ON LOCAL SUPER-PENALIZATION OF INTERIOR PENALTY DISCONTINUOUS GALERKIN METHODS

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Abstract. We prove in an abstract setting that standard (continuous) Galerkin finite element approximations are the limit of interior penalty discontinuous Galerkin approximations as the penalty parameter tends to infinity. We apply this result to equations of non-negative characteristic form and the non-linear, time dependent system of incompressible miscible displacement. Moreover, we investigate varying the penalty parameter on only a subset of a triangulation and the effects of local super-penalization on the stability of the method, resulting in a partly continuous, partly discontinuous method in the limit. An iterative automatic procedure is also proposed for the determination of the continuous region of the domain without loss of stability of the method.

Key words. discontinuous Galerkin methods, finite elements, interior penalty.

1. Introduction

The discontinuous Galerkin (dG) finite element method has become widely used in recent years for a variety of problems as it possesses several desirable qualities, such as: good stability properties due to the natural incorporation of upwinding techniques; flexible mesh design as hanging nodes and irregular meshes are admissible; and relatively easy implementation of hp-adaptive algorithms. These properties however come with the drawback of an increased number of degrees of freedom compared to a standard conforming method. For instance, when using an axi-parallel quadrilateral mesh in two dimensions with piecewise bilinear elements for which the standard continuous Galerkin (cG) finite element method has approximately n degrees of freedom (depending on boundary conditions) the dG method on the same mesh has approximately 4n degrees of freedom.

For advection-dominated advection-diffusion-reaction equations the standard cG method exhibits poor stability properties and non physical oscillations may pollute the approximation globally. Discontinuous Galerkin methods have generally better stability properties. In the case of interior penalty dG method, for instance, stability in the upwind direction has been shown in the inf-sup sense, e.g., in [1, 8], generalizing ideas from [21], where purely hyperbolic problems were considered.

Conceptually, somewhere between the standard cG and interior penalty dG methods lies the continuous-discontinuous Galerkin (cdG) finite element method [12], whereby one seeks a Galerkin solution on a finite element space $V_{\rm cdG}$ with $V_{\rm cG} \subset V_{\rm cdG} \subset V_{\rm dG}$, where $V_{\rm cG}$ and $V_{\rm dG}$ are the standard cG and dG finite element spaces. In the context of problems with layers or sharp fronts, continuous elements can be used away from the layers/fronts and discontinuous elements (accommodating appropriate upwinding) can be used in the region where the layers/fronts are present. This idea has been studied previously in the context of problems with layers by Dawson and Proft [16] using transmission conditions between regions where

different spaces are used. Cangiani, Georgoulis and Jensen [12] and Devloo, Forti and Gomez [17] have previously compared the cdG finite element method with alternative methods for advection-diffusion equations.

The control of discontinuities across element interfaces in the dG framework can be exercised by introducing and/or tuning the, so-called, jump penalization parameters. Using excessive penalization within a dG approximation will be referred to as the *super penalty method*. It is natural to expect that as the penalty parameter is increased the interelement jumps in the numerical approximation decrease. It has been shown by Larson and Niklasson [22] for stationary linear elliptic problems (using the interior penalty method) and by Burman, Quarteroni and Stamm [9] for stationary hyperbolic problems (penalising the jumps of the approximation for discontinuous elements and the jumps in the gradient of the approximation for continuous elements) that the dG approximation converges to the cG approximation as the jump penalization parameter tends to infinity.

In this work, our aim is twofold. Firstly, we present an alternative proof of the convergence of dG methods to cG methods, using a far more general framework covering the cases considered by [9, 22] and also non-linear and time dependent problems. Moreover, we show that super-penalization procedures can be localized to designated element faces, thereby arriving to partly continuous, partly discontinuous finite element methods. As particular examples we consider the limits of the interior penalty dG method for PDEs with non-negative characteristic form [20] and the mixed Raviart-Thomas-dG method for the miscible displacement system presented in [4].

Secondly, we continue the numerical investigations of [12] in the context of blending locally continuous and discontinuous methods. In particular, we investigate to what degree numerical oscillations appear as local super-penalization is applied. The aim, of course, is to find the extent to which degrees of freedom can be removed by using locally continuous finite element spaces without affecting the extra stability offered by dG methods. To this end, we consider an advection-dominated advection-diffusion problem containing boundary layer behaviour, where the continuous and the discontinuous regions of the finite element solution are tuned manually. A second example investigates the use of an iterative automatic procedure for the determination of the continuous region of the domain by local super-penalization without loss of stability of the method. The procedure is applied to the problem of incompressible miscible displacement.

This work is organized as followed. After introducing notation in Section 1.1 an abstract discussion of the limit of penalty methods is given in Section 2. We then show how this framework can be applied to equations of non-negative characteristic form in Section 3 and to the non-linear equations of incompressible miscible displacement in Section 4. Finally, Section 5 contains a number of numerical experiments and discussion of an iterative automatic procedure for determining the continuous regions of the approximation.

1.1. Notation. Let $\Omega \subset \mathbb{R}^d$ be a bounded, open polygonal domain. We denote by \mathcal{T}_h a subdivision of Ω into open non-overlapping d-simplices E. The diameter of $E \in \mathcal{T}_h$ is denoted by h_E . Let also $\mathcal{E}_h := \bigcup_{E \in \mathcal{T}_h} \partial E$ be the skeleton of the mesh \mathcal{T}_h , while $\mathcal{E}_h^o := \mathcal{E}_h \setminus \partial \Omega$. Finally, let Γ denote the set of elemental boundary faces, i.e., those which lie in $\partial \Omega$.

For $e \in \mathcal{E}_h^o$, with $e = \bar{E}^+ \cap \bar{E}^-$ for E^+ , $E^- \in \mathcal{T}_h$, we define $h_e := \min(h_{E^-}, h_{E^+})$. Given a generic scalar field $\nu : \Omega \to \mathbb{R}$ that may be discontinuous across e, we set $\nu^{\pm} := \nu|_{E^{\pm}}$, the interior trace on E^{\pm} . Similarly, for a generic vector field

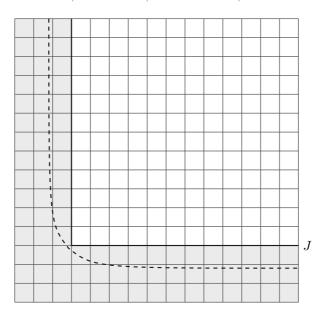


FIGURE 1.1. An example of a cdG decomposition. The region \mathcal{T}_{dG} is marked by gray elements, while \mathcal{T}_{cG} consists of the white elements. The interface between \mathcal{T}_{cG} and \mathcal{T}_{dG} is J. Such a decomposition could be used if left and below of the dashed line the solution exhibits a boundary layer.

 $\tau: \Omega \to \mathbb{R}^d$ possibly discontinuous across e we set $\tau^{\pm} := \tau|_{E^{\pm}}$. We define the average and jump for the generic scalar as

$$\{\!\!\{\nu\}\!\!\} := \frac{1}{2}(\nu^+ + \nu^-), \qquad [\![\nu]\!] := \nu^+ n^+ + \nu^- n^-, \qquad \text{on } e \in \mathcal{E}_h^o,$$

and for the generic vector field as

$$\{\!\!\{ m{ au} \}\!\!\} := rac{1}{2} (m{ au}^+ + m{ au}^-), \qquad [\![m{ au}]\!] := m{ au}^+ \cdot n^+ + m{ au}^- \cdot n^-, \qquad \text{on } e \in \mathcal{E}_h^o,$$

where n^{\pm} is the outward pointing normal from E^{\pm} on e. For $e \in \Gamma$ the definitions become

$$\{\!\!\{\nu\}\!\!\} := \nu, \qquad [\![\nu]\!] := \nu n, \qquad \{\!\!\{\boldsymbol{\tau}\}\!\!\} := \boldsymbol{\tau}, \qquad \text{on } e \in \Gamma.$$

Given a vector b, we denote the inflow and outflow boundaries of Ω by

$$\begin{split} \partial^{\text{in}}\Omega &:= \Gamma^{\text{in}} &:= \{ \boldsymbol{x} \in \partial\Omega : b \cdot n \leq 0 \}, \\ \partial^{\text{out}}\Omega &:= \Gamma^{\text{out}} &:= \{ \boldsymbol{x} \in \partial\Omega : b \cdot n > 0 \}, \end{split}$$

and for an element

$$\begin{split} \partial^{\text{in}} E & := \; \{ \boldsymbol{x} \in \partial E : b \cdot n \leq 0 \}, \\ \partial^{\text{out}} E & := \; \{ \boldsymbol{x} \in \partial E : b \cdot n > 0 \}. \end{split}$$

We denote the trace of a function ν on an edge by $\nu^{\rm in}$ (resp. $\nu^{\rm out}$) on the side of the edge where $b \cdot n \leq 0$ (resp. $b \cdot n > 0$). We construct the mesh so that the sign of $b \cdot n$ is the same for every $x \in e$.

For the cdG method, we will require the following additional notation. We identify a decomposition of our triangulation \mathcal{T}_h into two disjoint triangulations $\mathcal{T}_{\mathrm{dG}}$ and $\mathcal{T}_{\mathrm{cG}} := \mathcal{T}_h \setminus \mathcal{T}_{\mathrm{dG}}$, upon which continuous and discontinuous elements will be applied respectively, henceforth referred to as the continuous and discontinuous

regions of the triangulation. Define $J := \overline{\mathcal{T}}_{cG} \cap \overline{\mathcal{T}}_{dG}$ and let \mathcal{E}_{dG} be the skeleton of \mathcal{T}_{dG} and $\mathcal{E}_{cG} := \mathcal{E}_h \setminus \mathcal{E}_{dG}$. Note that with this definition the faces in J are part of the discontinuous skeleton \mathcal{E}_{dG} only. Define $\Gamma_{cG} := \mathcal{T}_{cG} \cap \Gamma$, the set of boundary faces of the continuous region, and similarly $\Gamma_{dG} := \mathcal{T}_{dG} \cap \Gamma$.

Finally, we will denote by ∇_h the elementwise divergence operator.

2. An abstract discussion

The operators defining discontinuous Galerkin formulations are typically made of two parts, herein denoted by \mathcal{B} and \mathcal{S} , respectively: \mathcal{B} is there to ensure consistency, while \mathcal{S} weakly enforces conformity thorough penalty and is transparent to the underlying continuous space. Here we show, in an abstract setting, under which conditions a discontinuous Galerkin solution will converge to a conforming solution under super-penalisation.

Consider a (possibly non-linear) operator $\mathcal{B}: W \times W \to \mathbb{R}$ where W is a finite dimensional vector space with norm $\|\cdot\|_W$. Suppose there exists a decomposition of W such that $V \oplus X = W$ for $V, X \subset W$. In particular this means we can write any $w \in W$ uniquely as w = v + x for some $v \in V$ and $x \in X$.

Assume that \mathcal{B} is coercive, i.e., there exists $\Lambda_W > 0$ (typically independent of the dimension of W), such that

(2.1)
$$\mathcal{B}(w,w) \ge \Lambda_W \|w\|_W^2 \qquad \forall \ w \in W.$$

Consider an additional linear operator $S: W \times W \to \mathbb{R}$, whose support is restricted to $X \times X$ in the sense that

(2.2)
$$S(v, \hat{v}) = 0 \qquad \forall \ v, \hat{v} \in V$$

and

(2.3)
$$S(v,x) = S(x,v) = 0 \quad \forall v \in V, x \in X.$$

We require coercivity on X, i.e., there exists $\Lambda_X > 0$ such that for all $x \in X$

$$\mathcal{S}(x,x) > \Lambda_X \|x\|_Y^2,$$

where $||x||_X : W \times W \to \mathbb{R}$ is a semi norm, which is a norm on the restriction to X. In view of (2.2) this gives $S(w, w) \geq \Lambda_X ||w||_X^2 = \Lambda_X ||x||_X^2$. We construct a further operator

$$\mathcal{B}_{\sigma} := \mathcal{B} + \sigma \mathcal{S}$$

where $0 \le \sigma \in \mathbb{R}$, and call this the *super penalised* bilinear form.

Let ℓ be an element of the dual space W^* of W, independent of σ . Then choose $w_{\sigma} \in W$ such that

(2.6)
$$\mathcal{B}_{\sigma}(w_{\sigma}, w) = \ell(w) \quad \forall \ w \in W.$$

Also choose $v_h \in V$ such that

(2.7)
$$\mathcal{B}(v_h, v) = \ell(v) \qquad \forall \ v \in V.$$

Observe that for all $\sigma \in \mathbb{R}$

(2.8)
$$\mathcal{B}_{\sigma}(v_h, v) = \mathcal{B}(v_h, v) = \ell(v) \quad \forall \ v \in V$$

using (2.2). Now with (2.1), (2.4) and (2.6) we have

$$\begin{split} \Lambda_W \|w_\sigma\|_W^2 + \sigma \Lambda_X \|w_\sigma\|_X^2 &\leq \mathcal{B}(w_\sigma, w_\sigma) + \sigma \mathcal{S}(w_\sigma, w_\sigma) \\ &= \mathcal{B}_\sigma(w_\sigma, w_\sigma) \\ &= \ell(w_\sigma) \\ &\leq \|\ell\|_{W^*} \|w_\sigma\|_W. \end{split}$$

Using Young's inequality we see

(2.9)
$$\frac{\Lambda_W^2}{\sigma} \|w_{\sigma}\|_W^2 + 2\Lambda_W \Lambda_X \|w_{\sigma}\|_X^2 \le \frac{1}{\sigma} \|\ell\|_{W^*}^2.$$

Each of Λ_W , Λ_X and $\|\ell\|_{W^*}$ are independent of σ . We write $w_{\sigma} = v_{\sigma} + x_{\sigma}$, the unique decomposition with $v_{\sigma} \in V$ and $x_{\sigma} \in X$. From (2.9) we see

(2.10)
$$\lim_{\sigma \to \infty} ||v_{\sigma} + x_{\sigma}||_{X} = \lim_{\sigma \to \infty} ||x_{\sigma}||_{X} = 0.$$

Therefore $x_{\sigma} \to 0$ as $\sigma \to \infty$.

Now assume that \mathcal{B} is continuous in the first argument in the following sense: If $\lim_{i\to\infty} w_i = w \in W$ then

(2.11)
$$\lim_{i \to \infty} \mathcal{B}(w_i, v) = \mathcal{B}(w, v) \qquad \forall \ v \in V.$$

Suppose $w_{\sigma} \nrightarrow v_h$ as $\sigma \to \infty$. Then there exists $\varepsilon > 0$ such that there is some sequence $\{w_{\sigma(i)}\}_i$ with $\sigma(i) \to \infty$ as $i \to \infty$ satisfying

$$(2.12) ||w_{\sigma(i)} - v_h||_W > \varepsilon \forall i \in \mathbb{N}.$$

Owing to (2.9) the sequence $\{w_{\sigma(i)}\}_i$ is a bounded subset of W. Then by the Heine-Borel theorem there exists a convergent subsequence, also denoted $\{w_{\sigma(i)}\}_i$, such that

(2.13)
$$\tilde{w} = \lim_{i \to \infty} w_{\sigma(i)}.$$

Considering (2.10) we know that $\tilde{w} \in V$. We have that for all $v \in V$

$$\mathcal{B}(\tilde{w}, v) = \mathcal{B}\left(\lim_{i \to \infty} w_{\sigma(i)}, v\right)$$

$$= \lim_{i \to \infty} \mathcal{B}(w_{\sigma(i)}, v) \qquad \text{by (2.11)}$$

$$= \lim_{i \to \infty} \mathcal{B}_{\sigma}(w_{\sigma(i)}, v) \qquad \text{by (2.3)}$$

$$= \lim_{i \to \infty} \ell(v) \qquad \text{by (2.6)}$$

$$= \ell(v).$$

Hence \tilde{w} satisfies (2.7) and by (2.13) we have

$$\lim_{i \to \infty} ||w_{\sigma(i)} - v_h||_W = 0.$$

This contradicts (2.12) and we conclude that all subsequences $\{w_{\sigma(i)}\}_i$ converge to v_h . Therefore

(2.14)
$$\lim_{\sigma \to \infty} (w_{\sigma} - v_h) = 0.$$

We finally remark on the potential loss of stability due to super-penalization. It can be seen from (2.9) that as $x_{\sigma} \to 0$ when $\sigma \to \infty$ the coercivity of \mathcal{B}_{σ} is increasingly compromised, which can lead to loss of stability and reduction on the rate of convergence in various settings. A proof of stability in the convection-dominated linear case is given in [11], where dG elements are used in boundary and internal layers and cG elements elsewhere.

3. Equations of Non-Negative Characteristic Form

We now examine the diffusion-advection-reaction equation (see [20])

(3.1)
$$-\nabla \cdot (\mathbb{A}(\boldsymbol{x})\nabla u) + b(\boldsymbol{x}) \cdot \nabla u + c(\boldsymbol{x})u = f(\boldsymbol{x}) \quad \text{in } \Omega,$$

$$u = 0 \quad \text{on } \partial\Omega$$

with b a \mathbb{R}^d valued function whose entries are Lipschitz continuous on $\overline{\Omega}$, $c \in L^{\infty}(\Omega)$ and $f \in L^2(\Omega)$ real valued functions. The diffusion coefficient \mathbb{A} is a $d \times d$ symmetric matrix with entries being bounded, piecewise continuous real-valued functions defined on $\overline{\Omega}$, with

$$\zeta^{\top} \mathbb{A} \zeta > 0 \qquad \forall \zeta \in \mathbb{R}^d, \text{ a.e. in } \overline{\Omega}.$$

Furthermore, we assume that $c - \frac{1}{2}\nabla \cdot b \ge 0$. Under these conditions, (3.1) is named a partial differential equation with non-negative characteristic form.

We define the dG space to be

$$(3.2) V_{\mathrm{dG}} := \{ v \in L^2(\Omega) : \forall E \in \mathcal{T}_h, v|_E \in \mathbb{P}^r \}.$$

We define the cdG space to be

$$(3.3) V_{\text{cdG}} := \{ v \in L^2(\Omega) : \forall E \in \mathcal{T}_h, v|_E \in \mathbb{P}^r, v|_{\Gamma_{\text{cG}}} = 0, v|_{\overline{\mathcal{T}}_{\text{cG}}} \in C(\overline{\mathcal{T}}_{\text{cG}}) \}$$

where \mathbb{P}^r is the space of polynomials of degree at most r supported on E. Finally we let V_{\perp} be a subspace of V_{dG} such that $V_{\mathrm{dG}} = V_{\mathrm{cdG}} \oplus V_{\perp}$ (corresponding to $W := V \oplus X$ in the notation of Section 2). Note that the standard continuous space is obtained by setting $\mathcal{T}_h = \mathcal{T}_{\mathrm{cG}}$.

Define $\mathcal{B}: V_{\mathrm{dG}} \times V_{\mathrm{dG}} \to \mathbb{R}$, the bilinear form for the interior penalty family of methods with $\vartheta \in \{-1,0,1\}$ for (3.1), by

(3.4)
$$\mathcal{B}(w,\hat{w}) := \mathcal{B}_d(w,\hat{w}) + \mathcal{B}_{ar}(w,\hat{w})$$

with

(3.5)
$$\mathcal{B}_{d}(w, \hat{w}) := \sum_{E \in \mathcal{T}_{h}} \int_{E} \mathbb{A} \nabla_{h} w \cdot \nabla_{h} \hat{w} \, d\boldsymbol{x} + \sum_{e \in \mathcal{E}_{h}} \int_{e} m \llbracket w \rrbracket \cdot \llbracket \hat{w} \rrbracket \, ds - \sum_{e \in \mathcal{E}_{h}} \int_{e} \left(\{\!\!\{ \mathbb{A} \nabla_{h} w \}\!\!\} \cdot \llbracket \hat{w} \rrbracket - \vartheta \{\!\!\{ \mathbb{A} \nabla_{h} \hat{w} \}\!\!\} \cdot \llbracket w \rrbracket \right) ds$$

and

(3.6)
$$\mathcal{B}_{ar}(w, \hat{w}) := \sum_{E \in \mathcal{T}_h} \int_E (b \cdot \nabla_h w) \hat{w} + cw \hat{w} \, \mathrm{d}\boldsymbol{x} \\ - \sum_{e \in \mathcal{E}_e^o} \int_e b \cdot [\![w]\!] \hat{w}^{\text{out}} \, \mathrm{d}s - \sum_{e \in \Gamma^{\text{in}}} \int_e (b \cdot n) w \hat{w} \, \mathrm{d}s.$$

We define $m := C_p\{\!\{\overline{\mathbb{A}}r^2\}\!\}/h_e$, $\overline{\mathbb{A}} := \||\sqrt{\mathbb{A}}|_2\|_{L^{\infty}(E)}$, with $|\cdot|_2$ denoting the matrix-2-norm, and $C_p(\vartheta) \geq 0$ fixed for a given ϑ . The linear form is given by

(3.7)
$$\ell(w) := \sum_{E \in \mathcal{T}_b} \int_E f w \, \mathrm{d} \boldsymbol{x}.$$

For $e \in \mathcal{E}_{cG}$ we have the additional term $\mathcal{S}: V_{dG} \times V_{dG} \to \mathbb{R}$ penalising the jumps where

(3.8)
$$\mathcal{S}(w, \hat{w}) := \sum_{e \in \mathcal{E}_{cG}} \int_{e} M[\![w]\!] \cdot [\![\hat{w}]\!] ds$$

and

$$M := \left(C_{ar} + C_d \frac{\{\!\!\{ \overline{\mathbb{A}} r^2 \}\!\!\}}{h_e} \right)$$

with C_{ar} and C_d fixed constants independent of σ . Then we define $\mathcal{B}_{\sigma}(w, \hat{w}) := \mathcal{B}(w, \hat{w}) + \sigma \mathcal{S}(w, \hat{w})$.

Observe that if we take $C_{ar} = C_d = 0$ (or $\sigma = 0$) we recover the usual interior penalty method. If we take $C_p = C_d = C_{ar} = 0$ and A = 0 we have the standard (unpenalised) bilinear form for the purely hyperbolic equation (assuming of course that we adjust the boundary conditions appropriately). Taking $C_d = 0$ and $C_{ar} \neq 0$ when A = 0 gives the method proposed in [7].

All functions in $V_{\rm cdG}$ are continuous on edges in $\mathcal{E}_{\rm cG}$ (recall that by definition edges in J are not included in $\mathcal{E}_{\rm cG}$). Therefore conditions (2.2) and (2.3) are satisfied for this \mathcal{S} . That is, for any $v, \hat{v} \in V_{\rm cdG}$ and $x \in V_{\perp}$

(3.9)
$$\mathcal{S}(v,\hat{v}) = \mathcal{S}(v,x) = \mathcal{S}(x,v) = 0.$$

We define the following norm for all $w \in V_{dG}$:

(3.10)
$$||w||_{\mathrm{dG}}^2 := \sum_{E \in \mathcal{T}_h} ||\sqrt{\mathbb{A}}\nabla_h w||_{L^2(E)}^2 + ||c_0 w||_{L^2(\Omega)}^2$$

$$+ \sum_{e \in \mathcal{E}_h} \frac{1}{2} |||b \cdot n|^{1/2} [\![w]\!]||_{L^2(e)}^2 + \sum_{e \in \mathcal{E}_h} ||\sqrt{m}[\![w]\!]||_{L^2(e)}^2$$

where $c_0 := \sqrt{c - 1/2\nabla_h \cdot b}$. We also define for $w \in V_{\mathrm{dG}}$

(3.11)
$$|w|_{\mathcal{S}}^{2} := \sum_{e \in \mathcal{E}_{cG}} \|\sqrt{M} [w]\|_{L^{2}(e)}^{2}.$$

Notice that $|\cdot|_{\mathcal{S}}$ is a semi-norm on V_{dG} but a norm on V_{\perp} ; it corresponds to the norm $\|\cdot\|_X$ of Section 2. To make this distinction clear we will write $\|x\|_{\mathcal{S}}$ for $x\in V_{\perp}$.

Lemma 3.12. If C_p is sufficiently large when $\vartheta = -1$ then \mathcal{B} is coercive on V_{dG} , i.e., for all $w \in V_{dG}$

(3.13)
$$\mathcal{B}(w,w) \ge \Lambda_{cc} \|w\|_{dG}^2$$

with $\Lambda_W = 1$ when $\vartheta = 1$ and $\Lambda_W = 1/2$ when $\vartheta = -1$.

From the definition it is clear that \mathcal{S} is coercive with constant one on V_{\perp} , i.e., for all $x \in V_{\parallel}$

(3.14)
$$S(x,x) = ||x||_{S}^{2}.$$

Definition 3.15. Define a dG approximation to (3.1) as $w_{\sigma} \in V_{dG}$ satisfying

(3.16)
$$\mathcal{B}_{\sigma}(w_{\sigma}, w) = \ell(w) \qquad \forall w \in V_{dG}.$$

Definition 3.17. Define a cdG approximation to (3.1) as $v_h \in V_{cdG}$ satisfying

(3.18)
$$\mathcal{B}_{\sigma}(v_h, v) = \ell(v) \qquad v \in V_{cdG}.$$

Using (3.9) we see that v_h also satisfies $\mathcal{B}(v_h, v) = \ell(v)$ for all $v \in V_{\text{cdG}}$.

Theorem 3.19. The dG finite element approximation w_{σ} converges to the cdG finite element approximation v_h as $\sigma \to \infty$, i.e.,

$$\lim_{\sigma \to \infty} (w_{\sigma} - v_h) = 0.$$

Proof. Following the argument of Section 2 we use Lemma 3.12 and (3.14) and note that (2.11) is satisfied as linear operators in finite-dimensional vector spaces are continuous.

4. Incompressible Miscible Displacement

We consider the problem of finding the numerical solution to the coupled equations for the pressure $p = p(t, \mathbf{x})$, Darcy velocity $u = u(t, \mathbf{x})$ and concentration $c = c(t, \mathbf{x})$ of one incompressible fluid in a porous medium being displaced by another. We consider the miscible case where both fluids are in the same phase.

Consider the domain $\Omega_T := (0, T) \times \Omega$. The equations for the miscible displacement are given by (e.g., [5, 6])

(4.1)
$$\varphi \frac{\partial c}{\partial t} + u \cdot \nabla c - \nabla \cdot (\mathbb{D}(u)\nabla c) + cq^{I} = \hat{c}q^{I},$$

$$(4.2) \nabla \cdot u = q^I - q^P,$$

(4.3)
$$u = -\frac{\mathbb{K}}{\mu(c)} \left(\nabla \cdot p - \rho(c)g \right)$$

with the boundary conditions on $\partial\Omega_T := (0,T) \times \partial\Omega$ given by

$$(4.4) u \cdot n = 0$$

$$(4.5) \qquad (\mathbb{D}(u)\nabla c) \cdot n = 0,$$

and the initial conditions

$$(4.6) c(0,\cdot) = c_0.$$

We denote by: $\varphi(\boldsymbol{x})$ the porosity of the medium; $q^I \geq 0$ and $q^P \geq 0$ the pressure at injected (source) and production (sink) wells; $\mathbb{K}(\boldsymbol{x})$ the absolute permeability of the medium; $\mu(c)$ the viscosity of the fluid mixture; $\rho(c)$ the density of the fluid mixture; g the constant vector of gravity; $\mathbb{D}(u, \boldsymbol{x})$ the diffusion-dispersion coefficient; \hat{c} the injected concentration; and c_0 the initial concentration, which we assume for simplicity to be 0. We define $a^{-1}(c) := \mathbb{K}^{-1}\mu$. The coupling is nonlinear through the coefficients $\mathbb{D}(u, \boldsymbol{x})$, $\mu(c)$ and the advection term. We make the following common specific choice for the diffusion dispersion tensor [15, 19, 24]:

(4.7)
$$\mathbb{D}(u, \mathbf{x}) = \varphi \left(d_m \mathbb{I} + |u| d_l \mathbb{E}(u) + |u| d_t (\mathbb{I} - \mathbb{E}(u)) \right)$$

where $\mathbb{E}(u) = uu^{\top}/|u|^2$ and \mathbb{I} is the identity matrix. We specify that the molecular, longitudinal and transverse diffusion coefficients d_m , d_l and d_t are positive real numbers.

We solve for the pressure and velocity using a Raviart-Thomas (RT) procedure [18, 23] and for the concentration using a cdG method. We refer to the whole scheme as a RT-cdG method. For $k \geq 0$ we define

$$U := \{ v \in (L^2(\Omega))^2 : v|_E \in (\mathbb{P}_k(E))^2 + \boldsymbol{x} \mathbb{P}_k(E) \ \forall E \in \mathcal{T}_h, \\ v \cdot n \text{ continuous on } e \in \mathcal{E}_h^o \}.$$

To avoid confusion for the pressure terms we define the space $P := V_{\rm dG}$ where $V_{\rm dG}$ is defined in (3.2). Then the velocity and pressure are approximated in $U \times P$. To simplify the presentation we use the same mesh \mathcal{T}_h to solve for u, p and c numerically at each time step and there is no refinement of the mesh or polynomial degree. However $\mathcal{T}_{\rm cG}$ and $\mathcal{T}_{\rm dG}$ are not fixed so the cdG space used to approximate c will vary with time. We define the time dependent cdG space by

$$V_{\mathrm{cdG}}^{j} := \{ v \in L^{2}(\Omega) : \forall E \in \mathcal{T}_{h}, v|_{E} \in \mathbb{P}^{r}, v|_{\Gamma_{\mathrm{cG}}^{j}} = 0, v|_{\overline{\mathcal{T}}_{\mathrm{cG}}^{j}} \in C(\overline{\mathcal{T}}_{\mathrm{cG}}^{j}) \}$$

where \mathcal{T}_{cG}^{j} and Γ_{cG}^{j} are the \mathcal{T}_{cG} region and external boundary of \mathcal{T}_{cG}^{j} at time t_{j} . As we assert that no change to the shape of the mesh occurs in time we define the time dependent dG space as in (3.2). Then we may set $V_{dG}^{j} := V_{cdG}^{j} \oplus V_{\perp}^{j}$ corresponding to the decomposition $W = V \oplus X$ of Section 2. Note that the degree k is the same for U and P but need not be equal to r, the degree of the polynomials used to approximate concentration.

Let $0 = t_0 < t_1 < \ldots < t_N = T$ be a partition of the time interval (0,T). For simplicity we assume that each time step is of equal length and define $\Delta t := t_j - t_{j-1}$ and the backward Euler operator $\mathfrak{d}_{\mathfrak{t}} c_h^j := (\Delta t)^{-1} (c_h^j - c_h^{j-1})$ for $j = 1, 2, \ldots, N$. For the diffusion part of the concentration equation define the bilinear form

$$(4.8) \qquad \mathcal{B}_{d}(c_{h}^{j}, d_{h}^{j}; u_{h}^{j}) = \sum_{E \in \mathcal{T}_{h}} (\mathbb{D}(u_{h}^{j}) \nabla_{h} c_{h}^{j}, \nabla_{h} d_{h}^{j})_{E} + \sum_{e \in \mathcal{E}_{h}^{o}} m(\llbracket c_{h}^{j} \rrbracket, \llbracket d_{h}^{j} \rrbracket)_{e} \\ - \sum_{e \in \mathcal{E}_{h}^{o}} \left[(\llbracket c_{h}^{j} \rrbracket, \{\!\!\{ \mathbb{D}(u_{h}^{j}) \nabla_{h} d_{h}^{j} \}\!\!\})_{e} + (\llbracket d_{h}^{j} \rrbracket, \{\!\!\{ \mathbb{D}(u_{h}^{j}) \nabla_{h} c_{h}^{j} \}\!\!\})_{e} \right]$$

for all $d_h^j \in V_{\text{cdG}}^j$. The penalty parameter m is defined by [4]

$$m^2: \mathcal{E}_h \to \mathbb{R}, \quad \boldsymbol{x} \mapsto C_{\mathrm{pen}} \frac{\max\{n_{\mathcal{E}_h}^{\top} \mathbb{D}(u_h^{j,+}, \boldsymbol{x}) n_{\mathcal{E}_h}, n_{\mathcal{E}_h}^{\top} \mathbb{D}(u_h^{j,-}, \boldsymbol{x}) n_{\mathcal{E}_h}\}}{h}$$

and C_{pen} is chosen such that it is larger than

$$\sup \left\{ h \max \left\{ \frac{\|\nu_h\|_{\partial E}^2}{\|\nu_h\|_E^2}, \frac{\|D^{1/2}\nabla_h\nu_h\|_{\partial E}^2}{\|D^{1/2}\nabla_h\nu_h\|_E^2} \right\} : \nu_h \in \mathbb{P}^s, D \in [\mathbb{P}^s]^{d \times d}, E \text{ shape regular} \right\}.$$

The bilinear form for convection, production and injection is given by the nonstandard form

$$(4.9) \qquad \mathcal{B}_{cq}(c_h^j, d_h^j; u_h^j) \\ = \frac{1}{2} \sum_{E \in \mathcal{T}_h} \left[(u_h^j \cdot \nabla_h c_h^j, d_h^j)_E - (u_h^j c_h^j, \nabla_h d_h^j) + ((q^I + q^P)c_h^j, d_h^j)_E \right] \\ + \frac{1}{2} \sum_{e \in \mathcal{E}_l^o} (u_h^j \cdot \llbracket c_h^j \rrbracket, d_h^{j,*})_e$$

where $d_h^{j,*}$ is defined by

(4.10)
$$d_h^{j,*} = \begin{cases} d_h^{j,-} & \text{if } u_h^j \cdot n^+ > 0, \\ d_h^{j,+} & \text{if } u_h^j \cdot n^+ \le 0. \end{cases}$$

This formulation ensures that \mathcal{B}_{cq} is semi-definite regardless of the properties of u_h^j . We do not need to restrict sums over edges to cells in \mathcal{T}_{cG} as in this region elements of V_{cdG} are continuous. Also note that the dG method is a special case of the cdG method where $\mathcal{T}_{cG} = \emptyset$.

Define
$$\mathcal{B}(c_h^j, d_h^j; u_h^j) := \mathcal{B}_d(c_h^j, d_h^j; u_h^j) + \mathcal{B}_{cq}(c_h^j, d_h^j; u_h^j)$$
 and

$$\mathcal{S}(c_h^j, d_h^j) := \sum_{e \in \mathcal{E}_{cG}^j} \int_e M[\![c_h^j]\!] \cdot [\![d_h^j]\!] \, \mathrm{d}s$$

where

$$M := \left(C_d \frac{r^2}{h_e} \right).$$

Then for any $c_h^j, d_h^j \in V_{\text{cdG}}^j$ and $x^j \in V_{\perp}^j$

$$S(c_h^j, d_h^j) = S(c_h^j, x^j) = S(x^j, c_h^j) = 0.$$

We define the following norm for $u_h^j \in U$ and $c_h^j \in V_{\mathrm{dG}}^j$:

$$|||c_h^j|||^2 := \sum_{E \in \mathcal{T}_h} ||\sqrt{\mathbb{D}(u_h^j)} \nabla_h c_h^j||_{L^2(E)}^2 + \frac{1}{2} ||q_0 c_h^j||_{L^2(\Omega)}^2$$

$$+ \sum_{e \in \mathcal{E}_h^o} \frac{1}{2} |||u_h^j \cdot n|^{1/2} [\![c_h^j]\!]||_{L^2(e)}^2 + \sum_{e \in \mathcal{E}_h^o} ||\sqrt{m} [\![c_h^j]\!]||_{L^2(e)}^2$$

where $q_0 := \sqrt{q^I + q^P}$. For $c_h^j \in V_{\mathrm{dG}}^j$ define

(4.13)
$$|c_h^j|_{\mathcal{S}}^2 := \sum_{e \in \mathcal{E}_{rC}^j} \|\sqrt{M} [c_h^j]\|_{L^2(e)}.$$

Notice that (4.13) is a semi-norm on V_{dG}^{j} but a norm on V_{\perp}^{j} .

Lemma 4.14. If C_{pen} is chosen large enough then \mathcal{B} is coercive for all $c_h^j \in V_{dG}^j$ and $u_h^j \in U$, i.e.,

(4.15)
$$\mathcal{B}(c_h^j, c_h^j; u_h^j) \ge \Lambda_W ||c_h^j||^2.$$

Proof. Combine equations (4.3) and (4.6) from [4].

We have by construction that $\mathcal S$ is coercive with constant one on $V_\perp^j,$ i.e., for all $x_h^j \in V_\perp^j$

(4.16)
$$S(x_h^j, x_h^j) = ||x_h^j||_{\mathcal{S}}^2.$$

We discretise the time derivative with the backward Euler operator. Summing over each discrete time step gives

$$\begin{split} \sum_{j=1}^{N} (\varphi \mathfrak{d}_{\mathfrak{t}} c_{h}^{j}, c_{h}^{j}) &= \sum_{j=1}^{N} \frac{1}{\Delta t} (\varphi c_{h}^{j}, c_{h}^{j}) - \frac{1}{\Delta t} (\varphi c_{h}^{j-1}, c_{h}^{j}) \\ &\geq \sum_{j=1}^{N} \frac{1}{\Delta t} \| \varphi^{1/2} c_{h}^{j} \|_{L^{2}(\Omega)}^{2} - \frac{1}{2\Delta t} \left(\| \varphi^{1/2} c_{h}^{j-1} \|_{L^{2}(\Omega)}^{2} + \| \varphi^{1/2} c_{h}^{j} \|_{L^{2}(\Omega)}^{2} \right) \\ &= \frac{1}{2\Delta t} \left(\| \varphi^{1/2} c_{h}^{N} \|_{L^{2}(\Omega)}^{2} - \| \varphi^{1/2} c_{h}^{0} \|_{L^{2}(\Omega)}^{2} \right) \end{split}$$

where we have used Young's Inequality. We have assumed that the initial concentration is 0 and so $\|\varphi^{1/2}c_h^0\|_{L^2(\Omega)}^2=0$.

Definition 4.17. Define the RT-dG approximation

$$(u_h, p_h, c_\sigma) \in \prod_{i=1}^N U \times \prod_{j=1}^N P \times \prod_{i=1}^N V_{dG}^j$$

to (4.1)-(4.6) as that generated by the algorithm: For $1 \leq j \leq N$ and $c_{\sigma}^{j-1} \in V_{dG}^{j}$ find $(u_{h}^{j}, p_{h}^{j}, c_{\sigma}^{j}) \in U \times P \times V_{dG}^{j}$ such that

$$(4.18) (\nabla_h \cdot u_h^j, w_h^j) = (q^I - q^P, w_h^j),$$

$$(4.19) (a^{-1}(c^j_\sigma)u^j_h, v^j_h) - (p^j_h, \nabla_h \cdot v^j_h) = (\rho(c^j_\sigma)g, v^j_h)$$

for all $(v_h^j, w_h^j) \in U \times P$ and

(4.20)
$$\left(\varphi \mathfrak{d}_{t} c_{\sigma}^{j}, d_{h}^{j} \right) + \mathcal{B}(c_{\sigma}^{j}, d_{h}^{j}; u_{h}^{j}) + \sigma \mathcal{S}(c_{\sigma}^{j}, d_{h}^{j}) = (\hat{c}q^{I}, d_{h}^{j})$$

for all $d_h^j \in V_{dG}^j$.

Definition 4.21. Define the RT-cdG approximation

$$(u_h, p_h, c_h) \in \prod_{j=1}^N U \times \prod_{j=1}^N P \times \prod_{j=1}^N V_{cdG}^j$$

to (4.1)-(4.6) as that generated by the algorithm: For $1 \leq j \leq N$ and $c_h^{j-1} \in V_{cdG}^j$ find $(u_h^j, p_h^j, c_h^j) \in U \times P \times V_{cdG}^j$ such that

$$(\nabla_h \cdot u_h^j, w_h^j) = (q^I - q^P, w_h^j),$$

$$(4.23) (a^{-1}(c_h^j)u_h^j, v_h^j) - (p_h^j, \nabla_h \cdot v_h^j) = (\rho(c_h^j)g, v_h^j)$$

for all $(v_h^j, w_h^j) \in U \times P$ and

$$\left(\varphi \mathfrak{d}_{\mathfrak{t}} c_h^j, d_h^j\right) + \mathcal{B}(c_h^j, d_h^j; u_h^j) + \sigma \mathcal{S}(c_h^j, d_h^j) = (\hat{c}q^I, d_h^j)$$

for all $d_h^j \in V_{cdG}^j$.

Theorem 4.25. The solution $c_{\sigma} \in \Pi_{j=1}^{N} V_{dG}^{j}$ defined in Definition 4.17 converges to $c_{h} \in \Pi_{j=1}^{N} V_{cdG}^{j}$ defined in Definition 4.21 as $\sigma \to \infty$, i.e.,

$$\lim_{\sigma \to \infty} (c_{\sigma} - c_h) = 0.$$

Proof. Following the argument of Section 2 we use Lemma 4.14 and (4.16). In order to complete the proof using this argument we must show that for every sequence $\{c_i^j\}_i$ with elements in V_{dG}^j and $\lim_{i\to\infty}c_i^j=c^j\in V_{\mathrm{dG}}^j$ we have

(4.27)
$$\lim_{i \to \infty} \mathcal{B}(c_i^j, d_h^j; u^j(c_i^j)) = \mathcal{B}(c^j, d_h^j; u^j(c^j)) \qquad \forall d_h^j \in V_{\mathrm{dG}}^j$$

as in (2.11), where $u^j(\cdot)$ is the element in U solving (4.18)-(4.19) for a given element of V_{dG}^j . Note that $u^j:V_{\mathrm{dG}}^j\to U$ is a continuous map and so $\lim_{i\to\infty}u^j(c_i^j)=u^j(\lim_{i\to\infty}c_i^j)=u^j(c^j)$. This also holds for derivatives as they are taken piecewise. Thus (4.27) holds at each timestep and for the whole discrete solution in time. \square

5. Numerical Experiments

We present numerical experiments to illustrate Theorems 3.19, 4.25 and investigate further the performance of the cdG method.

The results were produced using the C++ library deal.ii [2, 3] using both the super penalty approach (as $\sigma \to \infty$) and a direct cdG method, i.e., where the test functions are in $V_{\rm cdG}$ and therefore by construction there will be no jumps across edges in $\mathcal{E}_{\rm cG}$. For further details of the implementation we refer to [10].

5.1. Equations of Non-negative Characteristic Form. Let $\Omega = (0,1)^2$. We seek to solve

$$-\varepsilon \Delta u + (1,1) \cdot \nabla u = f.$$

Given homogeneous Dirichlet boundary conditions f is chosen such that the solution is given by

$$u(x,y) := \left(x - \frac{\mathrm{e}^{(x-1)/\varepsilon} - \mathrm{e}^{-1/\varepsilon}}{1 - \mathrm{e}^{-1/\varepsilon}}\right) \left(y - \frac{\mathrm{e}^{(y-1)/\varepsilon} - \mathrm{e}^{-1/\varepsilon}}{1 - \mathrm{e}^{-1/\varepsilon}}\right).$$

For $0 < \varepsilon \ll 1$ this problem exhibits exponential boundary layers along the outflow boundaries x = 1 and y = 1 of width $\mathcal{O}(\varepsilon)$. We consider a uniformly refined mesh of squares and set r = 1 (piecewise bilinear polynomials).

We first look at an example without a layer by setting $\varepsilon = 10$. We set $\mathcal{T}_{cG} = \mathcal{T}_h$, i.e., the cG method. Figure 5.1 shows the behaviour of the difference between the dG and cG approximations in the $L^2(\mathcal{T}_h)$ norm, $H^1(\mathcal{T}_h)$ semi-norm and the $L^2(\mathcal{T}_h)$

norm of the jumps across edges (represented by $[\![\cdot]\!]$). As σ grows the difference in each norm decreases linearly. The jumps in either approximation are already very small, i.e., the dG approximation is very close to an element in the cG space. We do not see oscillations polluting the continuous approximation.

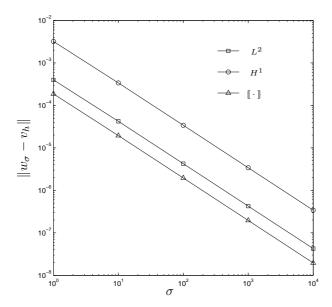
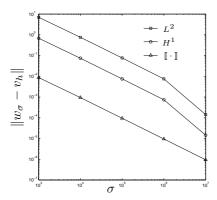
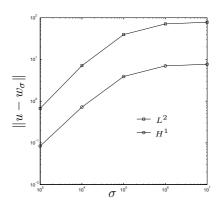


FIGURE 5.1. Example 5.1 with $\varepsilon = 10$ and $\mathcal{T}_{cG} = \mathcal{T}_h$. As the penalty parameter is increased the difference between the cG and dG approximations decreases linearly in the given norms.

We now motivate the cdG method by choosing $\varepsilon = 10^{-4}$ and again setting $\mathcal{T}_{cG} = \mathcal{T}_h$. The example now has a sharp layer at the outflow boundaries. We see in Figure 5.2(a) that increasing σ gives a linear response to the error as in Figure 5.1. When we look at the error in the dG approximation in Figure 5.2(b) we see that the approximation becomes worse as the penalty is increased. The layer causes non-physical oscillations to pollute the approximation. Although we see convergence of the dG approximation to the cG approximation this property is not desirable.

We now consider the cdG method with $\varepsilon = 5 \times 10^{-3}$ and 5×10^{-4} , values chosen so that the layer is partially resolved in the first case and not resolved in the second. The behaviour as σ is increased is the same as in the case $\mathcal{T}_h = \mathcal{T}_{cG}$ and so we do not plot this. We set $h = 2^{-5}$ and $\mathcal{T}_{cG} = (0, 1 - ah)^2$. Varying $a \in \mathbb{Z}$ determines the number of rows in the dG region at the outflow boundary. For $\varepsilon = 5 \times 10^{-3}$ oscillations are apparent in the cG approximation but the mesh is sufficiently refined so that they are not large. Decreasing a (that is, moving from a fully discontinuous approximation towards a fully continuous approximation) results in a small increase in the error of the approximation which can be seen in Figure 5.3. The continuous, discontinuous and continuous-discontinuous approximations are very close in the H^1 semi-norm, including when \mathcal{T}_{cG} covers the layer. When $\varepsilon = 5 \times 10^{-4}$ the layer is sufficiently sharp to induce large oscillations in the fully continuous approximation. By choosing the \mathcal{T}_{cG} region to allow discontinuities in at least the final element we may achieve a reduction in degrees of freedom to approximately 30% of the





- (a) The difference between the cG and dG approximations.
- (b) The error in the dG approximation.

FIGURE 5.2. Example 5.1 with $\varepsilon = 10^{-4}$ and $\mathcal{T}_{cG} = \mathcal{T}_h$. Now the problem has a layer and the error in the dG approximation grows as σ is increased. Non-physical oscillations pollute the approximation.

discontinuous approximation with very little effect on the error (4096 degrees of freedom for the dG approximation, 1024 for the cdG approximation and 1276 for the cdG approximation with one row of dG elements at the outflow boundary).

The dependence of the cdG approximation on the choice of \mathcal{T}_{cG} is further studied in [13]. For the linear problem we in particular refer to the numerical experiments in Section 8.3.

5.2. Incompressible Miscible Displacement: The "Quarter of Five Spot" Problem. As well as verifying Theorem 4.25 we wish to show that if the region where continuous elements are used is chosen appropriately there is little difference in the approximations via the RT-cdG or RT-dG method (where the concentration is approximated in the dG space).

We study a standard example [4, 14, 24] to illustrate the performance of the cdG method for the incompressible miscible displacement problem (4.1)-(4.6). With $\Omega=(0,1)^2$ the injection (resp. extraction) well is located at (1,1) (resp. (0,0)). The injection and extraction strength are represented over one element by piecewise constant functions such that $\int_{\Omega} q^I d\mathbf{x} = \int_{\Omega} q^P d\mathbf{x} = 0.018$. In (4.7) we set $d_l = 1.8 \times 10^{-4}$, $d_m = 1.8 \times 10^{-6}$ and $d_t = 1.8 \times 10^{-5}$. The porosity is set to 0.1. The concentration dependent viscosity is given by $\mu(c) = \mu(0)(1 + (\mathcal{M}^{1/4} - 1)c)^{-4}$ where $\mathcal{M} = 41.0$ is the mobility ratio (the ratio of the viscosity of the fluids), and $\mu(0) = 1$. For the initial concentration we set $c_0 = 0$ corresponding to Ω uniformly filled with one fluid. Set $\mathbb{K} = 0.0288\mathbb{I}$. We consider a uniform refinement of Ω into squares of side $h = 2^{-4}$ with timestep 4×10^{-3} and time interval (0.0, 2.0). With these values a sharp front in the concentration component spreads from the injection to extraction point. As can be seen in Figure 5.6(d) this causes oscillations in the continuous approximation.

First we present the difference between the dG approximation and the cG approximation (i.e., with $\mathcal{T}_{cG} = \mathcal{T}_h$) as $\sigma \to \infty$. In Figure 5.4 we show $||c_{\sigma} - c_h||$ in both the L^2 norm against time and the $L^2((0,T);L^2(\Omega))$ norm against increasing

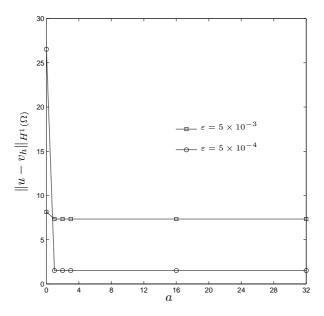


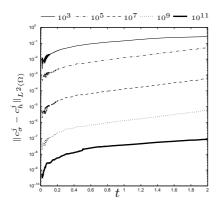
FIGURE 5.3. Example 5.1 varying \mathcal{T}_{cG} for a partially resolved and unresolved layer. In both cases the increase in the error as \mathcal{T}_{cG} covers the layer (a=0) is apparent. When the layer is sharper the increase is more pronounced as severe oscillations pollute the approximation.

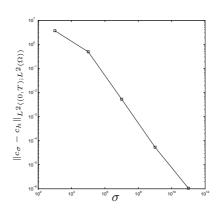
 σ . In Figure 5.4(a) we see a sharp increase in the error over the first few iterations. The initial conditions are in the continuous approximation space so the cG and dG approximations are close. As the layer spreads through the domain the difference between the cG and dG approximations for a given σ in the L^2 norm increases slowly. This is because the number of edges in the vicinity of the layer increases. Figure 5.4(b) shows the same behaviour as the stationary examples in Section 5.1.

Picking \mathcal{T}_{cG} in Section 5.1 was done via knowledge of the true solution and hence knowledge of any layers. We do not have this luxury for the problem considered in this section. We therefore undertake the following procedure for determining \mathcal{T}_{cG} :

- (1) Determine the initial pressure and velocity given c_h^0 and the injection profile.
- (2) Solve for the first time step using a RT-dG method to find a discontinuous c_h^1 .
- (3) For all edges determine $\|[c_h]\|_{L^2(e)}$.
- (4) Flag every cell where each edge satisfys $||[c_h]||_{L^2(e)} < \text{tol}$.
- (5) If every edge of a cell is flagged set that cell to be part of \mathcal{T}_{cG} in the next iteration. Otherwise the element will be in \mathcal{T}_{dG} .
- (6) For n iterations use the cdG mesh defined in the previous step.
- (7) For the $(n+1)^{th}$ iteration reset the mesh to be entirely dG, i.e., $\mathcal{T}_{cG}^{n+1} = \emptyset$ for the concentration component, then return to step (3).

The number of iterations between each cdG refinement and the tolerance should take the expected motion of the fluid and the time step into account. We do not





- (a) Evolution of the difference between the cG and dG approximations for $\sigma=10^3$ to 10^{11} in L^2 norm.
- (b) Plot of the difference between the cG and dG approximations in the $L^2(L^2)$ norm as σ is increased.

FIGURE 5.4. The effect of increasing σ for Example 5.2 with $\mathcal{T}_{cG} = \mathcal{T}_h$.

consider increasing σ for the cdG method, but rather study the performance of the method as the tolerance is increased by comparing the cdG approximation with a dG approximation where $C_{\rm pen}=10$ and $\sigma=0$. With these parameters we set the number of iterations between redefining the cdG space to be 5. We note that this procedure is computationally efficient as each iteration has comparable cost to an iteration of a mesh-adaptive algorithm; indeed, the above procedure can be included in such adaptive algorithm.

In Figure 5.5 we see that as the tolerance is decreased the difference between the dG and cdG approximations in the L^2 norm gets smaller. With a smaller tolerance fewer cells are marked as being continuous. The difference introduced by using some continuous elements does not seem to propagate in time.

In Table 5.1 we see that the number of degrees of freedom saved over the simulation (500 steps with $T=2.0,~\Delta t=4\times 10^{-3}$) is considerable. The effect on the approximation is however small measured in the $L^2(L^2)$ norm. The number of degrees of freedom for the cG method is not 128,000 as would be expected (one degree of freedom per vertex on a 16 × 16 square mesh for 500 timesteps) due to every fifth iteration being computed with the dG method.

In Figure 5.6 we show the dG, cG and cdG approximations after 380 timesteps. There is no visible difference between the plots for dG and cdG at each tolerance (Figures 5.6 (a), (b) and (c)). However for the fully continuous approximation the oscillations induced by the layer are clearly visible and distort the plot.

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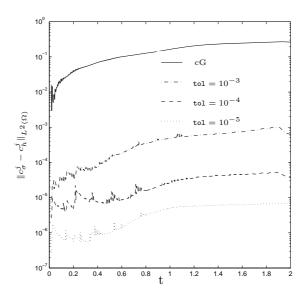
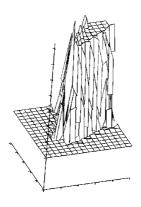


FIGURE 5.5. The behaviour of the cdG approximation for Example 5.2. Using some continuous elements does not dramatically increase the error of the cdG approximation compared to the dG approximation.

Table 5.1. The number of degrees of freedom used for 500 timesteps in Example 5.2. When $tol = 10^{-5}$ only 74.6% of the degrees of freedom are used compared to 43% for the continuous approximation.

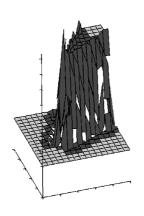
tol	dofs	$ c_{\sigma}-c_{h} _{L^{2}((0,T);L^{2}(\Omega))}$
cG	219,470	3.9970×10^{0}
10^{-3}	323,488	1.2073×10^{-2}
10^{-4}	355,328	7.0904×10^{-4}
10^{-5}	382,384	1.0455×10^{-4}
dG	512,000	0.0000×10^{0}

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(a) The fully discontinuous approximation.

(b) The cdG approximation with tol = 10^{-4} .



(c) The cdG approximation with tol = 10^{-3} .

(d) The fully continuous approximation.

FIGURE 5.6. A plot of the cG method, cdG method and dG method at time 1.52 (380 time steps). The discontinuous region is marked in dark grey for the cdG method. There is no appreciable difference between the first three plots. The oscillations are clearly visible in the fully continuous plot.

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