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COURSE OF NUMERICAL ANALYSIS FOR PARTIAL DIFFERENTIAL EQUATIONS

DISCONTINUOUS GALERKIN APPROXIMATION FOR CARDIAC ELECTROPHYSIOLOGY

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1 Introduction

1.1 Abstract

The aim of the project is to study and implement a suitable numerical scheme for the resolution of the *Bidomain Problem*, a famous system of equations that has been developed in the context of the electrophysiology of human heart. This work is basically the continuation of a two-years-long study carried out by three past course projects ([2], [1], [4]). In particular, the very goal of this project is to improve the results obtained in [4] (Marta and Perego) for the Bidomain model. In fact, even if a *Discontinuous Galerkin* discretization has been successfully implemented, results are not satisfactory from the point of view of stability and convergence. We think this notice is noteworthy as this work is primarily based on these provided data and codes. Through this article, it will be illustrated how we managed to solve these problems extending, optimizing and correcting these past numerical strategies.

1.2 The physical problem

We intend to present the physical meaning of the Bidomain equations very briefly since it has already been widely shown in the previous project (Marta and Perego). For a more complete explanation, we instead refer to [5].

The mechanical contraction and expansion of human heart has its origin in the *electrical activation* of the cardiac cells. At every heart-beat, myocytes are activated and deactivated following a characteristic electrical cycle (fig 1).

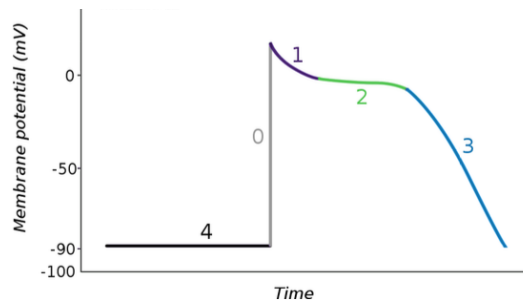


Figure 1: Membrane potential in function of time (one cardiac cycle)

The cell is initially at rest ($-90mV$, step 4). At a certain point, its potential increases rapidly ($\approx 2ms$) and reaches the value of $+20mV$: the cell is activated. Later, a plateau near $0mV$ is observed and then a slow repolarization to the initial potential.

From a microscopical point of view, we could study the dynamics acting in each single cell (as a consequence of the passage of chemical ions through specific channels, e.g. Ca^{2+} , Na^+ , K^+). From a macroscopical point of view, instead, one can think about it as a continuous electrical diffusion over the entire cardiac surface. Even if this consists in a very rapid phenomenon, the study of such propagation could be very interesting in order, for instance, to detect diseases in sick patients.

1.3 Mathematical models

Starting from the circuit in figure 2, applying some general electromagnetism laws and some calculations, the Bidomain model has been formulated (see [5] for more details and/or [3] for the complete passages).

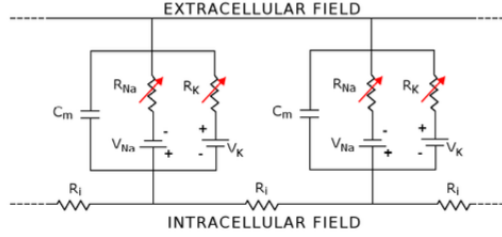


Figure 2: Simplified circuit to model the intracellular and extracellular potentials dynamics

The general formulation is then:

Definition 1 (Bidomain model).

$$\begin{cases} \chi_m C_m \frac{\partial V_m}{\partial t} - \nabla \cdot (\Sigma_i \nabla \phi_i) + \chi_m I_{ion} = I_i^{ext} & \text{in } \Omega_{mus} \times (0, T] \\ -\chi_m C_m \frac{\partial V_m}{\partial t} - \nabla \cdot (\Sigma_e \nabla \phi_e) - \chi_m I_{ion} = -I_e^{ext} & \text{in } \Omega_{mus} \times (0, T] \end{cases}$$

where:

- ϕ_i, ϕ_e are the *Intracellular and Extracellular Potentials* (unknowns)
- $V_m = \phi_i - \phi_e$ is the *Trans-membrane Potential*
- $\chi_m, C_m, \Sigma_i, \Sigma_e$ are known constants
- I_i^{ext}, I_e^{ext} are applied currents
- I_{ion} is the *Ionic Current*

- Ω_{mus} is the cardiac domain (myocardium + endocardium + epicardium)

Actually, this system is not complete since it misses boundary and initial conditions and a suitable model for I_{ion} . Initial conditions and Neumann boundary conditions for ϕ_i and ϕ_e are then imposed. For the definition of I_{ion} , instead, a *reduced ionic model* is chosen, in particular the *FitzHugh-Nagumo model*. Summing up:

Definition 2 (Bidomain + FitzHugh-Nagumo model with Neumann boundary conditions).

$$\left\{ \begin{array}{ll} \chi_m C_m \frac{\partial V_m}{\partial t} - \nabla \cdot (\Sigma_i \nabla \phi_i) + \chi_m I_{ion}(V_m, w) = I_i^{ext} & \text{in } \Omega_{mus} \times (0, T] \\ -\chi_m C_m \frac{\partial V_m}{\partial t} - \nabla \cdot (\Sigma_e \nabla \phi_e) - \chi_m I_{ion}(V_m, w) = -I_e^{ext} & \text{in } \Omega_{mus} \times (0, T] \\ I_{ion}(V_m, w) = k V_m (V_m - a)(V_m - 1) - w & \text{in } \Omega_{mus} \times (0, T] \\ \frac{\partial w}{\partial t} = \epsilon (V_m - \gamma w) & \text{in } \Omega_{mus} \times (0, T] \\ \Sigma_i \nabla \phi_i \cdot n = b_i & \text{on } \partial \Omega_{mus} \times (0, T] \\ \Sigma_e \nabla \phi_e \cdot n = b_e & \text{on } \partial \Omega_{mus} \times (0, T] \\ \text{Initial conditions for } \phi_i, \phi_e, w & \text{in } \Omega_{mus} \times \{t = 0\} \end{array} \right.$$

where:

- w is the *gating variable* (unknown)
- k, a, ϵ, γ are known constants
- b_i, b_e are the boundary conditions data
- n is the outward normal vector

From now on, the system of definition 2 will be the reference analytical problem for the development of numerical schemes.

To conclude, there exist other famous and useful models, such as the *Monodomain model*. But this is just a simplification of the Bidomain as in this case it is assumed that ϕ_i and ϕ_e are proportional. However, thanks to its simplicity, we often tested the code starting from the Monodomain implementation of the project [1] instead of analyzing directly the Bidomain.

1.4 Formulations

1.4.1 DG discrete formulation

We have seen the bidomain model in a complete form in 2. We introduce now a triangulation τ_h over Ω , with $\mathcal{F}_h = \mathcal{F}_h^I \cup \mathcal{F}_h^B$ set of the faces of the elements, which includes the internal and boundary faces respectively, and the DG space $V_h^k = \{v_h \in L^2 : v_h|_{\mathcal{K}} \in \mathbb{P}^k(\mathcal{K}) \quad \forall \mathcal{K} \in \tau_h\}$, where k is the degree of the piecewise continuous polynomial.

We obtain the semi discrete DG formulation:

$$\begin{aligned}
& \text{For any time } t \in [0, T] \text{ find } \Phi_h(t) = [\phi_i^h(t), \phi_e^h(t)]^T \in [V_h^k]^2 \text{ and } \omega_h(t) \in V_h^k : \\
& \sum_{K \in \tau_h} \int_K \chi_m C_m \frac{\partial V_m^h}{\partial t} V_h d\omega + a_i(\phi_i^h, v_h) + \sum_{K \in \tau_h} \int_K \chi_m k (V_m^h - 1)(V_m^h - a) V_m^h v_h d\omega + \\
& \quad + \sum_{K \in \tau_h} \int_K \chi_m w_h v_h d\omega = (I_i^{ext}, v_h) \quad \forall v_h \in V_h^p \\
& - \sum_{K \in \tau_h} \int_K \chi_m C_m \frac{\partial V_m^h}{\partial t} V_h d\omega + a_e(\phi_e^h, v_h) - \sum_{K \in \tau_h} \int_K \chi_m k (V_m^h - 1)(V_m^h - a) V_m^h v_h d\omega + \\
& \quad - \sum_{K \in \tau_h} \int_K \chi_m w_h v_h d\omega = (-I_e^{ext}, v_h) \quad \forall v_h \in V_h^p \\
& \sum_{K \in \tau_h} \int_K \frac{\partial w_h}{\partial t} v_h d\omega = \sum_{k \in \tau_h} \int_K \epsilon (V_m^h - \gamma w_h) v_h d\omega \quad \forall v_h \in V_h^p
\end{aligned}$$

where:

$$\begin{aligned}
a_k(\phi_k^h, v_h) &= \sum_{K \in \tau_h} \int_K (\Sigma_k \nabla_h \phi_k^h) \cdot \nabla_h v_h d\omega - \sum_{F \in \mathcal{F}_h^I} \int_F \{ \{ \Sigma_k \nabla_h \phi_k^h \} \} \cdot [[v_h]] d\sigma + \\
& - \delta \sum_{F \in \mathcal{F}_h^I} \int_F \{ \{ \Sigma_k \nabla_h v_h \} \} \cdot [[\phi_k^h]] d\sigma + \sum_{F \in \mathcal{F}_h^I} \int_F \gamma [[\phi_k^h]] \cdot [[v_h]] d\sigma \quad \text{for } k=i,e \\
(I_i^{ext}, v_h) &= \sum_{K \in \tau_h} \int_K I_i^{ext} v_h d\omega + \int_{\partial\Omega} b v_h d\sigma \\
(-I_e^{ext}, v_h) &= - \sum_{K \in \tau_h} \int_K I_e^{ext} v_h d\omega + \int_{\partial\Omega} b v_h d\sigma
\end{aligned}$$

Moreover, according to the choice of the coefficient δ we can define:

- $\delta = 1$: Symmetric Interior Penalty method (SIP)
- $\delta = 0$: Incomplete Interior Penalty method (IIP)

- $\delta = -1$: Non Symmetric Interior Penalty method (NIP)

To see in more detail the procedure see [4].

1.4.2 Algebraic formulation

Taking $\{\varphi_j\}_{j=1}^{N_h}$ base of V_h^k , so that we can write

$$\begin{aligned}\Phi_h(t) &= \begin{bmatrix} \phi_i^h(t) \\ \phi_e^h(t) \end{bmatrix} = \begin{bmatrix} \sum_{j=1}^{N_h} \phi_{i,j}(t) \varphi_j \\ \sum_{j=1}^{N_h} \phi_{e,j}(t) \varphi_j \end{bmatrix} \\ w_h(t) &= \sum_{j=1}^{N_h} w_j(t) \varphi_j \\ V_m^h(t) &= \sum_{j=1}^{N_h} V_{m,j}(t) \phi_j = \sum_{j=1}^{N_h} (\phi_{i,j}(t) - \phi_{e,j}(t)) \varphi_j\end{aligned}$$

Also we can introduce the matrices:

$$\left. \begin{aligned} V_{ij} &= \int_{\Omega} \nabla \varphi_j \cdot \nabla \varphi_i \\ I_{i,j}^T &= \sum_{F \in F_h^I} \int_F \{ \{ \nabla \varphi_j \} \} \cdot [[\varphi_i]] \\ I_{i,j} &= \sum_{F \in F_h^I} \int_F [[\varphi_j]] \cdot \{ \{ \nabla \varphi_i \} \} \\ S_{i,j} &= \sum_{F \in F_h^I} \int_F \gamma [[\varphi_j]] \cdot [[\varphi_i]] \end{aligned} \right\} A = (V - I^T - \theta I + S) \quad (1)$$

$$A_i = \Sigma_i A \quad \text{Intra-cellular stiffness matrix} \quad (2)$$

$$A_e = \Sigma_e A \quad \text{Extra-cellular stiffness matrix} \quad (3)$$

$$M_{ij} = \sum_{K \in \tau_h} \int_K \text{Mass matrix } \varphi_j \varphi_i \quad (4)$$

$$C(u_h)_{ij} = \sum_{K \in \tau_h} \int_K \chi_m k(u_h - 1)(u_h - a) \varphi_j \varphi_i \quad \text{Non linear matrix} \quad (5)$$

$$F_k = \begin{bmatrix} F_{i,k} \\ F_{e,k} \end{bmatrix} = \begin{bmatrix} \int_{\Omega} I_i^{ext} \varphi_k - \sum_{F \in F_h^B} \int_F b_i \varphi_k \\ - \int_{\Omega} I_e^{ext} \varphi_k - \sum_{F \in F_h^B} \int_F b_e \varphi_k \end{bmatrix} \quad (6)$$

Therefore our algebraic formulation is:

Find $\Phi_h(t) = [\phi_i^h(t), \phi_e^h(t)]^T \in [V_h^k]^2$ and $\omega_h(t) \in V_h^k$ for any $t \in (0; T]$ such that:

$$\begin{aligned}\chi_m C m M \dot{V}_m^h + A_i \phi_i^h + C(V_m^h) V_m^h + \chi_m M \omega_h &= F_i^h \\ -\chi_m C m M \dot{V}_m^h + A_e \phi_e^h - C(V_m^h) V_m^h - \chi_m M \omega_h &= F_e^h \\ M \dot{\omega}_h(t) &= \epsilon M (V_m^h(t) - \gamma \omega_h(t))\end{aligned} \quad (7)$$

Rewrite it with the block matrix and assuming that M is non singular:
Find $\Phi_h(t) = [\phi_i^h(t), \phi_e^h(t)]^T \in [V_h^k]^2$ and $\omega_h(t) \in V_h^k$ for any $t \in (0; T]$ such that:

$$\begin{aligned} & \chi_m C_m \begin{bmatrix} M & -M \\ -M & M \end{bmatrix} \begin{bmatrix} \dot{\Phi}_i^h(t) \\ \dot{\Phi}_e^h(t) \end{bmatrix} + \begin{bmatrix} A_i & 0 \\ 0 & A_e \end{bmatrix} \begin{bmatrix} \Phi_i^h(t) \\ \Phi_e^h(t) \end{bmatrix} + \\ & \begin{bmatrix} C(V_m^h) & -C(V_m^h) \\ -C(V_m^h) & C(V_m^h) \end{bmatrix} \begin{bmatrix} \Phi_i^h(t) \\ \Phi_e^h(t) \end{bmatrix} + \chi_m \begin{bmatrix} M & 0 \\ 0 & -M \end{bmatrix} \begin{bmatrix} w_h(t) \\ w_h(t) \end{bmatrix} = \begin{bmatrix} F_i^h \\ F_e^h \end{bmatrix} \quad (8) \\ & \omega_h(t) = \epsilon(V_m^h(t) - \gamma \omega_h(t)) \end{aligned}$$

2 Dubiner Basis

The most popular orthonormal basis on the reference triangle

$$\hat{K} = \{(\xi, \eta) : \xi, \eta \geq 0, \xi + \eta \leq 1\} \quad (9)$$

is the Dubiner polynomial basis [antonietti]. We consider the transformation in figure 3 between the reference square

$$\hat{Q} = \{(a, b) : -1 \leq a \leq 1, -1 \leq b \leq 1\} \quad (10)$$

and the reference triangle given by

$$\xi := \frac{(1+a)(1-b)}{4}, \eta := \frac{(1+b)}{2} \quad (11)$$

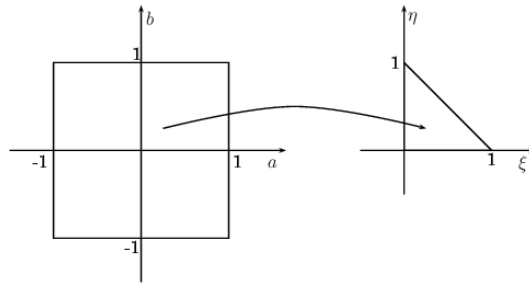


Figure 3: Transformation between the reference square to the reference triangle

The Dubiner basis is then constructed by a generalized tensor product of the Jacobi polynomials on the interval $(-1, 1)$ to form a basis on the reference square, which is then transformed by the above “collapsing” mapping to a basis on the reference triangle.

Definition 3 (Jacobi polynomials). *The Jacobi polynomials evaluated in $z \in \mathbb{R}^n$ are:*

– $n = 0$

$$J_0^{\alpha,\beta}(z) = \overbrace{[1 \quad 1 \quad \dots \quad 1]}^{n \text{ times}} \quad (12)$$

– $n = 1$

$$J_1^{\alpha,\beta}(z) = \frac{1}{2}(\alpha - \beta + (\alpha + \beta + 2) \cdot z); \quad (13)$$

– $n \geq 2$

$$\begin{aligned} J_n^{\alpha,\beta}(z) = \sum_{k=2}^n & \left[\frac{(2k + \alpha + \beta - 1)(\alpha^2 - \beta^2)}{2k(k + \alpha + \beta)(2k + \alpha + \beta - 2)} + \right. \\ & \frac{(2k + \alpha + \beta - 2)(2k + \alpha + \beta - 1)(2k + \alpha\beta)}{2k(k + \alpha + \beta)(2k + \alpha + \beta - 2)} J_{k-1}^{\alpha,\beta}(z) + \\ & \left. - \frac{2(k + \alpha - 1)(k + \beta - 1)(2k + \alpha + \beta)}{2k(k + \alpha + \beta)(2k + \alpha + \beta - 2)} J_{k-2}^{\alpha,\beta}(z) \right] \end{aligned} \quad (14)$$

An important property of this polynomials is:

Proposition 1. $J_i^{\alpha,\beta}(\cdot)$ is orthogonal w.r.t. the Jacobi weight $w(x) = (1 - x)^\alpha(1 + x)^\beta$:

$$\int_{-1}^1 (1 - x)^\alpha (1 + x)^\beta J_m^{\alpha,\beta} J_q^{\alpha,\beta}(x) dx = \frac{2}{2m + 1} \delta_{mq} \quad (15)$$

Thanks to this definition, we can now look more accurately the formula of the Dubiner basis

Definition 4 (Dubiner Basis).

$$\begin{aligned} \phi_{ij}(\xi, \eta) &:= c_{ij}(1 - b)^j J_i^{0,0}(a) J_j^{2i+1,0}(b) = \\ &= c_{ij} 2^j (1 - \eta)^j J_i^{0,0}\left(\frac{2\xi}{1 - \eta} - 1\right) J_j^{2i+1,0}(2\eta - 1) \end{aligned} \quad (16)$$

for $i, j = 1, \dots, p$ and $i + j \leq p$, where

$$c_{ij} := \sqrt{\frac{2(2i + 1)(i + j + 1)}{4^i}} \quad (17)$$

and $J_i^{\alpha,\beta}(\cdot)$ is the i -th Jacobi polynomial

We note that

Proposition 2. $\phi_{ij}(\xi, \eta)$ is orthogonal w.r.t. Legendre internal product:

$$\iint_{\hat{K}} \phi_{ij}(\xi, \eta) \phi_{mq}(\xi, \eta) d\xi d\eta = \delta_{im} \delta_{jq} \quad (18)$$

Indeed, we obtain a mass matrix diagonal as we can see in figure 4

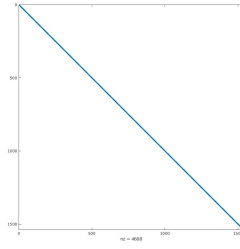


Figure 4: Elements non zero in the mass matrix w.r.t Dubiner basis

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