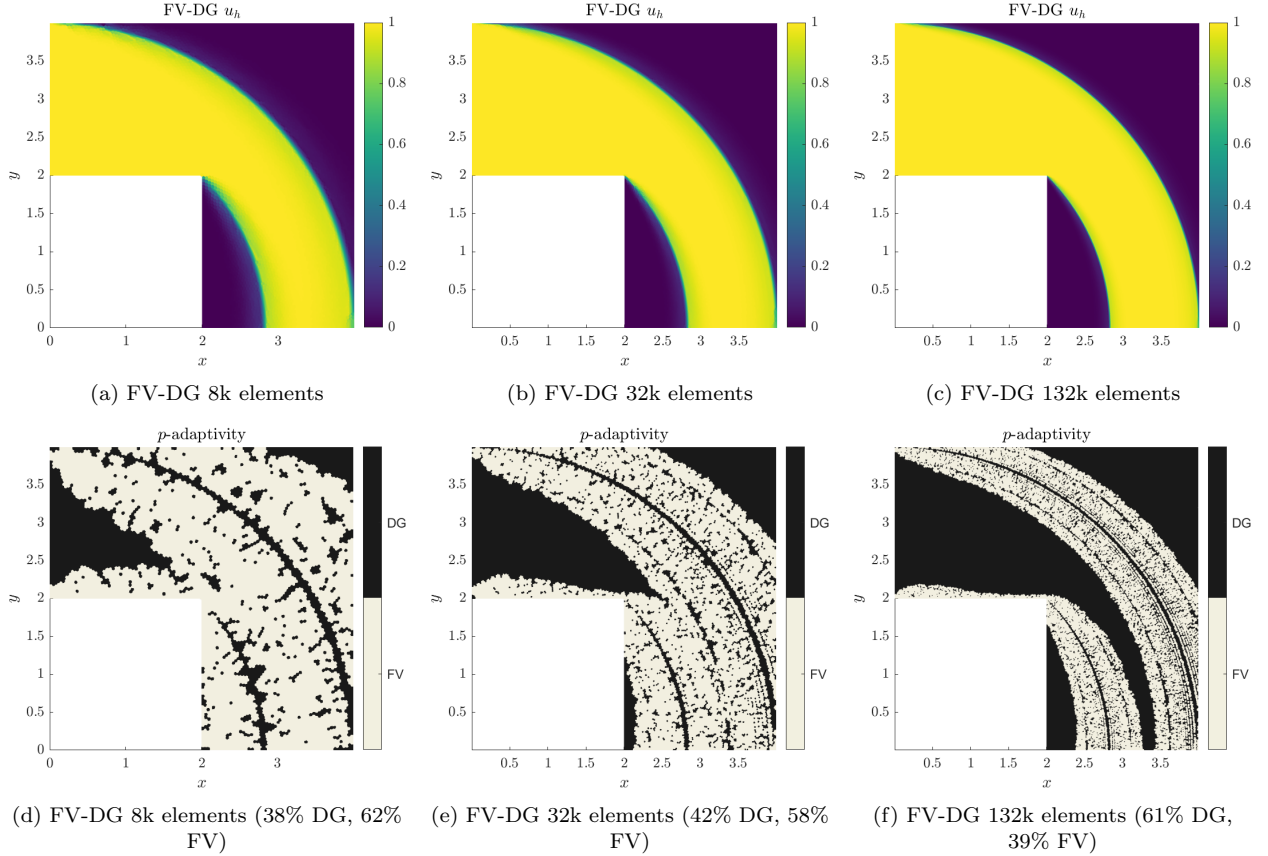


Figure 8: Adaptive scheme for FV-DG with $\delta = 10^{-6}$ and various mesh sizes.

8.2. FV-DG adapted scheme with stricter tolerances

With $\delta = 10^{-9}$ and $\delta = 10^{-13}$, comparable results are obtained by the FV-DG coupled scheme (see Fig. 9). At identical mesh resolution, the technique provides a much less diffusive solution than CCFV (see the third row of Fig. 9). Specifically, Figs. 9f (132k elements) and 9e (532k elements) indicate that CCFV remains markedly more diffusive than the adaptive FV-DG method.

The adaptive strategy delivers good results for this challenging benchmark. The DG cell count is controlled by δ ; typically, decreasing δ reduces the DG region. For instance, Fig. 9b with $\delta = 10^{-9}$ has a space composition of (61% DG, 39% FV), whereas Fig. 9d with $\delta = 10^{-13}$ has a space composition of (53% DG, 47% FV). A stricter tolerance for δ increases the FV region.

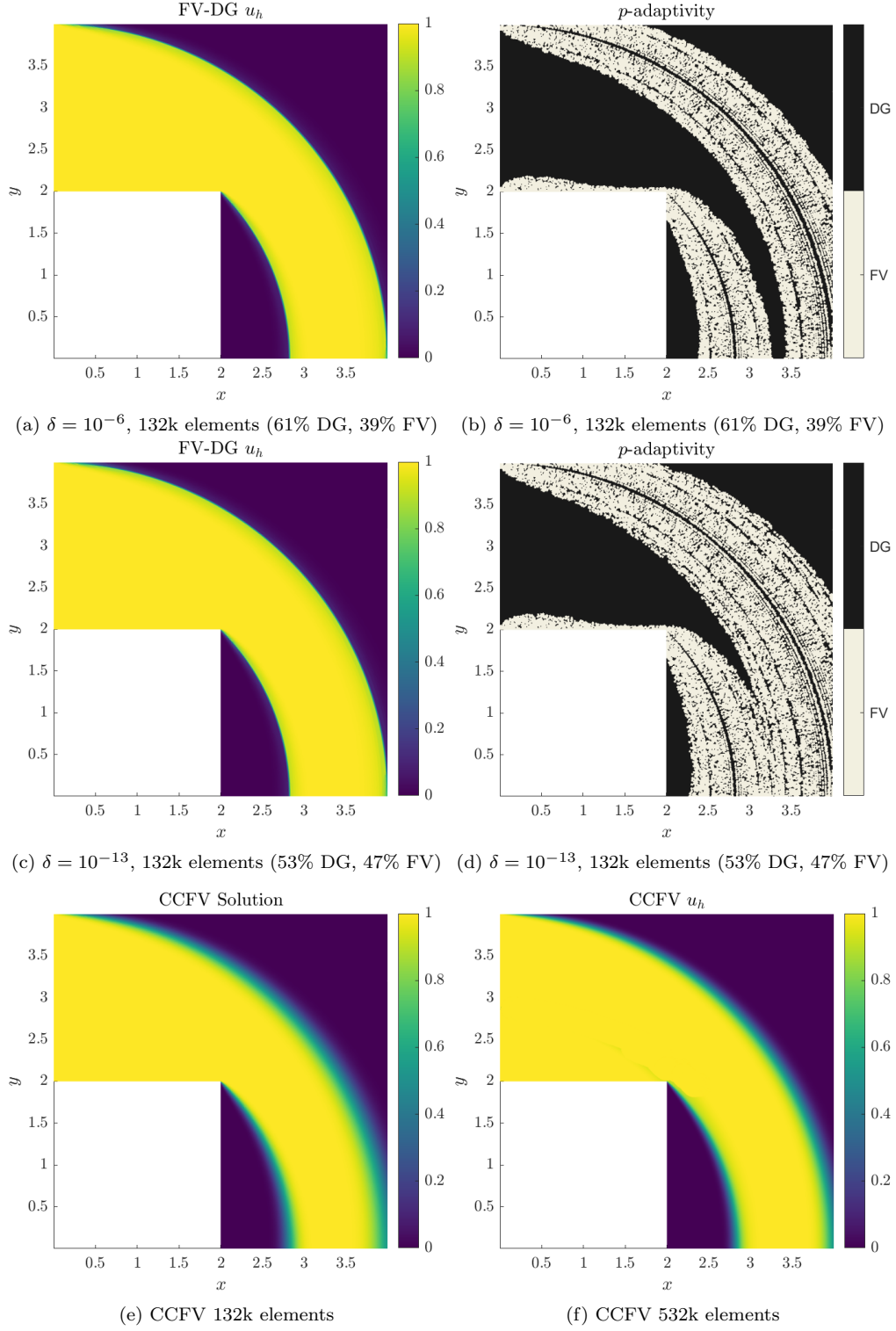


Figure 9: CCFV vs adaptive scheme for FV-DG with $\delta = 10^{-9}$ (row 1) and $\delta = 10^{-13}$ (row 2).

8.3. FV–DG adapted scheme with h -adaptivity

To demonstrate that the adaptive technique of Section 5 accommodates standard h -adaptivity, we set $\delta = 10^{-13}$. Due to the nature of the selection of DG and FV regions, we assume that the entire mesh consists of Voronoi cells. This ensures that the TPFA mesh assumption holds (see Assumption 1). After marking elements for refinement, we randomly place Voronoi sites in a small neighborhood of the marked elements' centroids. We then pass the new sites, together with the existing ones, to a Voronoi mesh generator [33, 34, 37].

Given u_h on an initial mesh $\mathcal{T}_h$, we drive mesh refinement with a jump-based indicator of the solution across element faces [20]:

$$\eta_E^2 := \sum_{\gamma \in \partial E} \int_{\gamma} \frac{[u_h]_{\gamma}^2}{h_{\gamma}} ds, \quad \eta_h(\mathcal{T}_h) = \left(\sum_E \eta_E^2 \right)^{1/2}.$$

Fig. 10 shows: (a) the initial mesh; (b) the mesh after h -refinement; (c) a zoom near $x = 2$; (d) the FV–DG solution on the adapted mesh; (e) the FV and DG subregions; and (f) the cells where the full DG method violates the cell-average constraint. Thus, the adaptive technique in Section 5 fits naturally within standard hp -adaptive finite element strategies.

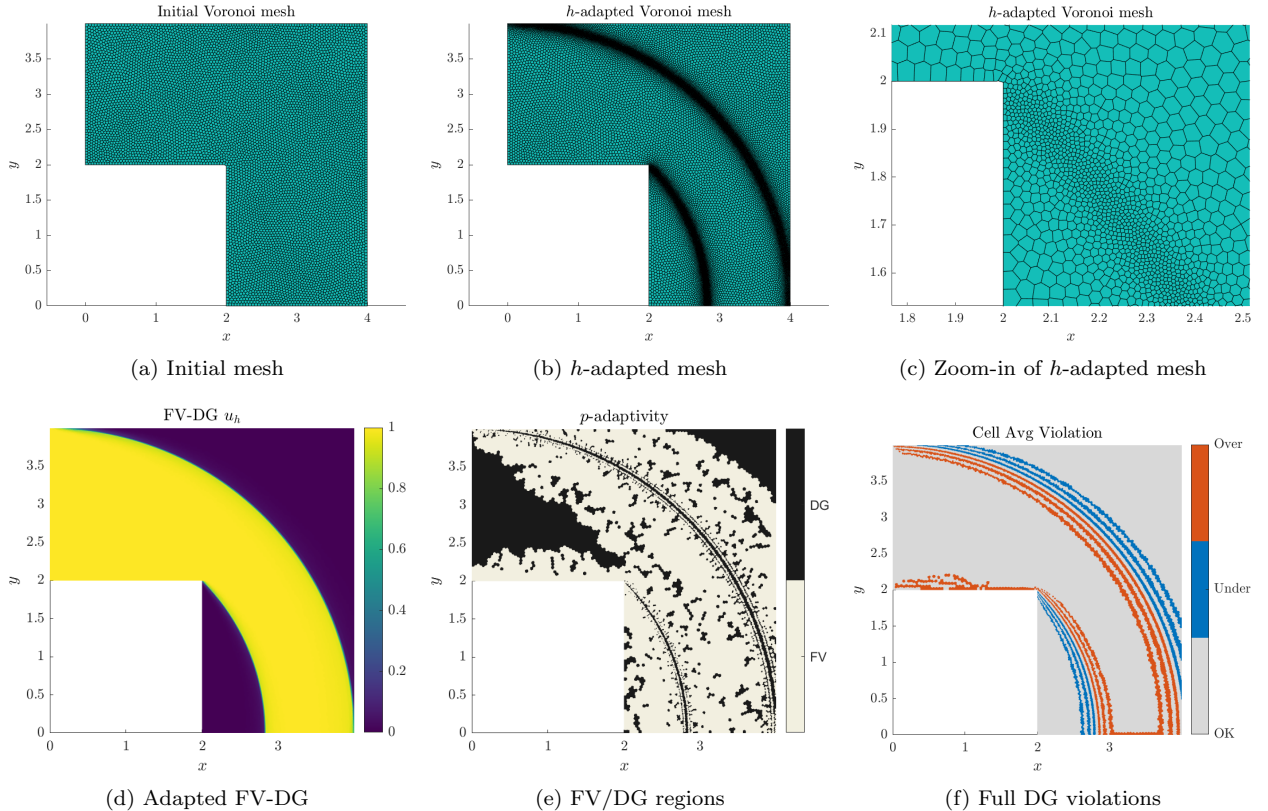


Figure 10: Results for an h -adapted mesh.

9. Internal and boundary layers

We adopt the test case from [38]. Because the problem is convection-dominated, the solution should carry the profile imposed at $x = 0$ downstream in the direction of $\vec{\beta}$. The diffusion coefficient is set as $K = 10^{-4}$, and $\vec{\beta} = [\cos(\pi/3), -\sin(\pi/3)]^T$.

$$\begin{aligned} u &= 0, & \text{on } \{(x, y) : y = 0, x \in (0, 1)\}, \\ u &= 1, & \text{on } \{(x, y) : y = 1, x \in (0, 1)\}, \\ u &= \frac{1}{2} + \frac{1}{\pi} \arctan(10^4(y - 0.7)), & \text{on } \{(x, y) : x = 0, y \in (0, 1)\}, \\ u &= 0 & \text{on } \{(x, y) : x = 1, y \in (0, 1)\}. \end{aligned}$$

For DG1, we put $\epsilon = -1$ and $\sigma = 14$. Fig. 11 contrasts the full CCFV and DG1 schemes. CCFV shows greater numerical diffusion, while DG1 yields sharper features at the cost of superious oscillations.

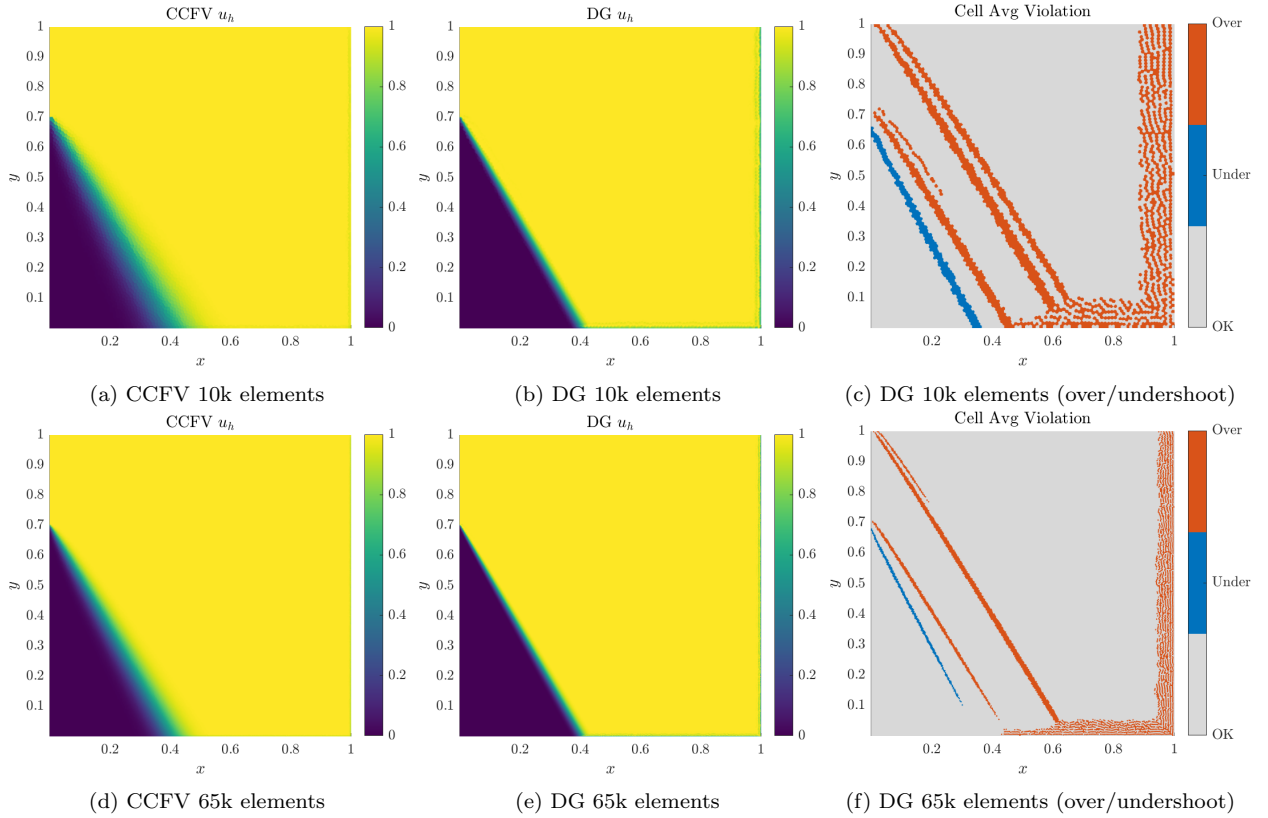


Figure 11: Baseline tests from Section 9 - use CCFV everywhere or DG everywhere with various mesh sizes.

In Fig. 12 we compare FV-DG1 and FV-DG2 using the OBB method [39] with $\epsilon = 1$, $\sigma = 0$ on interior edges, and $\sigma = 1$ on boundary edges. In both configurations, DG covers most of the domain, while boundary layers near $x = 1$ and $y = 0$ are adaptively assigned

FV cells to limit overshoot. Mesh refinement and increasing the DG degree from $k = 1$ to $k = 2$ both enlarge the DG regions.

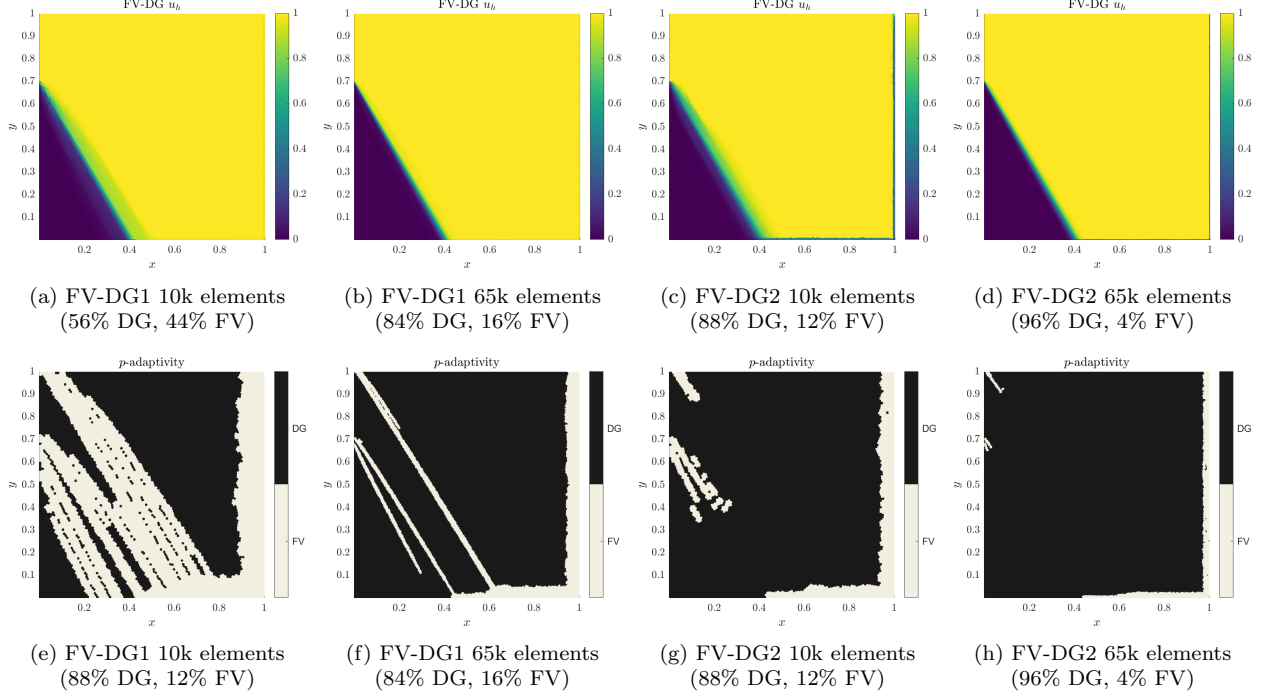


Figure 12: Adaptive scheme for FV-DG with $\delta = 10^{-9}$.

The adaptive technique is quite effective for this test case.

10. Hemker problem (convection-diffusion)

As introduced in [40], the Hemker problem has become a standard benchmark for convection–diffusion models, exhibiting many of the salient features encountered in applications. It describes the transport of energy from a body through a channel. The solution remains within $[0, 1]$. Sharp boundary layers form at the interior boundary, and interior layers propagate from the body in the direction of convection.

The domain is a rectangle with a circular puncture at the origin with radius one: $[(-3, 9) \times (-3, 3)] \setminus \{(x, y) : x^2 + y^2 \leq 1\}$. A Dirichlet condition of one is supplied on the circle boundary. Homogeneous Dirichlet boundary conditions are supplied at $x = -3$. The remaining boundaries are no flow (zero Neumann conditions). A value of $K = 10^{-8}$ is taken to ensure a strongly convection-dominated regime. Here and $\vec{\beta} = [1, 0]^T$.

Consistent with the other benchmarks, we consider baseline approximations using full CCFV and full DG. Fig. 13 presents results at multiple mesh resolutions. Full DG exhibits considerable bound violations (undershoot/overshoot); notably, the DG cell averages are predominantly overshooting. By contrast, CCFV is markedly more diffusive.

With $\delta = 10^{-9}$, the FV–DG adaptive scheme (Fig. 14) delivers clear gains over Figs. 13a and 13d: boundary layers are resolved more sharply, and continued mesh refinement expands

the DG coverage. Regions of high DG coverage coincide with areas of higher solution regularity.

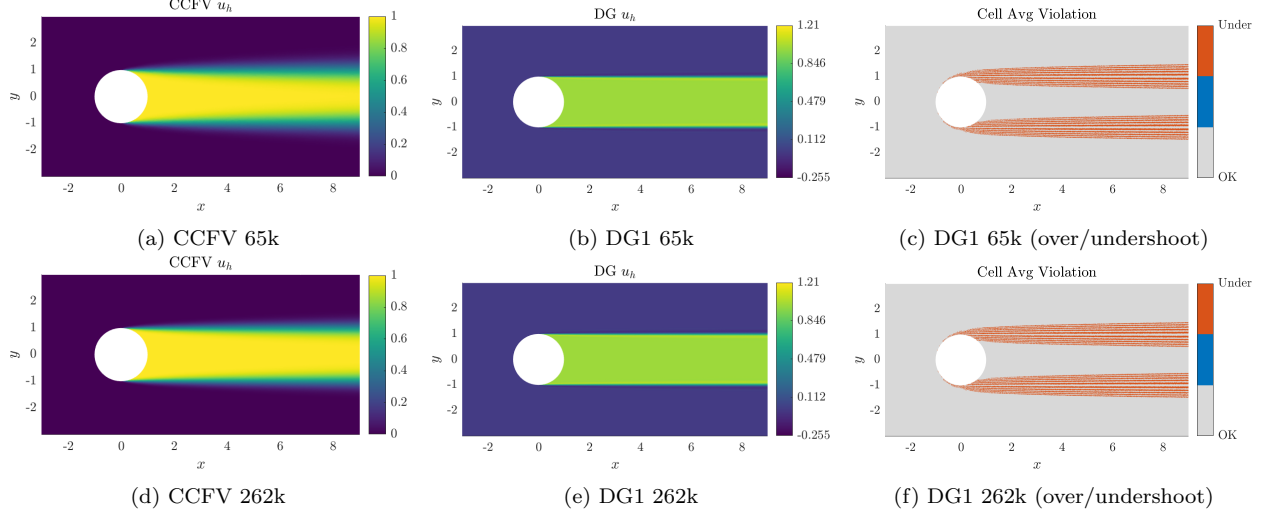


Figure 13: Baseline tests for the Hemker problem.

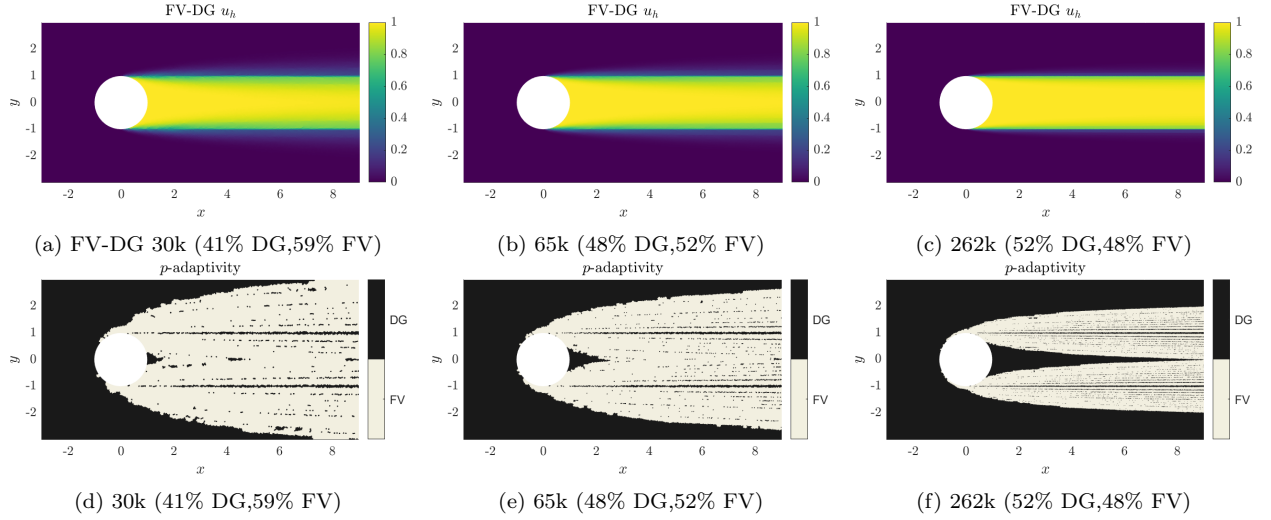


Figure 14: FV-DG1 results for the Hemker problem with $\delta = 10^{-9}$.

Figs. 15 and 16 visualizes the various approximations from a different vantage point. A view of the solution cross section along $x = 2$ is given in Fig. 16, along with magnification. The overshoot and undershoot from the DG scheme is more noticeable. The solutions from CCFV and FV-DG are more diffusive, but bound-preserving. Between the CCFV and FV-DG solutions, the latter is less diffusive.

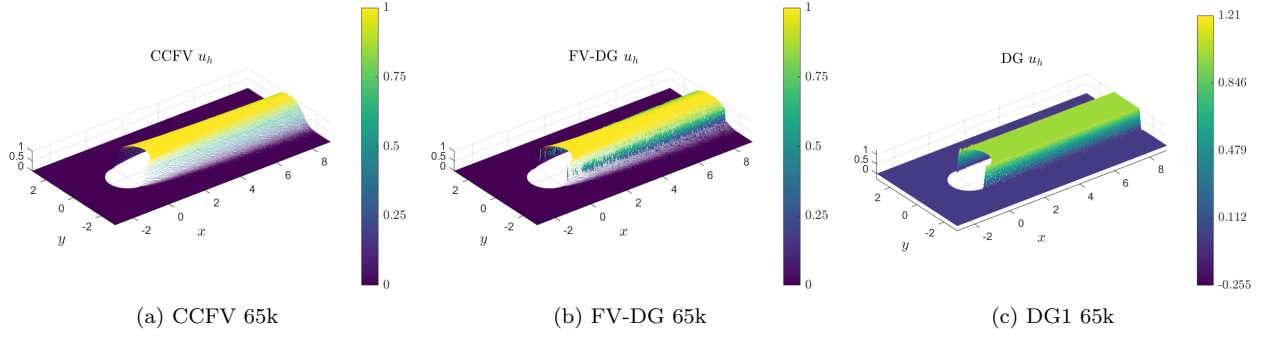


Figure 15: 3D visualization the Hemker problem with CCFV, DG1, and FV-DG ($\delta = 10^{-9}$).

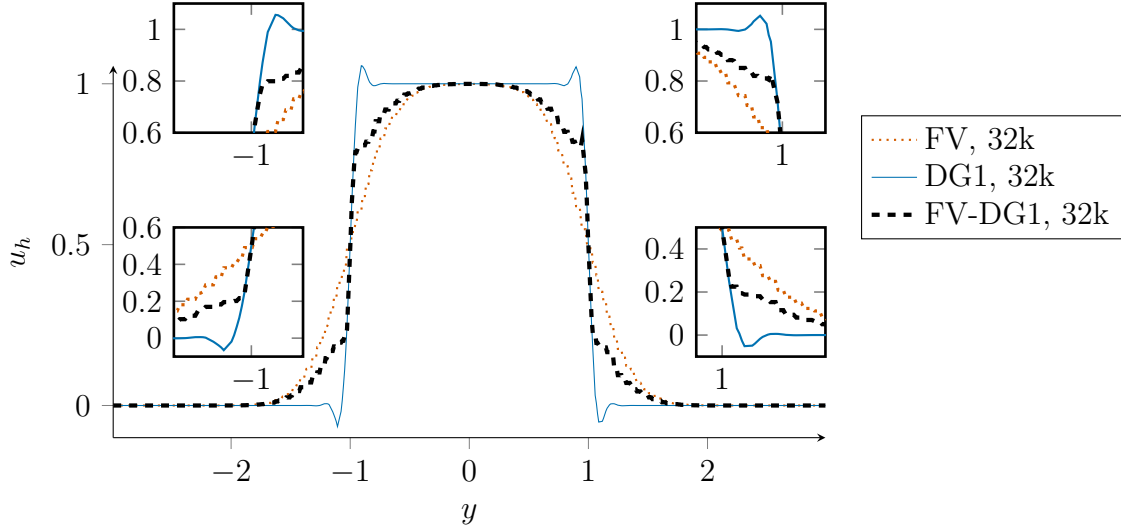


Figure 16: Profile along $x = 2$ for the Hemker problem.

11. Iteration counts, measure for the size of the spurious oscillations

Here we detail some metrics for the adaptive technique from Section 5. In [18], a metric for quantifying spurious oscillations is proposed; we employ that measure here:

$$\text{osc}(u_h) = \frac{1}{|\mathcal{T}_h|} \left[\sum_{E \in \mathcal{T}_h} \max\{0, -u_{\max} + \max_{(x,y) \in E} u_h(x,y)\} + \max\{0, u_{\min} - \max_{(x,y) \in E} u_h(x,y)\} \right],$$

where u_h is understood to be the solution to (2), and the exact solution to (1) satisfies $u_{\min} \leq u \leq u_{\max}$. As noted in Section 7.1, $\text{osc}(u_h)$ reports only deviations beyond $\min u$ and $\max u$; in-range oscillations are not captured. A value of $\text{osc}(u_h)$ near zero indicates that the approximation does not exhibit large extremal oscillations.

Table 1 reports the results. The **iters** column gives the total count of adaptation cycles (activations of the procedure in Section 5). Across all cases, we fix the mesh. For the full DG scheme, $\text{osc}(u_h)$ stays on the order of 10^{-1} and shows little reduction under mesh

refinement. The adaptive strategy of Section 5 yields a bound-preserving FV-DG scheme up to the prescribed tolerance δ . As δ is decreased, the measured oscillation magnitude drops accordingly, with $\text{osc}(u_h) \approx \delta$.

In the **iters** column, we see that the adaptation cycles required to meet a certain tolerance remains under 10. As one might expect, as the tolerance δ is made more strict, addition iterations are required. A more aggressive patch neighborhood $\mathcal{N}$ can also reduce the number of iterations. Adaptation cycles contribute negligibly to the total runtime, whereas the linear system solve and Voronoi remeshing are the primary bottlenecks.

Table 1: Summary of metrics for each test case.

Problem	Case	$ \mathcal{T}_h $	osc_{avg}	iters	δ	
Triple layer (Section 7)	DG1	262k	$1.280 \cdot 10^{-1}$	—	—	—
	FV-DG1	262k	$2.128 \cdot 10^{-6}$	4	10^{-6}	
	FV-DG1	262k	$1.102 \cdot 10^{-9}$	6	10^{-9}	
	FV-DG1	262k	$3.120 \cdot 10^{-13}$	7	10^{-13}	
L-Shaped (Section 8)	DG1	262k	$1.332 \cdot 10^{-1}$	—	—	—
	FV-DG1	262k	$1.923 \cdot 10^{-11}$	3	10^{-6}	
	FV-DG1	262k	$1.923 \cdot 10^{-11}$	3	10^{-9}	
	FV-DG1	262k	$1.817 \cdot 10^{-13}$	4	10^{-13}	
Internal layer (Section 9)	DG1	262k		—	—	—
	FV-DG1	262k	$2.832 \cdot 10^{-6}$	2	10^{-6}	
	FV-DG1	262k	$9.952 \cdot 10^{-9}$	3	10^{-9}	
	FV-DG1	262k	$2.256 \cdot 10^{-13}$	4	10^{-13}	
Hemker (Section 10)	DG1	262k	$7.040 \cdot 10^{-2}$	—	—	—
	FV-DG1	262k	$3.195 \cdot 10^{-6}$	5	10^{-6}	
	FV-DG1	262k	$1.330 \cdot 10^{-9}$	6	10^{-9}	
	FV-DG1	262k	$1.162 \cdot 10^{-13}$	8	10^{-13}	

12. Conclusion

In this paper we solve the steady convection–diffusion equation using a multinumetrics coupling of cell-centered finite volume and discontinuous Galerkin methods. We introduce a novel adaptive partitioning strategy that automatically selects FV and DG subdomains: whenever the cell average on an element falls outside the admissible range, that element and a small neighborhood are reassigned to the FV partition. The procedure integrates naturally with standard hp -adaptivity and is iterated until all cell averages satisfy the bounds up to a user-specified tolerance. Termination is guaranteed because the CCFV discretization is monotone. Since the DG and FV methods are not modified in any way during adaption,

standard stability and error analysis ensure the scheme is grounded in a rigorous framework. Numerical results on convection-dominated benchmarks demonstrate the effectiveness of the approach.

Future work will investigate alternative selection of the patch neighborhoods, slope limiters for Voronoi meshes, and algorithmic changes to Variants 1–2. Proving early termination of Variant 1 remains an open question.

Acknowledgments

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