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A software benchmark for cardiac elastodynamics

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Abstract

In cardiovascular mechanics, reaching consensus in simulation results within a physiologically relevant range of parameters is essential for reproducibility purposes. Although currently available benchmarks contain some of the features that cardiac mechanics models typically include, some important modeling aspects are missing. Therefore, we propose a new set of cardiac benchmark problems and solutions for assessing passive and active material behaviour, viscous effects, and pericardial boundary condition. The problems proposed include simplified analytical fiber definitions and active stress models on a monoventricular and biventricular domains, allowing straightforward testing and validation with already developed solvers.

1 Introduction

In computational biomechanics in general, efforts of defining benchmarks for verification and validation have been sparse throughout the years and are application dependent [1–4].

In particular, in the context of cardiovascular mechanics, reaching consensus in simulation results is an important task, since, for a given set of physical constants, different numerical solutions can be obtained, e.g., due to discretization strategies, polynomial degree of basis functions, numerical quadratures and time integrators [5–14]. Specially, when parameters are optimized from clinical data, it is crucial that these parameters may be valid for other groups and hence could be reused, given the high complexity involved in solving these inverse problems [15–17]

In [2] was proposed a first benchmark containing some of the features that cardiac mechanics models typically include. However some important features are lacking, such as the inclusion of state-of-the-art passive and active models, idealized geometrical dimensions, boundary conditions as well as time dependent effects (i.e. inertia and viscosity).

Therefore, we propose here a new set of cardiac benchmark problems and computed solutions for assessing passive and active material behaviour as in [18] with viscous effects and pericardial boundary conditions as in [7]. The problems proposed in this work include a simplified analytical fiber definition with active stress model by [19], allowing straightforward testing and validation with already developed solvers. The benchmark definition is agreed upon nine different research groups, who computed their solutions with numerical methods and software of their choice. A comparison is carried out among the solutions computed by the different groups, whose results demonstrate a substantial agreement between the participating teams.

The remainder of this article is organized as follows. Section 2 describes the mechanical problem in continuous form, its material properties and boundary conditions. Section 3 proposes the first benchmark problem in a monoventricular domain, with analytical geometry, fibers orientation and simulations setup for blinded and non-blinded phases among participants. Section 4 proposes a second benchmark for a biventricular domain, with state-of-the-art fiber orientation, constitutive law and analogous setups to Section 3 comprising a non-blinded phase only. Section 5 describes all participant software as well as their solver

strategies. Section 6 contains computed results with a qualitative and quantitative analysis of the first and second benchmark. Section 7 provides a discussion of the different approaches and results, finally in Section 9 the conclusion.

2 The mathematical model

2.1 Strong form

We define the problem in a domain $\Omega \subset \mathbb{R}^3$ with boundary $\partial\Omega := \Gamma_{top} \cup \Gamma_{epi} \cup \Gamma_{endo}$. Let us denote by $\mathbf{u} : \Omega \rightarrow \mathbb{R}^3$ the displacement field to be found, $\mathbf{u}(\mathbf{X})$ its evaluation in \mathbf{X} for $\mathbf{X} \in \Omega$, by $\mathbb{F} := \mathbb{I} + \mathbf{Grad}(\mathbf{u}) := \mathbb{I} + \frac{\partial \mathbf{u}}{\partial \mathbf{X}}$ the deformation gradient, $\mathbf{Div}(\mathbf{u}) := \frac{\partial}{\partial \mathbf{X}} \cdot \mathbf{u}$ the divergence, $J := \det(\mathbb{F}(\mathbf{u})) := \det(\mathbb{F})$ the jacobian, $\mathbb{E} = \frac{1}{2}(\mathbb{C} - \mathbb{I})$ the Green-Lagrange tensor, $\mathbb{C} := \mathbb{F}^\top \mathbb{F}$ the right Cauchy tensor and \mathbb{I} and the identity matrix respectively.

Let us denote $\mathbb{T} := \mathbb{T}(\mathbf{u})$ the Cauchy stress tensor associated to the unknown displacement field \mathbf{u} and the second Piola Kirchhoff stress tensor denoted by $\mathbb{S} := J\mathbb{F}^{-1}\mathbb{T}\mathbb{F}^{-\top}$, the problem to solve over the time-interval $(0, 1]$, is described by the equations:

$$\begin{aligned} \rho \ddot{\mathbf{u}} - \mathbf{Div}(J\mathbb{T}\mathbb{F}^{-\top}) &= \mathbf{0} \quad \text{in } \Omega \\ J\mathbb{T}\mathbb{F}^{-\top} \mathbf{N} &= pJ\mathbb{F}^{-\top} \mathbf{N} \quad \text{on } \Gamma_{endo} \\ J\mathbb{T}\mathbb{F}^{-\top} \mathbf{N} \cdot \mathbf{N} + \alpha_{epi} \mathbf{u} \cdot \mathbf{N} + \beta_{epi} \dot{\mathbf{u}} \cdot \mathbf{N} &= 0 \quad \text{on } \Gamma_{epi} \\ J\mathbb{T}(\mathbb{F}^{-\top} \mathbf{N}) \times \mathbf{N} &= \mathbf{0} \quad \text{on } \Gamma_{epi} \\ J\mathbb{T}\mathbb{F}^{-\top} \mathbf{N} + \alpha_{top} \mathbf{u} + \beta_{top} \dot{\mathbf{u}} &= \mathbf{0} \quad \text{on } \Gamma_{top} \end{aligned} \tag{1}$$

with \mathbf{N} the unit wall normal vector and \top the transpose if used as superscript.

2.2 Material model

The material behaviour is characterized via \mathbb{S} including the anisotropic, viscous and active parts, namely

$$\mathbb{S}(t) := \frac{\partial \Psi_{aniso}}{\partial \mathbb{E}} + \frac{\partial \Psi_{visco}}{\partial \dot{\mathbb{E}}} + \tau(t) \mathbf{f} \otimes \mathbf{f}, \tag{2}$$

with each term described below[‡]:

- The anisotropic material energy Ψ_{aniso} describes the nearly incompressible Holzapfel-Ogden material [18] with isochoric-volumetric split, via the isotropic invariant $I_1 = J^{-2/3} \text{tr}(\mathbb{C})$, the transverse isotropic invariants $I_{4f} := \mathbf{f} \cdot \mathbb{C} \mathbf{f}$ and $I_{4s} := \mathbf{s} \cdot \mathbb{C} \mathbf{s}$ for

[‡]Models aiming at characterizing the strain-stress behavior have been substantially studied [18, 20–22]. The literature points at the work of *Guccione et al.* [23] and *Holzapfel et al.* [18], the latter widely used for human cardiac models. Nevertheless, not only one convention has been used to characterize the fiber orientation, especially with the choice of the sheet and sheet-normal directions. Several works implementing either case have shown consistent deformations, endocardial pressure as well as ejection volumes [6, 24–30].

the fiber directions at the reference domain $\mathbf{f}, \mathbf{s} : \Omega \rightarrow \mathbb{R}^3$ and anisotropic invariant $I_{8fs} := \mathbf{f} \cdot \mathbb{C}\mathbf{s}$. Explicitly Ψ_{aniso} is given by:

$$\begin{aligned} \Psi_{aniso} = & \frac{a}{2b} \exp \{b(I_1 - 3) - 1\} + \sum_{i \in \{f, s\}} \frac{a_i}{2b_i} \chi(I_{4i}) (\exp \{b_i(I_{4i} - 1)^2\} - 1) \\ & + \frac{a_{fs}}{2b_{fs}} (\exp \{b_{fs} I_{8fs}^2\} - 1) + \frac{\kappa}{4} (J^2 - 1 - 2 \ln(J)) \end{aligned} \quad (3)$$

with $\chi(x) = x$ if $x > 1$ and 0 otherwise, for $x \in \mathbb{R}_+$, denoting the fiber compression switch model. The last term denotes the incompressibility penalty proposed in [31] with parameter $\kappa > 0$.

A suggested approximation is given by $\chi(x) \approx \frac{1}{1+e^{-k(x-1)}}$, for $k > 0$ a fixed parameter specified later on.

- The viscoelastic energy is characterized with parameter η in the form [32]:

$$\Psi_{visc} := \frac{\eta}{2} \text{tr}(\dot{\mathbb{E}}^2) \quad (4)$$

- The active stress is taken as in [19], characterized by a time-dependent stress function τ , solution to the evolution equation

$$\dot{\tau}(t) = -|a(t)|\tau(t) + \sigma_0|a(t)|_+ \quad (5)$$

denoting $a(\cdot)$ the activation function and σ_0 contractility, and the remaining terms defined as:

$$\begin{aligned} |a(t)|_+ &= \max\{a(t), 0\} \\ a(t) &:= \alpha_{max} \cdot f(t) + \alpha_{min} \cdot (1 - f(t)) \\ f(t) &= S^+(t - t_{sys}) \cdot S^-(t - t_{dias}) \\ S^\pm(\Delta t) &= \frac{1}{2} (1 \pm \tanh(\frac{\Delta t}{\gamma})). \end{aligned} \quad (6)$$

2.3 Pressure model

We consider a time-dependent pressure for (1), derived from the active stress function. The solution $p = p(t)$ is characterized by the evolution equation

$$\dot{p}(t) = -|b(t)|p(t) + \sigma_{mid}|b(t)|_+ + \sigma_{pre}|g_{pre}(t)|_+ \quad (7)$$

with $b(\cdot)$ the activation function described as:

$$\begin{aligned} b(t) &= a_{pre}(t) + \alpha_{pre}g_{pre}(t) + \alpha_{mid} \\ a_{pre}(t) &:= \alpha_{max} \cdot f_{pre}(t) + \alpha_{min} \cdot (1 - f_{pre}(t)) \\ f_{pre}(t) &= S^+(t - t_{sys-pre}) \cdot S^-(t - t_{dias-pre}) \\ g_{pre}(t) &= S^-(t - t_{dias-pre}) \end{aligned} \quad (8)$$

and S^\pm defined as in (6).

3 Benchmark 1: monoventricular mechanics

3.1 Geometry

Using the same analytical formula as in [2], we define the domain via the parametrization (in \mathbb{R}^3) for a truncated ellipsoid, i.e., satisfying:

$$(x, y, z) = (r_{\text{long}} \cos(\mu), r_{\text{short}} \sin(\mu) \cos(\theta), r_{\text{short}} \sin(\mu) \sin(\theta)) \quad (9)$$

with the following dimensions:

- The endocardial surface

$$r_{\text{short}} = 2.5 \times 10^{-2}[m], \quad r_{\text{long}} = 9.0 \times 10^{-2}[m], \quad \mu \in [-\pi, -\arccos(\frac{5}{17})], \theta \in [-\pi, \pi] \quad (10)$$

- The epicardial surface

$$r_{\text{short}} = 3.5 \times 10^{-2}[m], \quad r_{\text{long}} = 9.7 \times 10^{-2}[m], \quad \mu \in [-\pi, -\arccos(\frac{5}{20})], \theta \in [-\pi, \pi] \quad (11)$$

The domain is created using the software *Gmsh* [33] and distributed to all participants in different formats, created with an element size[§] $h = 5 \times 10^{-3}[m]$. Supplemented material is provided with such data as well as a repository including implementation details[‡].

3.2 Fibers

The definition of fibers is based on a local coordinate system derived from the ellipsoid parametrization. Using the ellipsoid parametrization, a point \mathbf{x} in the domain Ω is described as:

$$\mathbf{x}(\mu, \theta, \bar{t}) = (r_l(\bar{t}) \cos(\mu), r_s(\bar{t}) \sin(\mu) \cos(\theta), r_s(\bar{t}) \sin(\mu) \sin(\theta)), \quad (12)$$

with μ, θ as defined previously and $\bar{t} : \Omega \rightarrow [0, 1]$ is defined as the solution to the problem:

$$\begin{aligned} \Delta \bar{t} &= 0 && \text{in } \Omega \\ \bar{t} &= 0 && \text{on } \Gamma_{\text{endo}} \\ \bar{t} &= 1 && \text{on } \Gamma_{\text{epi}} \\ \frac{\partial \bar{t}}{\partial \mathbf{N}} &= 0 && \text{on } \Gamma_{\text{top}}. \end{aligned} \quad (13)$$

[§]The element size is defined as the optimal edge length around any point node in the mesh with specified target size $h > 0$. Therefore, not a lower or upper edge limit, rather, an averaged value computed to match a user-provided target size. For further details, we refer to [33].

[‡]The repository cardiac-benchmark-toolkit stores the data provided to all teams in several formats `.geo`, `.msh`, `.xdmf`, `.h5`, as well as an user-friendly interface to recreate the monoventricular domain at different mesh sizes.

The tangent basis derived from (12), denoted as $[\mathbf{e}_{\bar{t}}, \mathbf{e}_{\mu}, \mathbf{e}_{\theta}]$, is defined as:

$$\begin{aligned}\tilde{\mathbf{e}}_{\bar{t}} &= \frac{\partial \mathbf{x}}{\partial \bar{t}}, & \tilde{\mathbf{e}}_{\mu} &= \frac{\partial \mathbf{x}}{\partial \mu}, & \tilde{\mathbf{e}}_{\theta} &= \frac{\partial \mathbf{x}}{\partial \theta} \\ \mathbf{e}_{\bar{t}} &= \frac{\tilde{\mathbf{e}}_{\bar{t}}}{\|\tilde{\mathbf{e}}_{\bar{t}}\|_{\mathbb{R}^3}}, & \mathbf{e}_{\mu} &= \frac{\tilde{\mathbf{e}}_{\mu}}{\|\tilde{\mathbf{e}}_{\mu}\|_{\mathbb{R}^3}}, & \mathbf{e}_{\theta} &= \frac{\tilde{\mathbf{e}}_{\theta}}{\|\tilde{\mathbf{e}}_{\theta}\|_{\mathbb{R}^3}},\end{aligned}\tag{14}$$

Using (14), the fiber, sheet-normal and sheet directions are defined as follows:

$$\begin{aligned}\mathbf{f}(\bar{t}, \mu, \theta) &= \sin(\alpha(\bar{t})) \mathbf{e}_{\mu} + \cos(\alpha(\bar{t})) \mathbf{e}_{\theta} \\ \mathbf{n}(\bar{t}, \mu, \theta) &= \frac{\mathbf{e}_{\mu} \times \mathbf{e}_{\theta}}{\|\mathbf{e}_{\mu} \times \mathbf{e}_{\theta}\|_{\mathbb{R}^3}} \\ \mathbf{s}(\bar{t}, \mu, \theta) &= \frac{\mathbf{f}(\bar{t}, \mu, \theta) \times \mathbf{n}(\bar{t}, \mu, \theta)}{\|\mathbf{f}(\bar{t}, \mu, \theta) \times \mathbf{n}(\bar{t}, \mu, \theta)\|_{\mathbb{R}^3}}\end{aligned}\tag{15}$$

with $\alpha(\bar{t}), r_l(\bar{t}), r_s(\bar{t})$ parameters defined as:

$$\begin{aligned}\alpha(\bar{t}) &= (\alpha_{endo} + (\alpha_{epi} - \alpha_{endo})\bar{t}) \frac{\pi}{180} \\ r_l(\bar{t}) &= r_{\text{long_endo}} + (r_{\text{long_epi}} - r_{\text{long_endo}})\bar{t} \\ r_s(\bar{t}) &= r_{\text{short_endo}} + (r_{\text{short_epi}} - r_{\text{short_endo}})\bar{t}\end{aligned}\tag{16}$$

for $r_{\text{long_endo}}, r_{\text{short_endo}}$ the long/short radius in (10) and $r_{\text{long_epi}}, r_{\text{short_epi}}$ the long/short radius in (11).

The computation of the fibers close to the apex is problematic. Given a point in the ellipsoid $\mathbf{x} = (x, y, z)$ and $\bar{t} = \bar{t}(\mathbf{x})$ we propose to compute the associated parameters μ, θ to such a point as $\mu = \text{atan2}(a, b)$ for $a = \frac{\sqrt{y^2 + z^2}}{r_s(\bar{t})}, b = \frac{x}{r_l(\bar{t})}$ and $\theta = 0$ if $\mu \leq 10^{-7}$ else $\theta = \pi - \text{atan2}(z, -y)$.

Depicted in Figure 1 is the labeled ellipsoid geometry, including the fiber and sheet directions.

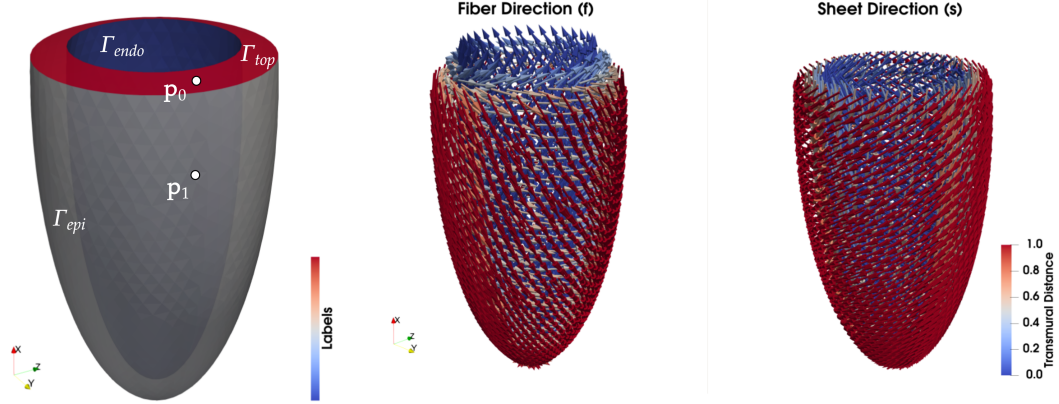


Figure 1: The labeled ellipsoid geometry (left) includes positions of particles $\mathbf{p}_0, \mathbf{p}_1$ for reference. The fiber (center) and sheet (right) directions described in (15) for a $\mp 60^\circ$ angle configuration, are colored using the transmural distance \bar{t} over the domain.

3.3 Step 0 (non-blinded): Splitting passive and active responses

We first perform a validation with teams having access to the solutions of the rest of participants. This served to refine the problem description and to encourage a larger number of participants.

3.3.1 Case A: Active response

Each group solves numerically the equations described in Section 2, with geometry and fibers as in Sections 3.1 and 3.2 respectively, parameters as in Tables 1, 2, 3 and a zero endocardial pressure, i.e. $p = 0$ on Γ_{endo} over all timesteps.

The groups are requested to provide the displacement field $\mathbf{u}_h(\mathbf{X})$ over time at two spatial locations, $\mathbf{p}_0 = (0.025, 0.03, 0)$, $\mathbf{p}_1 = (0, 0.03, 0)$. Such spatial locations do not describe points of the mesh provided to the participants, thus each team must have interpolation algorithms available.

Parameter	$\rho[\frac{kg}{m^3}]$	$\eta[Pa \cdot s]$	$\kappa[Pa]$	$k[-]$	$\alpha_{top}[\frac{Pa}{m}]$	$\alpha_{epi}[\frac{Pa}{m}]$	$\beta_{top}[Pa \cdot \frac{s}{m}]$	$\beta_{epi}[Pa \cdot \frac{s}{m}]$
Value	10^3	10^2	10^6	100	10^5	10^8	5×10^3	5×10^3

Table 1: Parameters describing the strong form of the problem defined in (1).

Parameter	$a[Pa]$	$a_f[Pa]$	$a_{fs}[Pa]$	$a_s[Pa]$	$b[\cdot]$	$b_f[\cdot]$	$b_{fs}[\cdot]$	$b_s[\cdot]$
Value	59.0	18472.0	216.0	2481.0	8.023	16.026	11.436	11.12

Table 2: Parameters of the constitutive law describing the directional behavior through fiber and sheet directions, described in (3).

Parameter	$\sigma_0[Pa]$	$\gamma[s]$	α_{\min}	α_{\max}	$t_{\text{sys}}[s]$	$t_{\text{dias}}[s]$	α_{endo}	α_{epi}
Value	1.5×10^5	0.005	-30	5	0.16	0.484	-60°	+60°

Table 3: Parameters defining the active stress activation function, solution to (5) and fibers' angles at endo/epi-cardium, as in [7].

Depicted in Figure 2 is the evolution of the stress function τ over time for physical parameters specified therein.

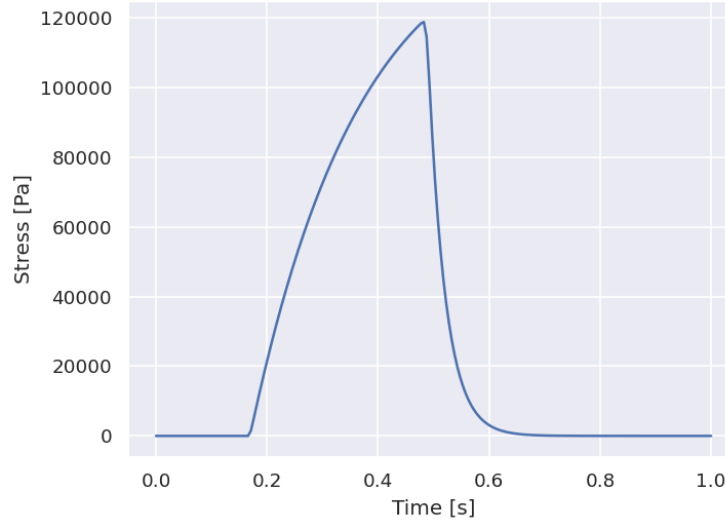


Figure 2: Evolution of the stress function τ described in (5) over the time interval $[0, 1]$ with physiological parameters proposed in Table 3 from [7]. It reaches a maximum value of 118817.07 [Pa].

3.3.2 Case B: Passive response

Each group solves numerically the equations described in Section 2 with geometry, fibers as in Sections 3.1, 3.2, and parameters as in Tables 1, 2, 4 and no active part, i.e. $\tau(t) = 0 \quad \forall t > 0$. The groups are requested to provide the displacement field $\mathbf{u}_h(\mathbf{X})$ over time at $\mathbf{p}_0, \mathbf{p}_1$.

Depicted in Figure 3 is the evolution of the pressure $p(t)$ over time for the parameters specified therein.

Parameter	Values
$\alpha_{min}[-]$	-30
$\alpha_{max}[-]$	5
$\alpha_{pre}[-]$	5
$\alpha_{mid}[-]$	1
$\sigma_{pre}[Pa]$	7000
$\sigma_{mid}[Pa]$	16000
$t_{sys-pre}[s]$	0.17
$t_{dias-pre}[s]$	0.484
$\gamma[s]$	0.005

Table 4: Parameters for the pressure model (7)

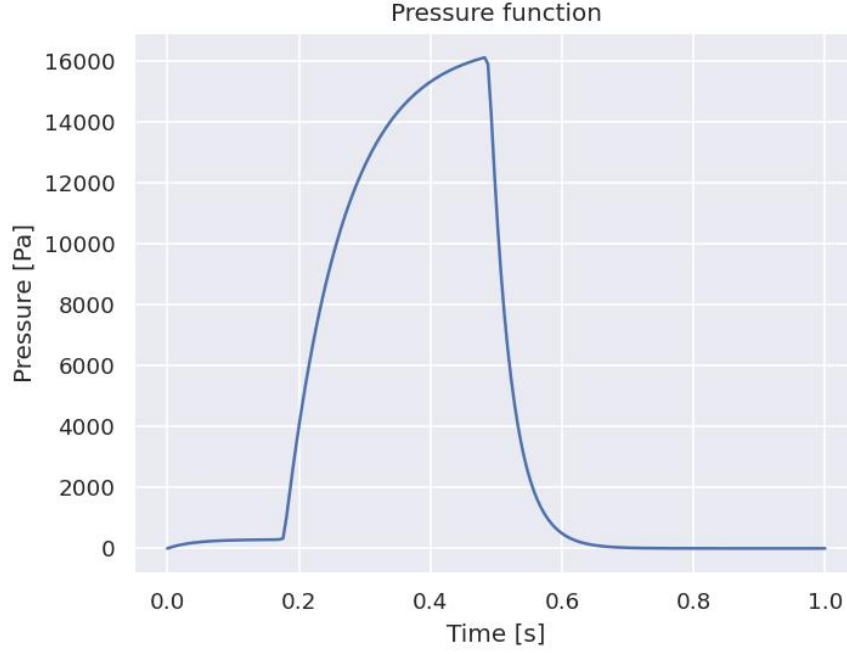


Figure 3: Evolution of the pressure $p(t)$ described in (7) over the time interval $[0, 1]$. Parameters as in Table 4. It reaches a maximum of 16117.52 [Pa].

3.4 Step 1 (non-blinded): active and passive response

Each group solves numerically the equations described in Section 2 with geometry, fibers as in Sections 3.1, 3.2 and parameters as in Tables 1, 2, 3, 4. The groups are requested to provide the displacement field $\mathbf{u}_h(\mathbf{X})$ over time at two spatial locations, $\mathbf{p}_0, \mathbf{p}_1$, as described in Subsection 3.3.1.

3.5 Step 2: Blinded variation of physical parameters

In the second step, all groups are requested to run computations – fully blinded from each other – with modified physical constants with respect to Section 3.4. Here we changed a specific combination of parameters, namely a, a_f, a_{fs}, a_s by a constant factor and σ_0 . The values taken for each parameter combination are given in Table 5, accounting for 3 different cases. The values have been chosen to get noticeably different results among the proposed cases and to challenge the robustness of the solvers.

Each group is requested to compute the displacement field at points $\mathbf{p}_0, \mathbf{p}_1$ for each case.

Setup	a	a_f	a_{fs}	a_s	σ_0
Case A	177	55416	648	7443	2×10^5
Case B	295	92360	1080	12405	1×10^5
Case C	19	6157	72	827	2×10^5

Table 5: Each case combines a change in stiffness parameters a, a_f, a_{fs}, a_s with changes in the contractibility parameter σ_0 .

4 Benchmark 2: biventricular mechanics (blinded)

4.1 Strong formulation

Let us consider an idealized biventricular domain $\Omega \subset \mathbb{R}^3$ with boundaries $\partial\Omega := \Gamma_{endo-lv} \cup \Gamma_{endo-rv} \cup \Gamma_{epi} \cup \Gamma_{top}$. We denote by $\mathbf{u} : \Omega \rightarrow \mathbb{R}^3$ the displacement field, $\mathbb{T} := \mathbb{T}(\mathbf{u})$ the stress tensor as in Section 2.2 and $p_{lv}(t), p_{rv}(t)$ for each $t > 0$ pressure terms solving (7), with parameters to be specified below. We define the remaining operators \mathbb{F}, J as in Section 2. The problem to solve is described by the equations:

$$\begin{aligned}
\rho \ddot{\mathbf{u}} - \text{Div}(J\mathbb{T}\mathbb{F}^{-\top}) &= \mathbf{0} && \text{in } \Omega \\
J\mathbb{T}\mathbb{F}^{-\top}\mathbf{N} &= p_{lv}J\mathbb{F}^{-\top}\mathbf{N} && \text{on } \Gamma_{endo-lv} \\
J\mathbb{T}\mathbb{F}^{-\top}\mathbf{N} &= p_{rv}J\mathbb{F}^{-\top}\mathbf{N} && \text{on } \Gamma_{endo-rv} \\
J\mathbb{T}\mathbb{F}^{-\top}\mathbf{N} \cdot \mathbf{N} + \alpha_{epi}\mathbf{u} \cdot \mathbf{N} + \beta_{epi}\dot{\mathbf{u}} \cdot \mathbf{N} &= 0 && \text{on } \Gamma_{epi} \\
J\mathbb{T}(\mathbb{F}^{-\top}\mathbf{N}) \times \mathbf{N} &= \mathbf{0} && \text{on } \Gamma_{epi} \\
J\mathbb{T}\mathbb{F}^{-\top}\mathbf{N} + \alpha_{top}\mathbf{u} + \beta_{top}\dot{\mathbf{u}} &= \mathbf{0} && \text{on } \Gamma_{top}
\end{aligned} \tag{17}$$

4.2 Geometry

To define the geometry we will introduce some notation. Given $\mathbf{x}_{cen} \in \mathbb{R}^3$ and $\{a, b, c\} \in \mathbb{R}_+$, we define $V(\mathbf{x}_{cen}, (a, b, c)) \subset \mathbb{R}^3$ an ellipsoidal domain, centered at \mathbf{x}_{cen} with (a, b, c) the length of each $(\hat{x}, \hat{y}, \hat{z})$ semiaxis, $\partial V(\mathbf{x}_{cen}, (a, b, c))$ its boundary.

The biventricular domain $\Omega \subset \mathbb{R}^3$ is characterized by four different surfaces:

- Epicardial surface (Γ_{epi}) described as a set of points $\mathbf{x} = (x, y, z) \in \mathbb{R}^3$ satisfying

$$\begin{aligned} \mathbf{x} \in \partial V(\mathbf{0}, (a_{lv-epi}, b_{lv-epi}, c_{lv-epi})) \Delta \partial V(\mathbf{x}_{rv}, (a_{rv-epi}, b_{rv-epi}, c_{rv-epi})) \\ \mathbf{x} \text{ s.t. } x < 0 \end{aligned} \quad (18)$$

for $\mathbf{x}_{rv} = (0, 0, 0.02)$ and centers $(a_{lv-epi}, b_{lv-epi}, c_{lv-epi}) = (0.08, 0.039, 0.039)$ and $(a_{rv-epi}, b_{rv-epi}, c_{rv-epi}) = (0.075, 0.038, 0.059)$.

- Left endocardial surface ($\Gamma_{endo-lv}$) described as the set $\mathbf{x} = (x, y, z) \in \mathbb{R}^3$ satisfying

$$\begin{aligned} \mathbf{x} \in \partial V(\mathbf{0}, (a_{lv-endo}, b_{lv-endo}, c_{lv-endo})) \\ \mathbf{x} \text{ s.t. } x < 0 \end{aligned} \quad (19)$$

for $(a_{lv-endo}, b_{lv-endo}, c_{lv-endo}) = (0.069, 0.025, 0.025)$.

- Right endocardial surface ($\Gamma_{endo-rv}$) described as the set $\mathbf{x} = (x, y, z) \in \mathbb{R}^3$ satisfying

$$\begin{aligned} \mathbf{x} \in \partial V(\mathbf{x}_{rv}, (a_{rv-endo}, b_{rv-endo}, c_{rv-endo})) \\ \mathbf{x} \text{ s.t. } x < 0 \end{aligned} \quad (20)$$

for $(a_{rv-endo}, b_{rv-endo}, c_{rv-endo}) = (0.07, 0.033, 0.054)$.

- Base (Γ_{top}) as the set $\mathbf{x} = (x, y, z) \in \bar{\Gamma}_{epi} \cup \bar{\Gamma}_{endo-lv} \cup \bar{\Gamma}_{endo-rv}$ s.t. $x = 0$

The proposed geometry is depicted in Figure 4.

4.3 Fibers

For the fiber directions, we use a Laplace-Dirichlet Rule-Based (BT-LDRB) algorithm [34], modified to adhere to the convention utilized for the cross-fiber orientations[‡]. We take a values of $\mp 60^\circ$ (with respect to a local coordinate system) for the left and right endo/epicardial fiber angles[§].

The fibers are created using the `lifex` software [41, 42]. Figure 5 depicts the step-by-step procedure to prescribe the fiber architecture in the biventricular geometry [34, 41]. For further details refer to [34].

[‡]In the last decades, myocardial orientation has been studied from histological data [18, 35] and *Diffusion Tensor Imaging* [36, 37], but their reconstructed noisy data suffers from low resolution, limiting its characterization, especially given the thickness of ventricles, which is usually smaller than the voxel size [38]. Several construction algorithms have been proposed to recreate the fiber orientation, ranging from complex registration data-dependent algorithms to *Rule-Based Methods*, which remains an active area of research [34, 39, 40].

[§]The convention in this work entails switching the directions \mathbf{s} and \mathbf{n} in relation to the formalism entailed in the state-of-the-art [34].

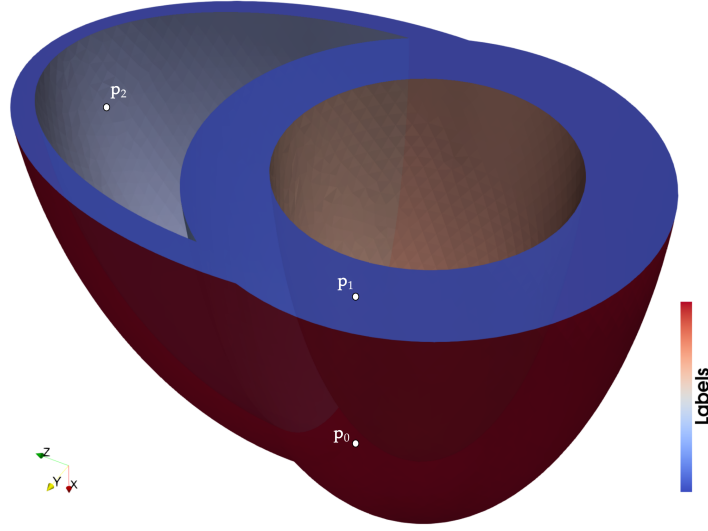


Figure 4: Geometry for the biventricular domain with colored boundaries: Γ_{epi} , $\Gamma_{endo-lv}$, $\Gamma_{endo-rv}$ and Γ_{top} . Positions of particles of interest \mathbf{p}_0 , \mathbf{p}_1 and \mathbf{p}_2 are depicted with circles for reference.

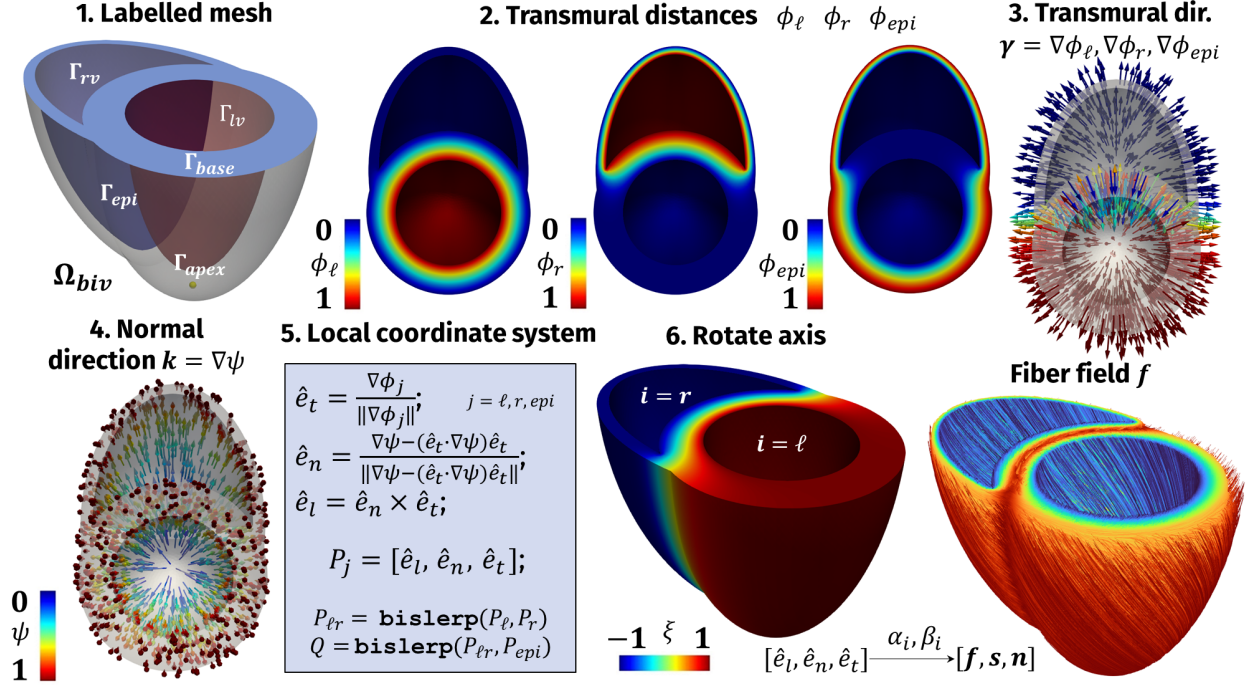


Figure 5: Step-by-step procedure for the fiber architecture. In 1. labelled mesh with boundaries, 2. transmural distances $\phi_\ell, \phi_r, \phi_{epi}$, 3. transmural directions $\gamma = \nabla\phi_\ell, \nabla\phi_r, \nabla\phi_{epi}$, 4. normal direction $k = \nabla\psi$, 5. local coordinate definition $\hat{e}_t, \hat{e}_n, \hat{e}_l$, 6. rotation of axis and fiber field system $[f, s, n]$

4.4 Physical constants and evaluation of results

Each group is requested to solve problem (17) with parameters as in Tables 2, 7, 8, 9, with geometry and fibers as in Sections 4.2, 4.3, for two refinement levels[‡]. Table 6 details the number of tetrahedra and nodes:

Mesh	Num. of tetrahedra	Num. of nodes
Ω_{h_1}	45,304	11,444
Ω_{h_2}	121,133	27,807

Table 6: Number of tetrahedra and nodes for two refinement levels, denoted by Ω_{h_1} and Ω_{h_2}

Extending Step 2 of Benchmark 1, groups are requested to provide displacement fields $\mathbf{u}_h(\mathbf{X})$ over time at three spatial locations $\mathbf{p}_0 = (0.025, 0.03, 0)$, $\mathbf{p}_1 = (0, 0.03, 0)$, $\mathbf{p}_2 = (0.025, 0, 0.072)$.

Parameter	$\rho[\frac{kg}{m^3}]$	$\eta[Pa\ s]$	$\kappa[Pa]$	$k[-]$	$\alpha_{top}[\frac{Pa}{m}]$	$\alpha_{epi}[\frac{Pa}{m}]$	$\beta_{top}[Pa\frac{s}{m}]$	$\beta_{epi}[Pa\frac{s}{m}]$
Value	10^3	10^2	10^6	100	10^6	10^8	5×10^3	5×10^3

Table 7: Parameters describing the strong form of the problem defined in (17).

Parameter	$\sigma_0[Pa]$	$\gamma[s]$	α_{\min}	α_{\max}	$t_{\text{sys}}[s]$	$t_{\text{dias}}[s]$
Value	1.5×10^5	0.005	-30	5	0.163	0.5

Table 8: Parameters defining the active stress activation function, solution to (5), for the biventricular model.

[‡]The repository cardiac-benchmark-toolkit stores the biventricular domain in several formats `.geo`, `.msh`, `.xdmf`, `.h5`.

Parameter (p_{lv})	Value	Parameter (p_{rv})	Value
$\alpha_{min}[\cdot]$	-30	$\alpha_{min}[\cdot]$	-30
$\alpha_{max}[\cdot]$	5	$\alpha_{max}[\cdot]$	5
$\alpha_{pre}[\cdot]$	5	$\alpha_{pre}[\cdot]$	1
$\alpha_{mid}[\cdot]$	15	$\alpha_{mid}[\cdot]$	10
$\sigma_{pre}[Pa]$	12000	$\sigma_{pre}[Pa]$	3000
$\sigma_{mid}[Pa]$	16000	$\sigma_{mid}[Pa]$	4000
$t_{sys-pre}[s]$	0.17	$t_{sys-pre}[s]$	0.17
$t_{dias-pre}[s]$	0.484	$t_{dias-pre}[s]$	0.484
$\gamma[s]$	0.005	$\gamma[s]$	0.005

Table 9: Left table: parameters used for $p_{lv}(t)$, so that it attains a maximum of $16491.14 [Pa] \approx 123 [mmHg]$. Right: parameters used for $p_{rv}(t)$ with a maximum of $4166.66 [Pa] \approx 31 [mmHg]$.

Depicted in Figures 6 and 7 are the time-evolution of the activation function and pressure curves with parameters specified therein.

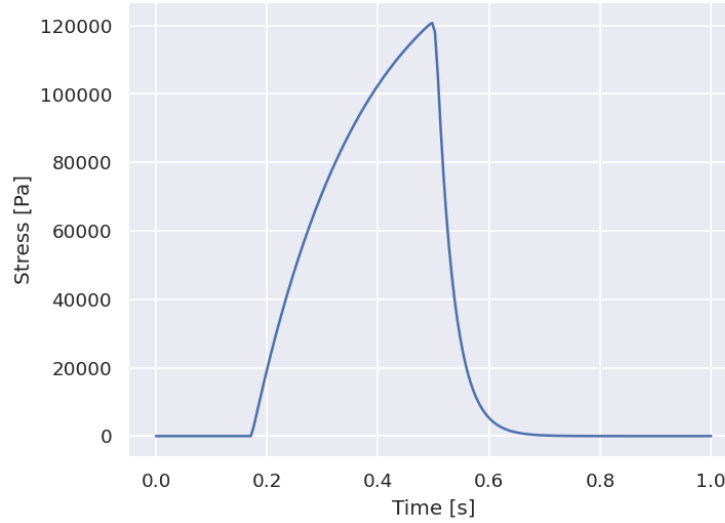


Figure 6: Evolution of the stress function τ described in (5) over the time interval $[0, 1]$ with parameters as in Table 8. It reaches a maximum value of $120775.56 [Pa]$.

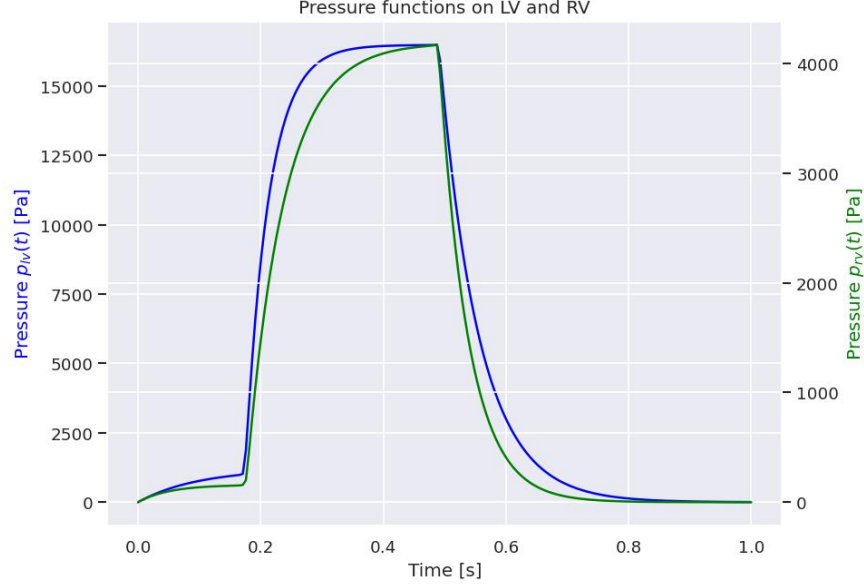


Figure 7: Evolution of pressure $p_{lv}(t), p_{rv}(t)$ are shown with blue and brown colors respectively over the time interval $[0, 1]$. Parameters as in Table 4. Maximum values of $16491.15 [Pa]$ and $4171.07 [Pa]$ for p_{lv}, p_{rv} respectively.

5 Numerical solvers and participants

Each group was requested to disclose their strategies to solve problems (1) and (17). Settings for the software, spatial and temporal discretization methods are described in Table 10. The notation \mathbb{P}_2 indicates that the incompressibility is handled via penalization (as described in the previous sections), and therefore only the displacements are discretized with quadratic basis functions. The notation \mathbb{P}_1 indicated the analogous case for linear basis functions. The notation $\mathbb{P}_1 - \mathbb{P}_1$ indicates that incompressibility is handled directly using the pressure as unknown, where the saddle-point problem is discretized with linear basis functions, including a stabilization term for the pressure field. This variable, defined in the muscle, differs from the pressure prescribed at the boundaries, describing chamber and epicardial effects.

Affiliation	Software	Spatial Discretization	Quad. rule and degree	Linear Solver	Time Discretization
Medical University of Graz	CARPentry [43]	\mathbb{P}_1 - \mathbb{P}_1 [44]	KL - 6	GMRES [45] with precondition	KGen- α with spectral radius $\rho_\infty = 0.5$, timestep at 1[ms]
King's College London	Ambit [46] (FEniCSx)	\mathbb{P}_2	GL - 6	LU [47] for bench. 1 and GMRES with AMG precondition. [48] for bench. 2	Gen- α with $\alpha_m = \alpha_f = 0$, $\beta = 0.25$, $\gamma = 0.5$, timestep at 1[ms]
Technische Universität München	4C [49]	\mathbb{P}_2	GL - 4 (stiffness), 11 (mass)	GMRES with AMG precondition.	Gen- α with $\alpha_m = \alpha_f = 0.5$, $\beta = 0.25$, $\gamma = 0.5$, timestep at 1[ms].
Simula Research	FEniCS [50]	\mathbb{P}_2	GL - 6	LU [51]	Gen- α with $\alpha_m = 0.2$, $\alpha_f = 0.4$, timestep at 1[ms].
University of Groningen	CHIMeRA (FEniCS)	\mathbb{P}_2	GL - 6	LU [47]	Gen- α with $\alpha_m = \alpha_f = 0$, $\beta = 0.25$, $\gamma = 0.5$, timestep at 1[ms].
University of Michigan	CHeart [52]	\mathbb{P}_2	KL - 4	LU [47]	Mid-point rule, timestep at 1[ms]
Politecnico di Milano	life ^x [41]	\mathbb{P}_2	GL - 4	GMRES with AMG precondition.	BDF1 Implicit, timestep at 1[ms].
Columbia University	SimVascular svFSI [53, 54]	\mathbb{P}_1 - \mathbb{P}_1	GL - 4	GMRES with Schwarz precondition. for bench. 1 and GMRES with iLU [55] for bench. 2	KGen- α with spectral radius $\rho_\infty = 0.5$, timestep at 1[ms]
Technische Universiteit Delft	COMSOL Multiphysics v.6.1 [56] [11, 57]	\mathbb{P}_2	GL - 4	LU [58]	Gen- α with $\alpha_m = \alpha_f = 0$, $\beta = 0.25$, $\gamma = 0.5$, timestep at 1[ms]

Table 10: Summary table of strategies. All problems are solved using the Newton method. Notation: Gen- α denotes the generalized- α scheme [59] and KGen- α an subclass of Gen- α with specific spectral properties [60]. GL denotes the Gauss-Legendre and KL the Keast Lyness quadrature rules, respectively. FEniCSx [61–63] is the successor of FEniCS [64–66]. In the linear solver column, teams differ between linear algebra backend such as MUMPS [47], SuperLU [51], PETSc [48], Trilinos [55] and PARDISO [58], to solve the linear system of the newton step.

6 Results

In this section, we provide comparison results for the two benchmark problems. Quantitative and qualitative assessment is done using displacement tracking and by defining a measure of discrepancy between teams. Solutions from different teams can be distinguished with different colors, which are provided in displacement curves for benchmark 1 and including visualizations for benchmark 2. Comparisons between \mathbb{P}_1 and \mathbb{P}_2 are also provided for benchmark 2.

6.1 Benchmark 1

6.1.1 Step 0 (non-blinded): Splitting passive and active response

The comparison of displacement curves at particles $\mathbf{p}_0, \mathbf{p}_1$ is depicted in Figure 8 for active response alone and in Figure 9 for passive response. Each figure presents displacements for each of the component, allowing a straightforward assessment of differences between teams. The largest differences can be observed primarily along the interval $(0.2, 0.6)$ [s], and especially for the x-component of both particles in the case of passive response.

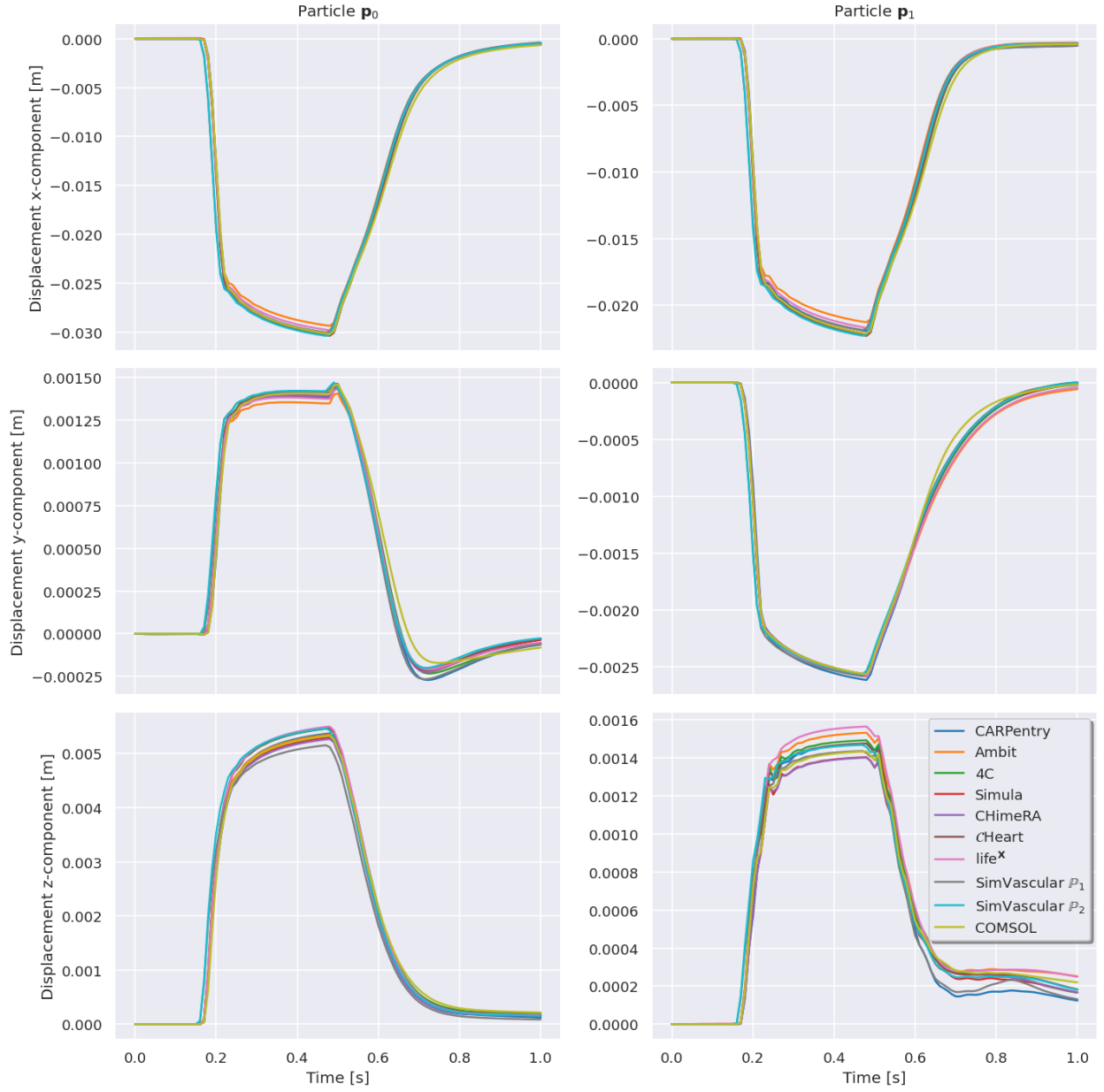


Figure 8: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$ for Step 0, case A - active response.

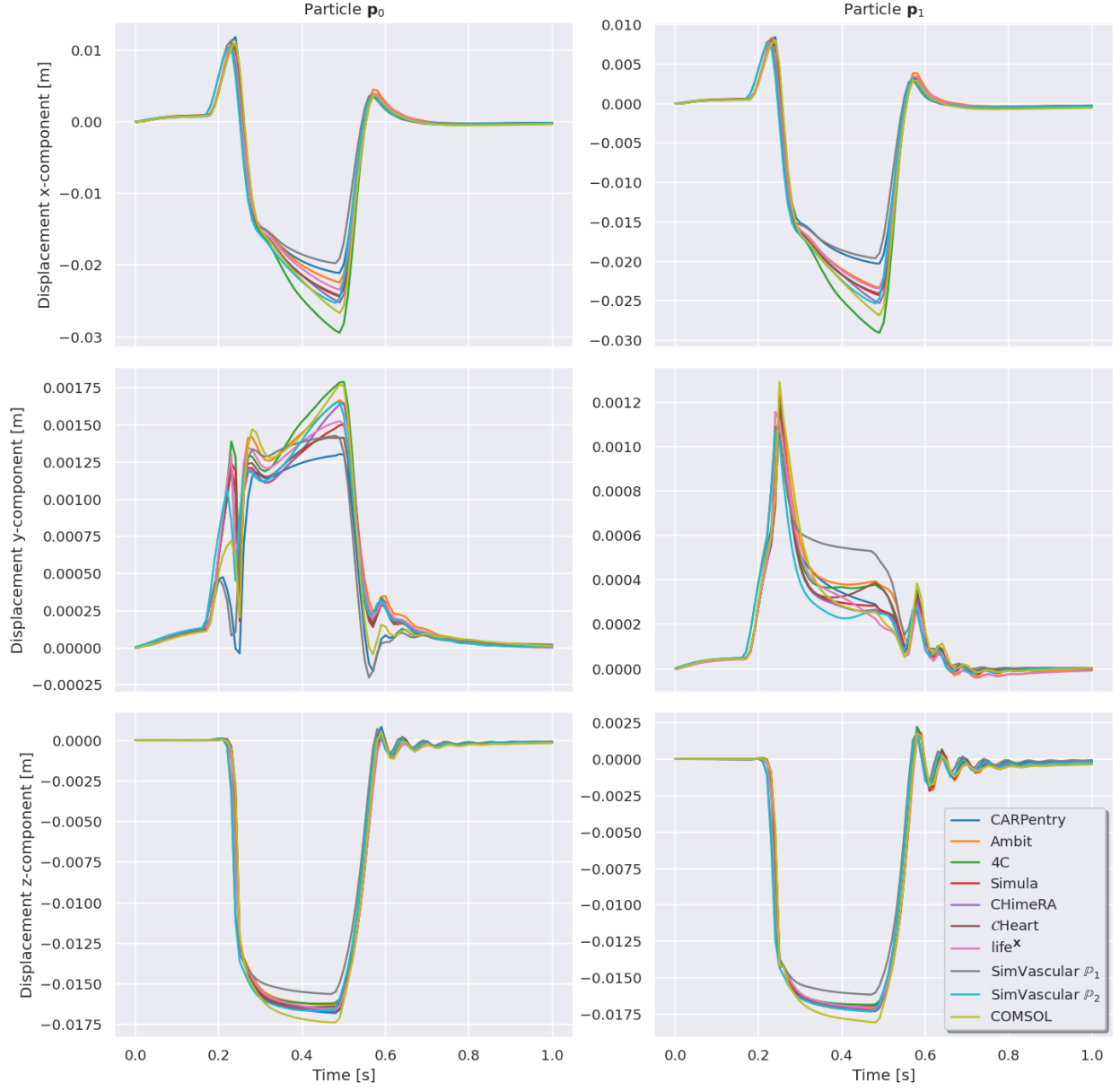


Figure 9: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$ for Step 0, case B - passive response.

6.1.2 Step 1 (non-blinded): active and passive response

The comparison results for each requested quantity are depicted in Figure 10. The componentwise representation of displacement showcases the differences in the order of magnitude of deformation. Displacements along the z-component are one order of magnitude smaller than those of the x-component. Maximum differences between teams remain smaller than $0.5 [mm]$ in the worst case, as seen in the z-component.

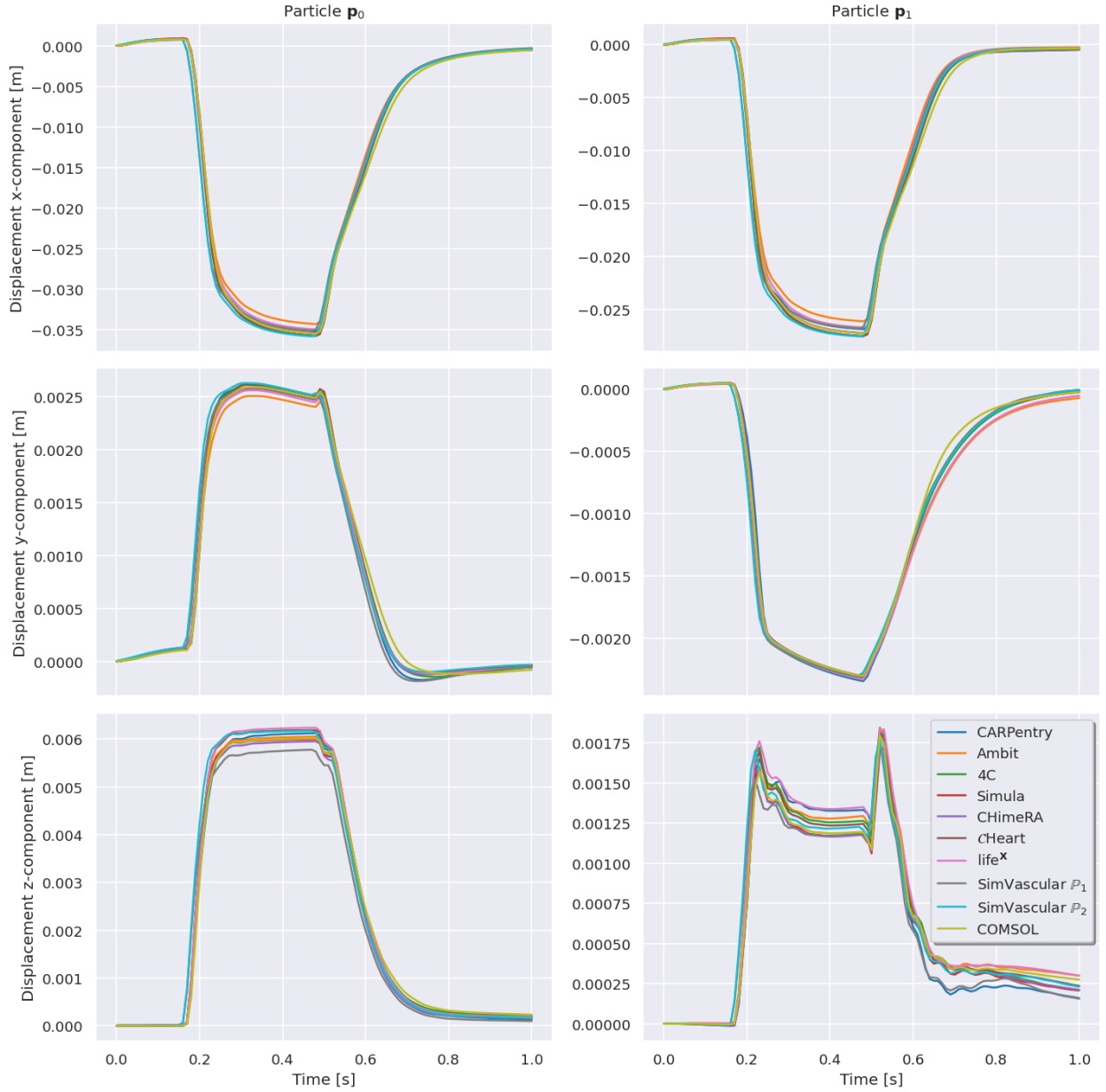


Figure 10: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$ for Step 1, case of joint active and passive response.

6.1.3 Step 2

A qualitative comparison of displacement curves at different particles is depicted in Figures 11-13. For a quantitative assessment of the curves, we propose a [RE]lative [D]iscrepancy between each dataset, denoted by RED , defined as:

$$\text{RED}(\mathbf{p}) = \frac{1}{T} \sum_{t_n=0}^T \frac{\|\mathbf{u}(t_n, \mathbf{p}) - \bar{\mathbf{u}}(t_n, \mathbf{p})\|_{\ell^2}}{\|\bar{\mathbf{u}}(t_n, \mathbf{p})\|_{\ell^2}} \quad \mathbf{p} \in \{\mathbf{p}_0, \mathbf{p}_1\} \quad (21)$$

with $\bar{\mathbf{u}}(t, \cdot) = \frac{1}{N} \sum_{i=1}^N \mathbf{u}^i(t, \cdot)$ for $t \in (0, 1)$, \mathbf{u}^i the displacement field of team i and N the total number of teams. The datasets are subsampled at $10[ms]$, i.e. $T = 101$ datapoints. If a group simulated with a different timestep, then linear interpolation is used to compute the corresponding displacement values. Intuitively, the relative discrepancy function provides a time-averaged discrepancy to an average result, using the ℓ^2 norm to add each direction. Table 11 summarizes the relative discrepancies for each team.

Setup	Case A		Case B		Case C	
	$\text{RED}(\mathbf{p}_0)$	$\text{RED}(\mathbf{p}_1)$	$\text{RED}(\mathbf{p}_0)$	$\text{RED}(\mathbf{p}_1)$	$\text{RED}(\mathbf{p}_0)$	$\text{RED}(\mathbf{p}_1)$
CARPentry	0.134	0.229	0.360	0.301	0.060	0.118
Ambit	0.115	0.185	0.79	0.243	0.060	0.168
4C	0.080	0.115	0.171	0.136	0.059	0.059
Simula	0.094	0.200	0.311	0.352	0.041	0.054
CHIMeRA	0.078	0.108	0.149	0.135	0.045	0.056
CHeart	0.202	0.198	0.300	0.250	0.048	0.045
life ^x	0.108	0.154	0.273	0.252	0.049	0.129
SimVascular \mathbb{P}_1	0.220	0.371	0.309	0.267	0.156	0.211
SimVascular \mathbb{P}_2	0.276	0.370	0.360	0.328	0.146	0.157
COMSOL	0.186	0.196	0.287	0.329	0.105	0.111

Table 11: Comparison of relative deviations for each participant group.

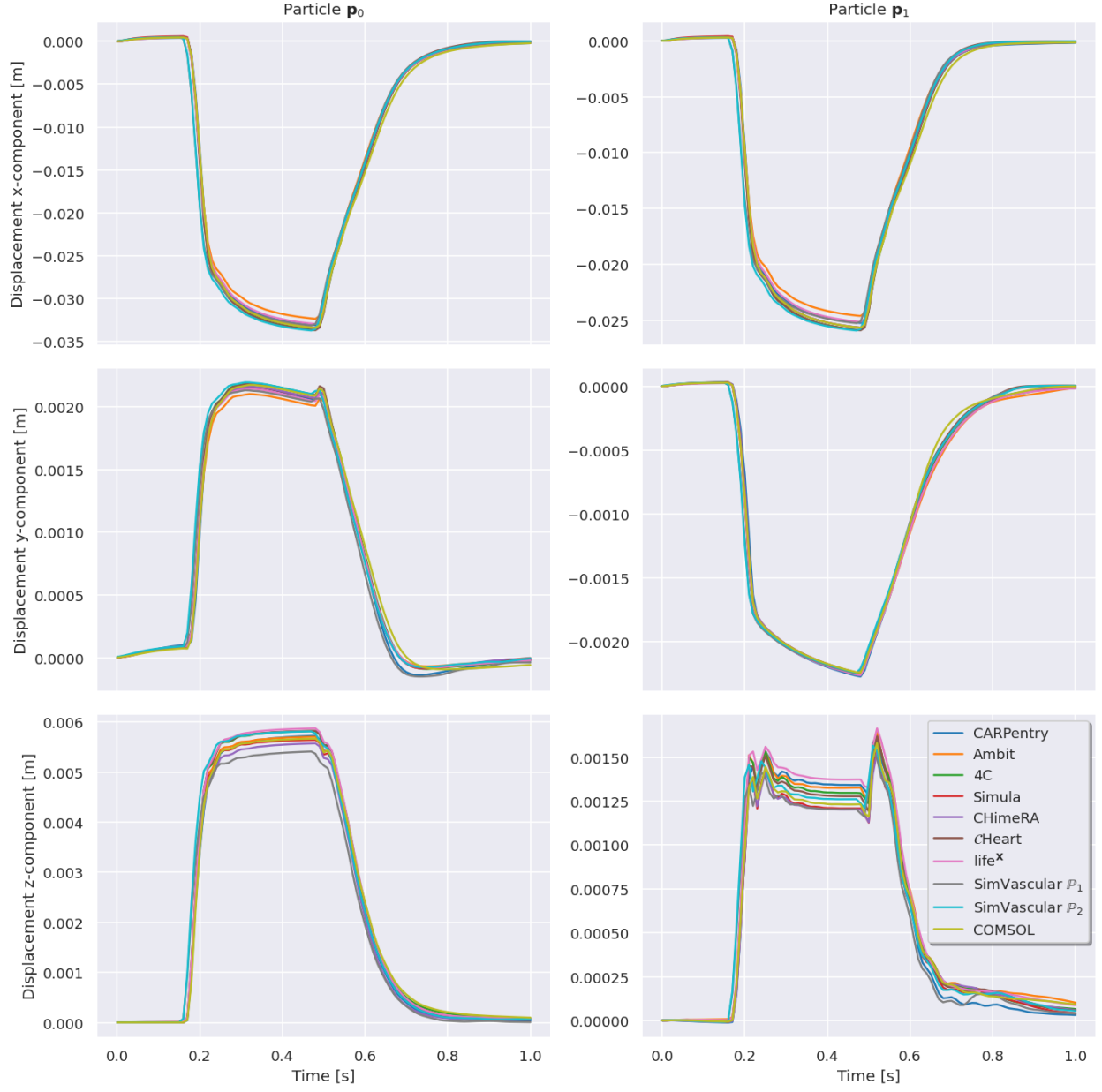


Figure 11: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$, case A.

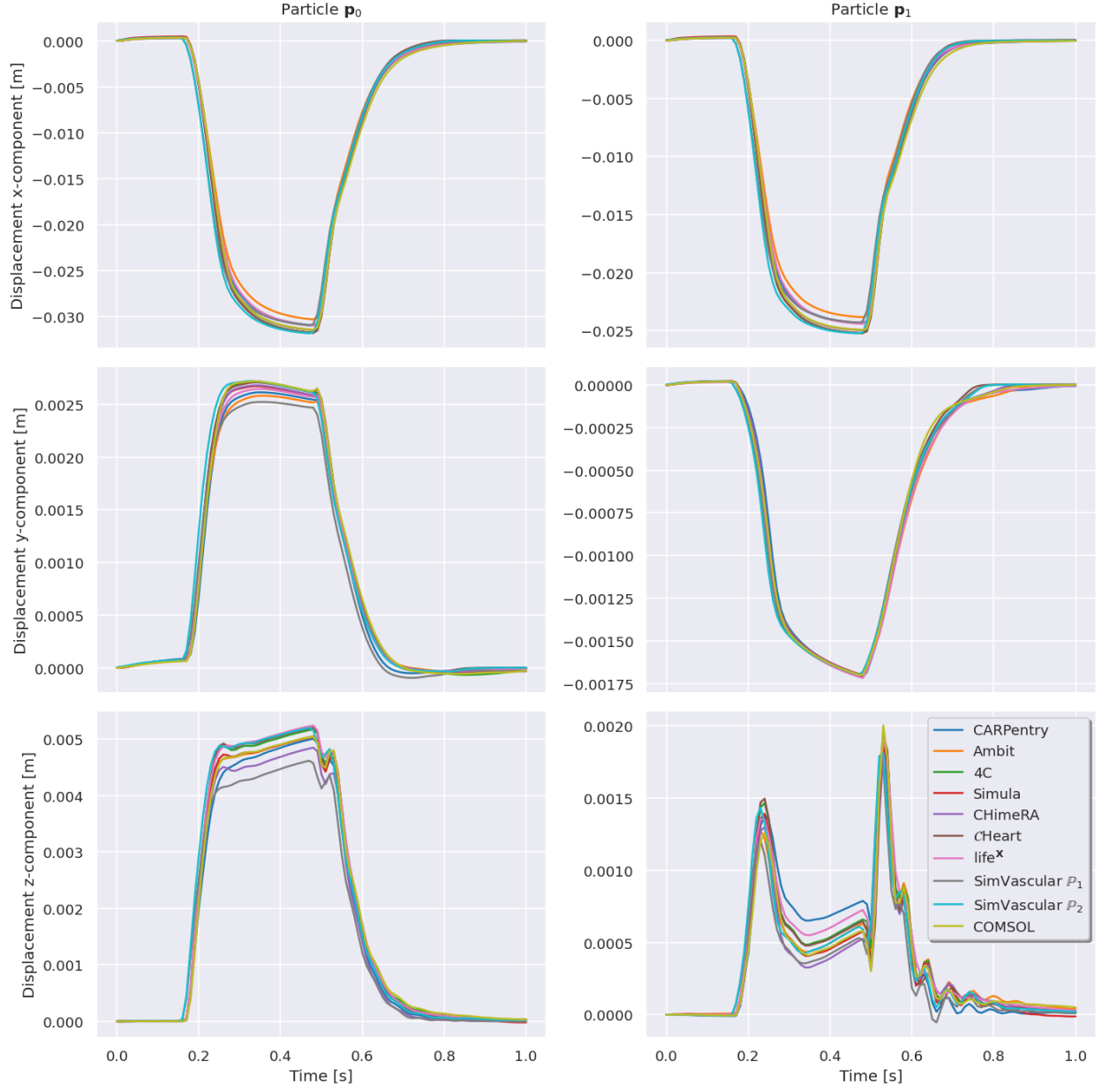


Figure 12: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$, case B.

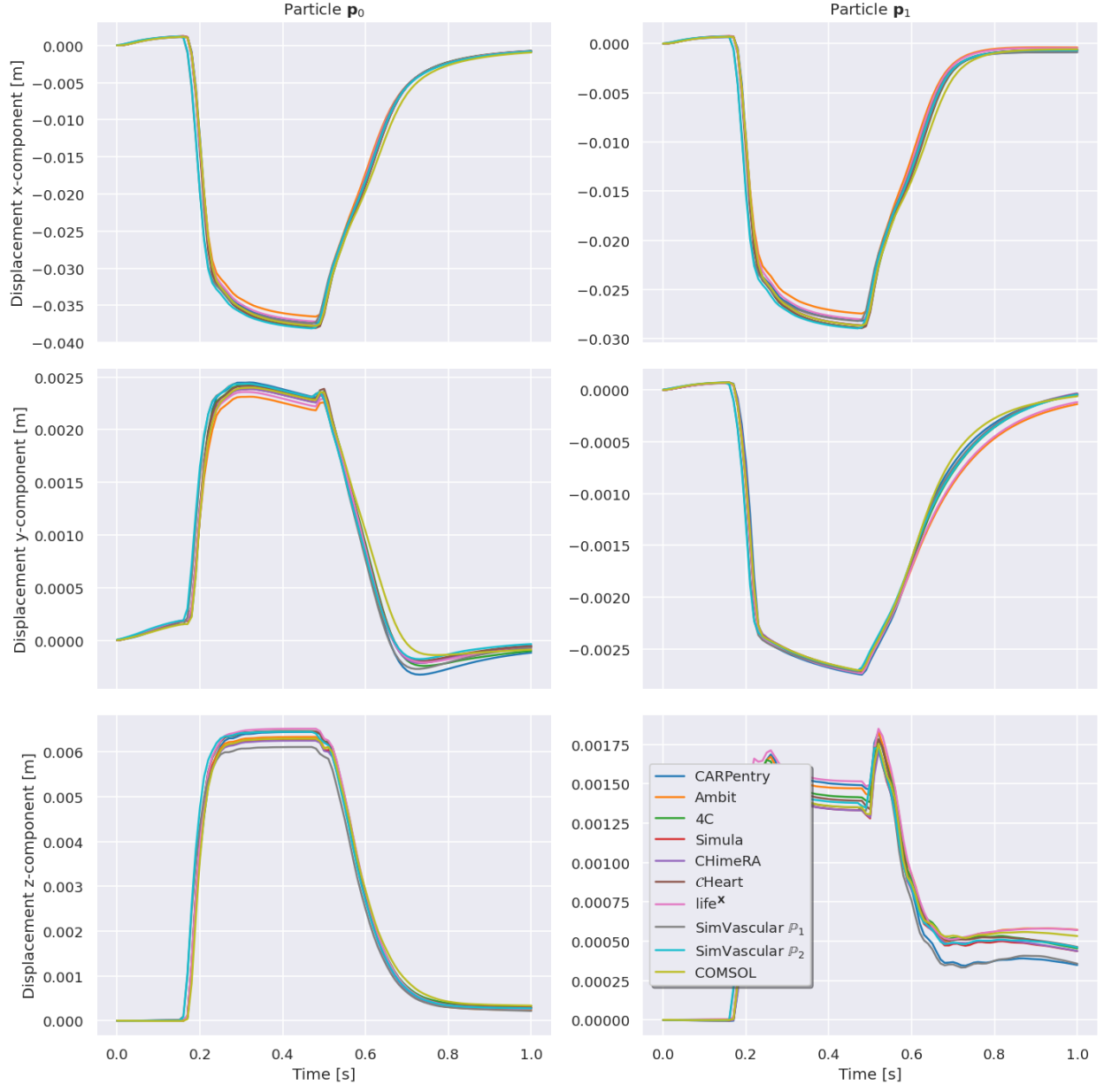


Figure 13: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ and $\mathbf{u}_h(\mathbf{p}_1)$, case C.

6.2 Benchmark 2

To analyze the results, qualitative assessment is done through visual inspection and the displacement tracking at three particles \mathbf{p}_0 , \mathbf{p}_1 and \mathbf{p}_2 . We provide quantitative assessment using the measure of discrepancy **RED**, as in Section 6.1.3, for all particles in both meshes. Visual comparison between solutions can be depicted in Figure 14 using overlapped views at two different times, namely, 0.3[s] and 0.5[s]. The views are defined using the two-chamber (long) axis and the base-to-apex (short) axis. Particle trajectories are depicted in Figures 15 for the coarse mesh Ω_{h_1} and 16 for the fine mesh Ω_{h_2} . Table 12 summarizes the discrepancies in each case. Comparison curves between spatial discretization in \mathbb{P}_1 and \mathbb{P}_2 are depicted in Figures 17 and 18, including only teams that provided both datasets.

Setup	Blinded on Ω_{h_1}			Blinded on Ω_{h_2}		
	RED(\mathbf{p}_0)	RED(\mathbf{p}_1)	RED(\mathbf{p}_2)	RED(\mathbf{p}_0)	RED(\mathbf{p}_1)	RED(\mathbf{p}_2)
CARPentry	0.915	0.545	0.415	1.019	0.504	0.452
Ambit	0.094	0.086	0.21	0.136	0.084	0.288
4C	0.104	0.129	0.221	0.108	0.094	0.278
Simula	0.564	0.848	1.472	0.446	0.513	1.769
CHIMeRA	0.121	0.108	0.182	0.111	0.079	0.347
CHeart	0.144	0.11	0.226	0.137	0.085	0.406
life ^x	0.125	0.099	0.144	0.103	0.077	0.318
SimVascular \mathbb{P}_1	0.483	0.295	0.95	0.294	0.184	0.508
COMSOL	0.14	0.158	0.335	0.183	0.155	0.326

Table 12: Comparison of relative deviations for each participant group in Benchmark 2.

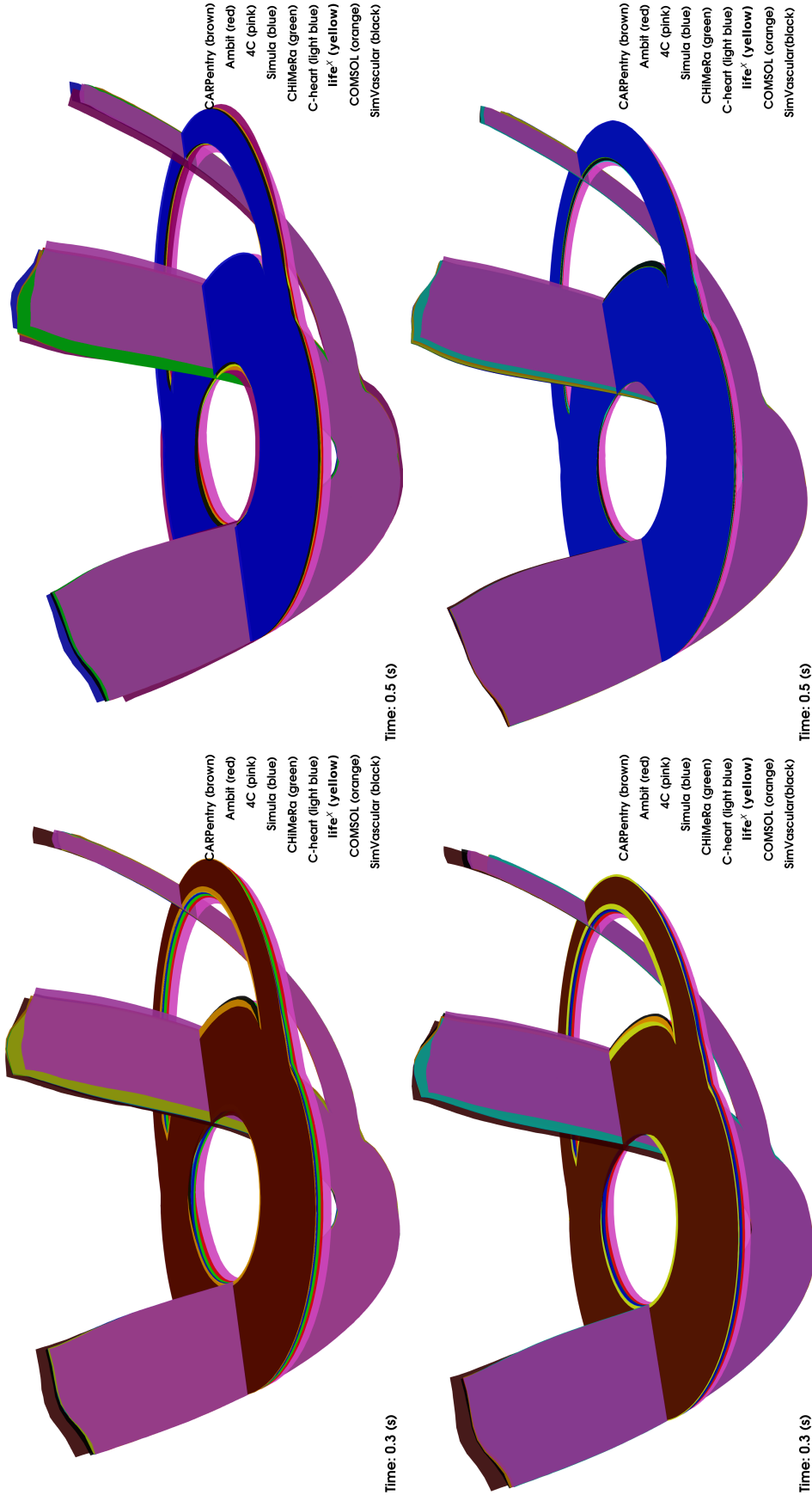


Figure 14: Visual overlapping of each team solution using two-chamber and short-axis views, at two different time instants $t = 0.3[s]$ (left column) and $t = 0.5[s]$ (right column) in the coarse mesh Ω_{h_1} (top row) and fine mesh Ω_{h_2} (bottom row).

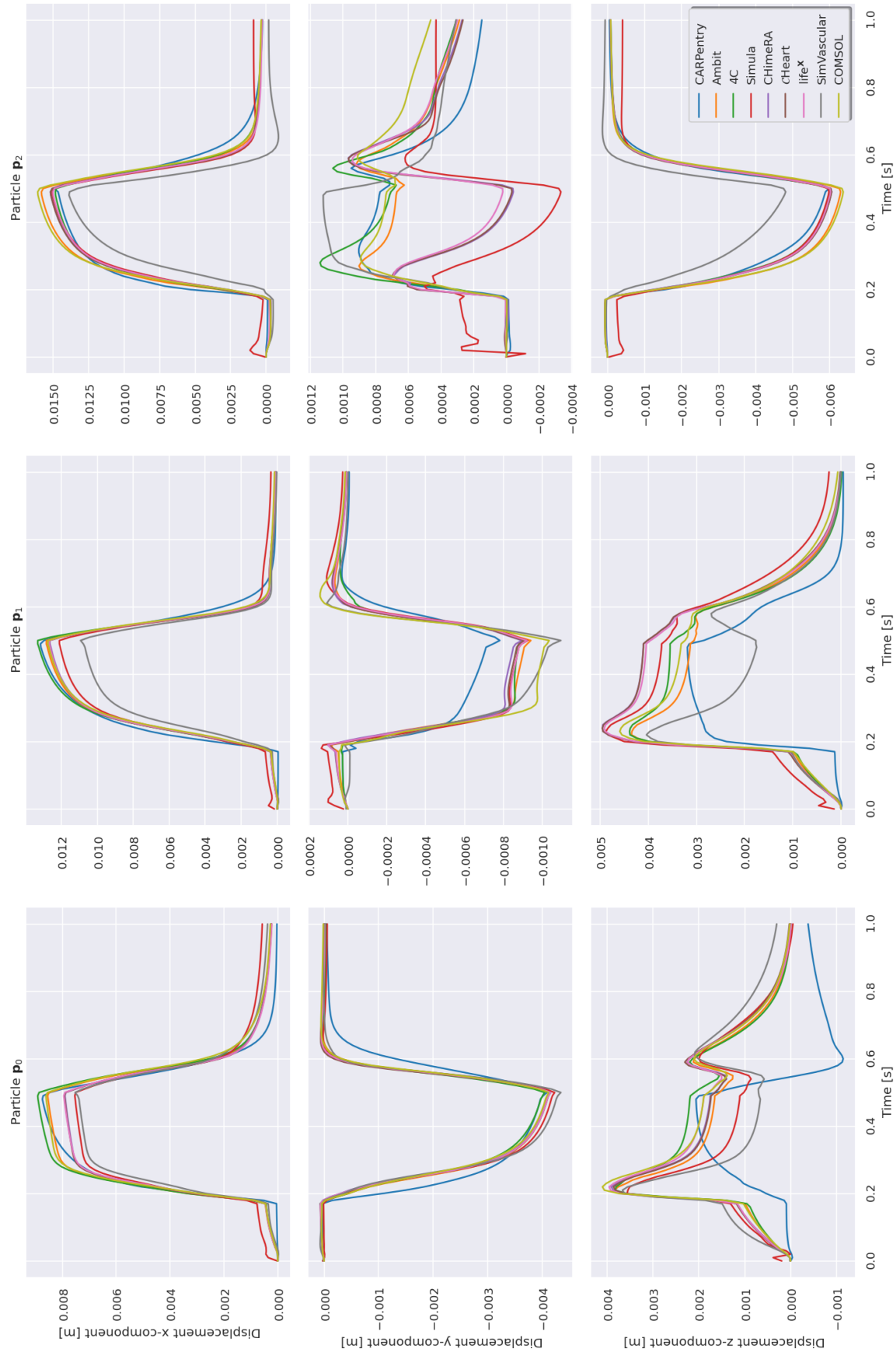


Figure 15: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ (left), $\mathbf{u}_h(\mathbf{p}_1)$ (center) and $\mathbf{u}_h(\mathbf{p}_2)$ (right) in the coarse mesh Ω_{h_1} .

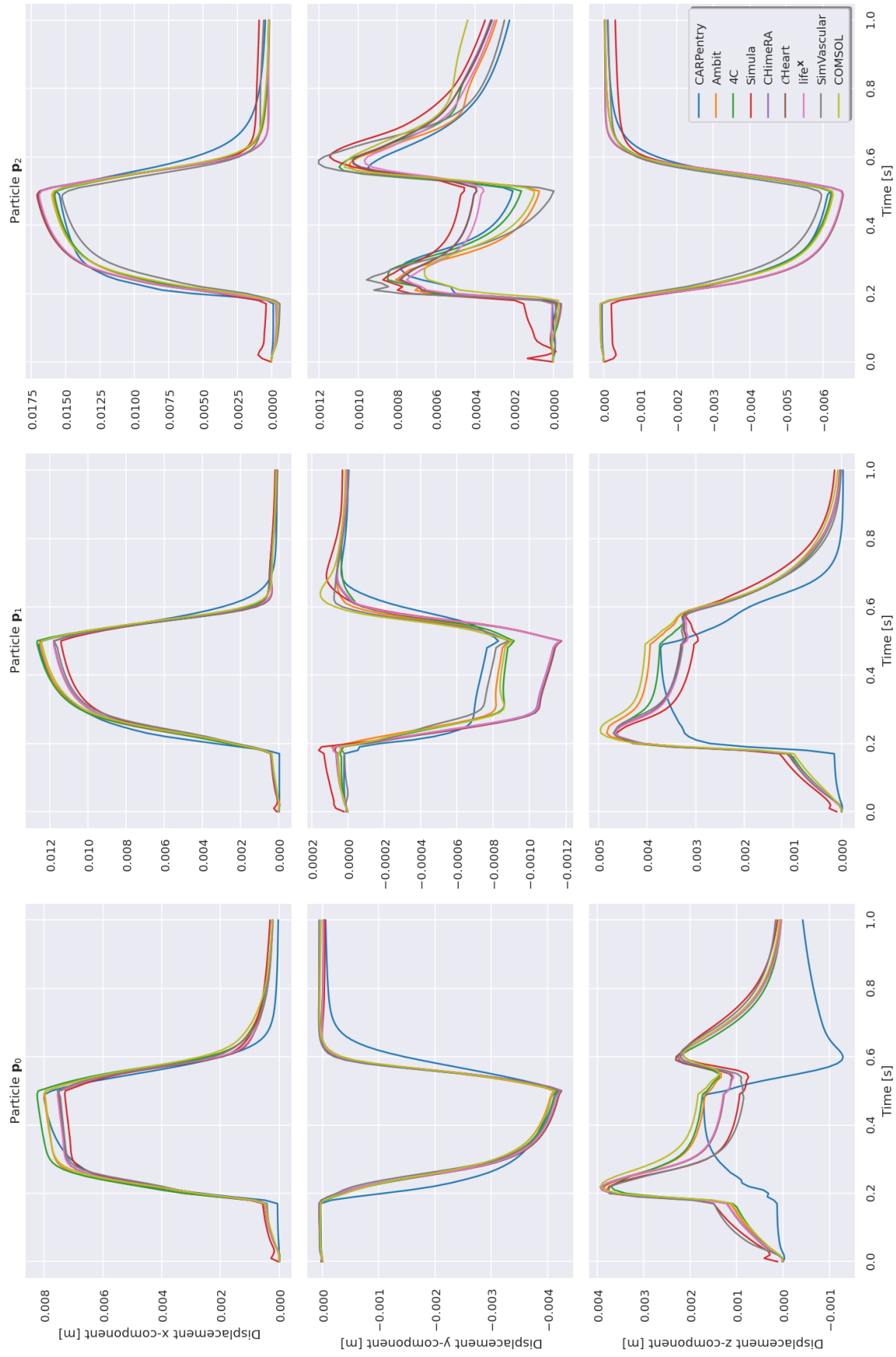


Figure 16: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ (left), $\mathbf{u}_h(\mathbf{p}_1)$ (center) and $\mathbf{u}_h(\mathbf{p}_2)$ (right) in the fine mesh Ω_{h_2} .

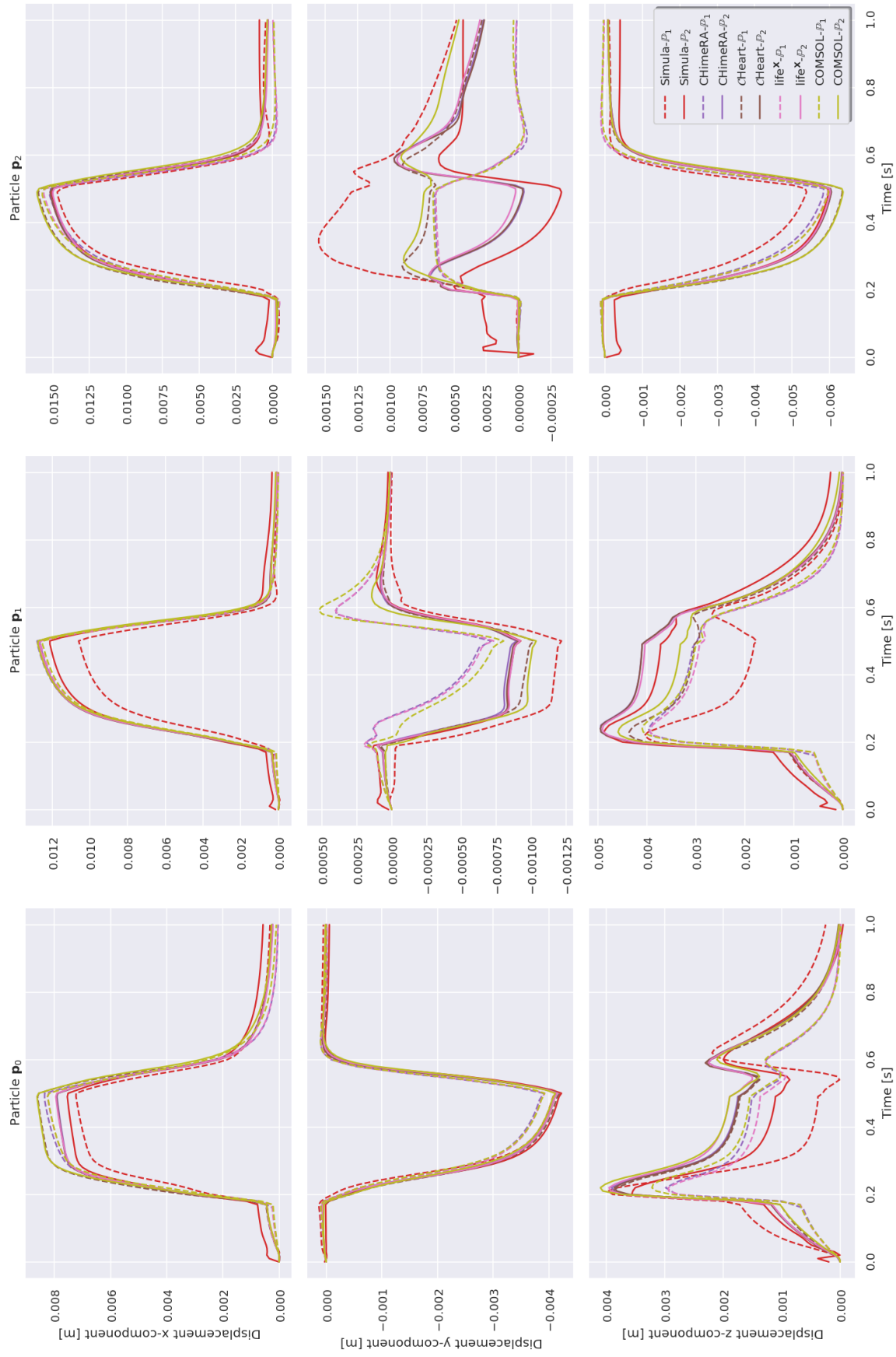


Figure 17: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ (left), $\mathbf{u}_h(\mathbf{p}_1)$ (center) and $\mathbf{u}_h(\mathbf{p}_2)$ (right) in the coarse mesh Ω_{h_1} . In dashed lines the \mathbb{P}_1 solutions and in full lines the \mathbb{P}_2 solutions.

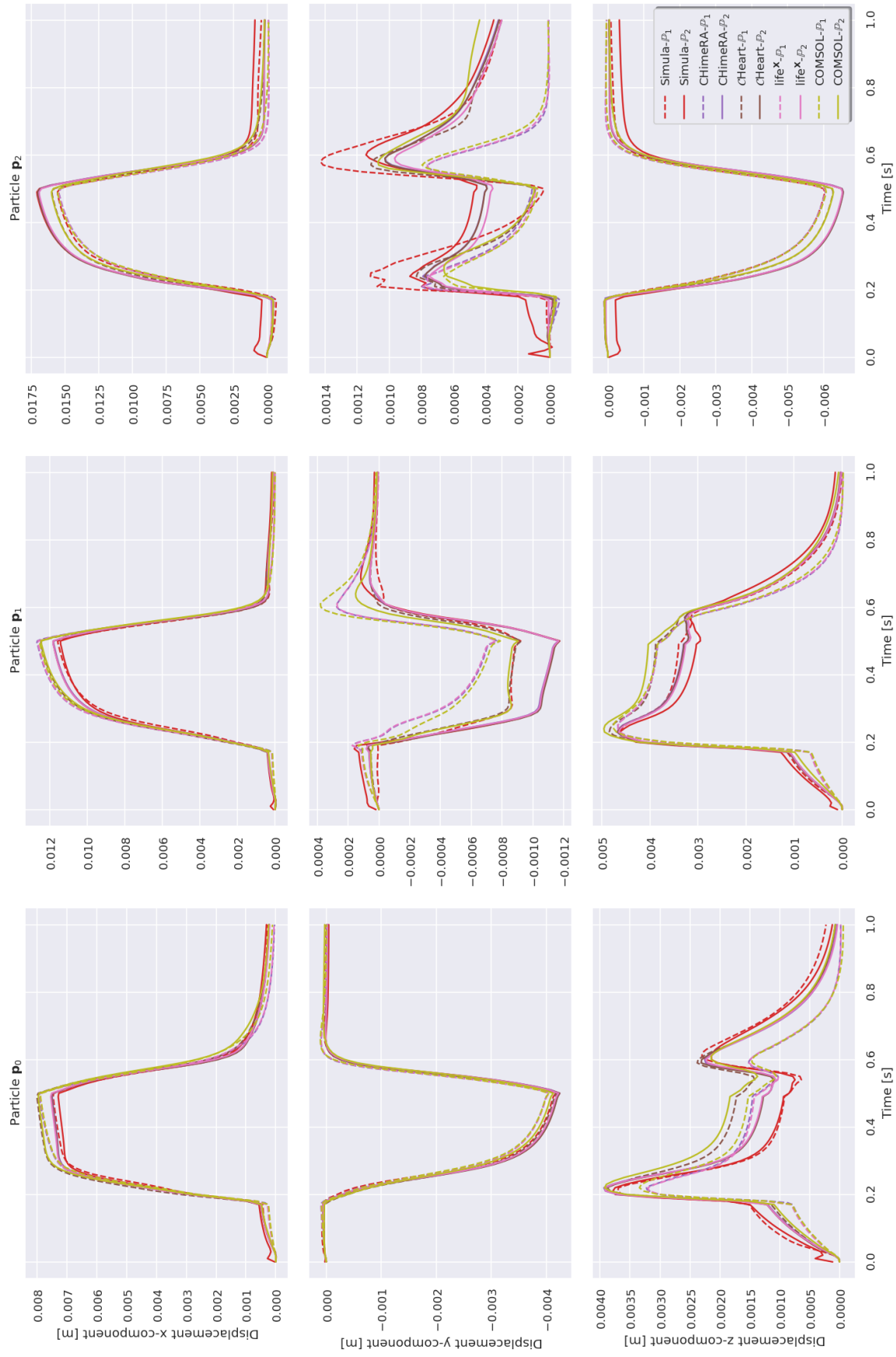


Figure 18: Comparison per component of displacement $\mathbf{u}_h(\mathbf{p}_0)$ (left), $\mathbf{u}_h(\mathbf{p}_1)$ (center) and $\mathbf{u}_h(\mathbf{p}_2)$ (right) in the fine mesh Ω_{h_2} . In dashed lines the \mathbb{P}_1 solutions and in full lines the \mathbb{P}_2 solutions.

7 Discussion

This work proposed a set of benchmark problems and solutions for cardiac elastodynamics, in both, monoventricular and biventricular geometries. In depth evaluation of the solutions is done using the discrepancy measure **RED**. For reproducibility and verification reasons, the results are available as supplemented material.

The benchmarks proposed here not only assess nonlinear elastodynamics but also test active material behavior and pericardial boundary conditions [7]. They also showcase the the variability of the numerical approaches used by the scientific community in cardiac biomechanics. This work not only provides the analytical description for the monoventricular case, as done in [2], but also utilizes the state-of-the-art fiber generation pipeline [34] for the biventricular domain.

This report provides an unambiguous mathematical description of cardiac benchmark problems, sufficient for reproducibility purposes with agreement of solutions between teams for all proposed problems. In total, nine different research groups submitted solutions to the benchmark problems. All computational setups relied on the finite-element method and three different approaches were considered, two by handling incompressibility via penalization with the displacement field unknown, discretized in \mathbb{P}_1 , \mathbb{P}_2 and incompressibility handled using stabilization with pressure and velocity (or displacement) as unknowns in $\mathbb{P}_1 - \mathbb{P}_1$.

The monoventricular benchmark case comprises three different problems that aim to assess passive 3.3.2 and active 3.3.1 responses of the cardiac contractility, as well as their combined effect 3.4. The splitting between each independent response and their combined effect allows for separated benchmarking of different model components.

In the non-blinded phase, teams had access to numerical solutions provided by other participants. In this phase, solutions agree closely on the active, passive and joint responses, as depicted in Figures 8-10. The difference between curves is below $0.5 [mm]$, only a small fraction of the typical element size employed ($3-5 [mm]$). The largest differences are observed when the discrete system is passively loaded as seen in Table 11.

In the blinded-phase, teams test their numerical setup in three new sets of parameters and assess agreement between solutions. The choice of parameters defines different material regimes, a high and low stiffness set of parameters (Case A and C, respectively), tuned to have physiological contraction and a third case tuned to have small deformations to test robustness of the solvers. In all cases, a reasonable agreement is observed among teams, as depicted in Figures 11-13. This agreement is present despite some groups (**SimVascular**, **CHIMeRA**, **CHeart** and **life^x**) using different ways to generate or interpolate the fiber directions, as summarized in Table 10.

The blinded biventricular case 4.4, aims to assess each solver in a more realistic scenario with a non-parametric fiber configuration and a generalized elastodynamic formulation. As depicted in Figures 15 and 16, there is a closer match between results when using the finer mesh Ω_{h_2} compared to Ω_{h_1} . A qualitative comparison of the solutions along the short and long axis planes is depicted in Figure 14. Greater discrepancy is observed between solutions based on $\mathbb{P}_1 - \mathbb{P}_1$ compared to \mathbb{P}_2 formulations. Discrepancies are noticeable across all particle displacements in the coarser mesh, particularly for the particle **p₂**, reflecting the variability of

the solutions across the mesh. Effects of the spatial discretization are also considered in this work. Comparisons between solution fields in \mathbb{P}_1 and \mathbb{P}_2 , depicted in Figures 17 and 18 for the coarse and fine mesh, showcase a dependency of the solution on the discretization space and to the fibers orientation, with differences larger than $5[mm]$ in the interval $(0.2, 0.5) [s]$.

A close agreement between most of the groups in the monoventricular case (Table 11) using different software and methods. However, in the biventricular case, an increased discrepancy is evident (Table 12), even among the teams that used similar software platform, e.g. FEniCS (*Ambit*, *Simula*, *CHIMeRA*). This is likely due not only to differences in geometry but also to the rule-based calculation of fibers, which were provided at the degrees of freedom and then interpolated to the quadrature points. This interpolation process may introduce additional variability in the simulation outputs across the groups.

8 Limitations

This work presents a number of limitations that could be tackled in future studies.

Though this study represents a considerable improvement in modeling complexity, it still addresses only one physical field, namely mechanics. Including additional fields in a multiphysics framework —such as fluid-solid interaction, poromechanics, electromechanics, and 0D-3D models —would likely be the most reasonable next steps. However, this approach may reduce the number of groups participating in each of these benchmarks.

In principle, the observed differences may disappear if the discretization is refined to the point where all solvers reach convergence, but no detailed convergence analysis was performed in this study. At that stage, the comparison would focus primarily on computational cost, assuming that, as one would hope, all the solvers converge to the same solution but with different accuracy orders.

The present work also has limitations regarding the realism of the fiber model. While the fiber-sheet-normal model is well established for left ventricular geometries, there is a lack of data for the right ventricle and interventricular region [67]. The benchmark could be updated with more realistic biventricular fiber models [34, 68]. However, this is likely to introduce additional numerical challenges and require more careful discretization due to the thinness of the right ventricular wall.

While incorporating a human or animal geometry, especially one including the atria, is feasible in principle, it falls outside the scope of the proposed benchmark and would overly complicate the setup. Additionally, generating atrial fibers remains a significant challenge [34, 69]. Therefore, given the focus of our work, we rely on idealized ventricular geometries paired with state-of-the-art fiber models. A more realistic and complex geometry could also lead to challenges in incorporating the fiber orientation and in the comparison and interpretation of the results. The current study seeks to achieve a balanced model complexity, which would render the results useful and relevant but still allow control of relevant model properties and facilitate the comparison of the results. The study represents a significantly increased complexity compared with previous benchmarks in the field, and the inclusion of a more realistic geometry is left for a potential follow-up.

We restricted ourselves to the study of the constitutive model in [18], which is the most commonly used one in continuum-based, organ-scale simulations. In any case, apart from the newest viscoelastic model [70], we are not aware of important more recent developments in this field. Therefore, we believe our choice remains highly relevant due to the widespread use of the presented model. Providing benchmarks for other models, such as the ones reviewed recently in [71, 72] is out of the scope of the present study.

Another limitation of this study consists in that the variability of the output to all model parameters (such that representing viscoelasticity and (in)compressibility) was not studied, though those effects were included in the model.

Though some insights for mesh sensitivity are given in Benchmark 2, this aspect was not fully explored in this article, and we consider relevant for future benchmarking efforts, together with reporting more quantities such as strains and stresses, which are often more sensitive to the discretization methods.

9 Conclusion

Consensus in simulation results is an important task, as several discretization parameters need to be selected. In this software benchmark for cardiac elastodynamics, a set of physiological test cases is proposed, comprising two different geometries. The methodology for assessment of results is based on a non-blinded calibration step and consecutive blinded steps. Nine research teams within the domain of cardiac mechanics participated in this benchmark. The benchmarks are structured as a series of steps with progressively increasing complexity, offering a step-by-step approach for verifying newly developed code. In the case of the monoventricular domain, which includes analytical fiber definition, consensus of solutions is observed in all displacement directions when changing the material parameters. Notably, different numerical methods and software implementations produced comparable results, with agreement between all participating teams. For the biventricular domain, an idealized geometry is introduced, with fibers based on the state-of-the-art in cardiac mechanics. Furthermore, tuned parameters for a physiological contraction are introduced, generalizing the previous monoventricular benchmark. The results for the biventricular model become more subject to differences arising from the incompressibility handling and space discretization as well as the fiber discretization. In this more challenging case, a few groups produced consistently comparable results (qualitatively and quantitatively), though using fully different software platform and libraries. However, it is important to note that since the test cases were deliberately chosen to be realistic and complex, it makes it difficult to determine the “reference” solutions. Nevertheless, overall the results will still serve as a range of values for valuable guidance to future authors and solver developers.

The input data and simulation results are publicly available in [Zenodo](#). Some of the computational codes are also openly available: [CARPentry](#), [Ambit](#), [Simula's](#), [SimVascular](#), [life^x](#).

A Appendix: Fiber convention

The most appropriate modeling choice for the fiber, sheet, and normal directions in the ventricular region remains an active area of debate within the community of cardiac mechanics. Variability in histological studies and computational methods to extract principal tissue directions exacerbate this discussion [20, 21, 34, 73–75]. Whereas the transmural evolution of the myocardial fiber direction across the transmural wall seems to be well accepted, two main modeling approaches can be distinguished with respect to the assigned sheet and normal directions in computational ventricle models. Following the works in [18, 20], various groups take the sheet direction (\mathbf{s}) to be oriented along the transmural direction and the normal direction (\mathbf{n}) to be orthogonal to both fiber and sheet directions [21, 73, 74, 76, 77]. Following the works in [6, 7], other groups assume the normal direction (\mathbf{n}) to be oriented along the transmural direction and the sheet direction (\mathbf{s}) to be orthogonal to both fiber and normal directions. With proper tuning of the constitutive parameters, both approaches can lead to realistic deformation profiles during diastolic loading and systolic contraction. Given our choice to use tuned constitutive parameters from a group using the second convention, we followed their myocardial architecture convention for our monoventricular benchmark cases. In reality, sheet and normal vector fields can be considered to have transmural radial-longitudinal angle variations [34, 75]. As such, both conventions provide a simplified but relevant approach towards simulating cardiac mechanics starting from the end-systolic and end-diastolic configuration, respectively.

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