

A curvilinear high order finite element framework for electromechanics: from linearised electro-elasticity to massively deformable dielectric elastomers

Roman Poya ^{a,b1}, Antonio J. Gil ^{a2}, Rogelio Ortigosa ^a, Ruben Sevilla ^a, Javier Bonet ^c, Wolfgang A. Wall ^b

^a*Zienkiewicz Centre for Computational Engineering, College of Engineering, Swansea University, Bay Campus,
SA1 8EN, United Kingdom*

^b*Institute for Computational Mechanics, Technische Universität München, Boltzmannstrasse 15, D-85748
Garching b. München, Germany*

^c*University of Greenwich, London, SE10 9LS, United Kingdom*

Abstract

This paper presents a high order finite element implementation of the convex multi-variable electro-elasticity for large deformations large electric fields analyses and its particularisation to the case of small strains through a staggered scheme. With an emphasis on accurate geometrical representation, a high performance curvilinear finite element framework based on an a posteriori mesh deformation technique is developed to accurately discretise the underlying displacement-potential variational formulation. The performance of the method under near incompressibility and bending actuation scenarios is analysed with extremely thin and highly stretched components and compared to the performance of mixed variational principles recently reported by Gil and Ortigosa [1, 2, 3]. Although convex multi-variable constitutive models are elliptic hence, materially stable for the entire range of deformations and electric fields, other forms of physical instabilities are not precluded in these models. In particular, physical instabilities present in dielectric elastomers such as pull-in instability, snap-through and the formation, propagation and nucleation of wrinkles and folds are numerically studied with a detailed precision in this paper, verifying experimental findings. For the case of small strains, the essence of the approach taken lies in guaranteeing the objectivity of the resulting work conjugates, by starting from the underlying convex multi-variable internal energy, whence avoiding the need for further symmetrisation of the resulting Maxwell and Minkowski-type stresses at small strain regime. In this context, the nonlinearity with respect to electrostatic counterparts such as electric displacements is still retained, hence resulting in a formulation similar but more competitive with the existing linearised electro-elasticity approaches. Virtual prototyping of many application-oriented dielectric elastomers are carried out with an eye on pattern forming in soft robotics and other potential medical applications.

Keywords: Monolithic & staggered electro-elasticity, high order curvilinear meshes, dielectric elastomers, material instability, wrinkling

¹Corresponding author: r.poya@swansea.ac.uk

²Corresponding author: a.j.gil@swansea.ac.uk

1. Introduction

In recent years, exploiting actuation and harvesting through the heterogenous class of Electro-Active-Polymers (EAP) has received considerable research focus. In particular, the electronic subgroup of EAP such as Dielectric Elastomers (DE) and electrostrictive relaxor ferroelectric polymers or Piezoelectric Polymers (PP) have become the subject of intensive mathematical and numerical analyses. Typically, the physically insightful information that could be gained from the numerical studies, depends on the capability of the underlying mathematical model utilised to simulate EAP. This has led to the development of a diverse range of mathematical models ranging from simplified to high fidelity models [4, 5].

On one end of the spectrum lies the class of simplified formulations, where one-dimensional idealisation in the form of rod and beam structures with mass-spring-damper-capacitor support using small strains and linear electrostatic assumption is utilised [4, 6, 7]. Particularly more popular in the experimental physics community, these models have been used successfully to characterise actuation and energy harvesting capabilities of a range of materials, inherently due to their proximity to the actual experimental set-up. The potential of a material, in exhibiting electrostriction is typically exploited using such simplified formulations [8, 9, 10, 11, 12, 13].

On the other end of the spectrum lies the class of mathematically more sophisticated formulations that exploit the large deformation characteristics of EAP [5, 14, 15, 1, 2, 16, 17, 18]. The point of departure for such formulations is an assumed energy functional for the coupled electromechanical system. Conceptually, essential and suitable mathematical requirements for the energy functional such as ellipticity [19, 20], multi-variable convexity [1, 2], coercivity [21] and material frame indifference [22] can only be studied in a large deformation context. From a phenomenological point of view, these requirements or rather restrictions have important physical implications, in particular in guaranteeing the positive definiteness of the generalised electromechanical acoustic tensor, existence of real wave speeds in the material in the vicinity of an equilibrium configuration and the electromechanical stability of the material [23, 2]. Apart from these requirements other forms of physical instabilities present in dielectric elastomers such as pull-in instability, snap-through and the formation, propagation and nucleation of wrinkles have also been reported for these models, numerically as well experimentally [24, 25, 26, 27, 28, 29, 30, 31].

In an important intermediate class for electromechanics, the large deformation characteristics of the system are neglected, whereas the nonlinearity still present in the material emanates from the electrostriction of the material through the Maxwell (for vacuum V_∞) or Minkowski (for material V) stress tensors [32, 33, 34, 35]. Theoretical aspects of these formulations were first introduced in Landau and Lifshitz [33]. The practical relevance of Maxwell stress tensor has led to a widespread utilisation of these formulations for exploiting electrostriction and magnetostriction. Unfortunately, electrostrictive models based on the utilisation of Minkowski stress tensor, in the generic case of anisotropy do not satisfy material frame indifference (i.e. objectivity or invariance of the energy with respect to rotations) of the electromechanical (total) stress tensor, due to the inherent non-symmetric nature of the Minkowski stress. Several authors in the past have used *ad-hoc* solutions, such as symmetrisation of the total stress tensor, or consideration of the conservation of angular momentum in the formulation, as a remedy [34, 36]. Nevertheless, the extended electromechanical Hessian still remains non-symmetric, which dictates the development of specialised non-symmetric finite element frameworks. Recently, Bustamente [37] has shown that physically admissible energy functionals can be constructed by choosing suitable constitutive restrictions such that their linearisation yields objective Minkowski-type stresses.

The present manuscript presents a computational framework suitable for both geometrically linearised and large deformation large electric field electromechanics. A convex multi-variable strain energy description based on the works of Gil and Ortigosa [1, 2, 3] is chosen for modelling EAPs under actuation and energy harvesting scenarios. For the case of small strains, following Bustamente [37], the present manuscript extends the framework developed by Gil and Ortigosa [1, 2] to the case of geometrically linearised electrostriction, to redress the aforementioned inconsistencies for the class of intermediate formulations. Importantly, all the aforementioned mathematical requirements are imposed at a large deformation level to arrive at a physically admissible energy functional. In this context, convex multi-variable energies typically expressed in terms of fundamental kinematic measures $\{\mathbf{F}, \mathbf{H}, J\}$ are re-expressed in terms of a set of symmetric kinematics $\{\mathbf{C}, \mathbf{G}, C\}$ to guarantee the objectivity of the energy functional. Linearisation with respect to geometrical fields is then performed by perturbing the energy in the vicinity of the reference configuration. Analogous to [38], this is achieved through a staggered scheme where the equations of electrostatics are solved for in a nonlinear fashion whereas the linearised mechanical equations are updated incrementally.

Admittedly, all the three class of formulations have been primarily applied to simplified geometries where the aim has been to verify the computational framework rather than to simulate realistic electromechanical components. In the present work, emphasis is put on accurate geometrical representation of electromechanical components by utilising an a posteriori curvilinear mesh deformation technique to accurately represent the true CAD boundaries. This is achieved by placing the high order nodes in the computational mesh on the true CAD boundary (curves/surfaces) through the solid mechanics based mesh deformation technique [38], but the underlying finite element functional spaces are not modified, i.e. a standard isoparametric finite element discretisation is employed. By relying on accurate geometrical representation and standard finite element technology, the current framework takes finite element analysis of electromechanical systems beyond the verification stage through virtual prototyping of a series of electromechanical components with potential applications in pattern forming, soft-robotics and other medical applications.

The paper is organised as follows. In [section 2](#), nonlinear continuum electromechanics is described. In [section 3](#), a variational framework for displacement-potential electromechanics is described. Aspects of particularisation to small strains and the corresponding staggered scheme are also discussed in this section. In [section 4](#), a series of numerical examples pertaining to the capability of the current framework in modelling DEs are analysed, starting from the h and p convergence properties of the curvilinear finite element framework presented in [subsection 4.1](#). The effect of accurate boundary representation using high order curvilinear finite elements is analysed in [subsection 4.2](#) and compared to high order planar elements (elements with planar faces/edges). In [subsection 4.3](#), examples of electromechanical actuation for linearised electromechanics using the presented staggered scheme are discussed. In [subsection 4.4](#), the performance of the current high order finite element displacement-potential approach is compared to those of mixed Hu-Washizu formulations presented in [1, 2, 3]. In [subsection 4.5](#), impact of code transformation and data parallelism for an efficient implementation of convex multi-variable electro-elasticity in the context of high order finite elements is presented. Finally, a series of examples pertaining to the massive deformation of dielectric elastomers are presented in [subsection 4.6](#). The inherent instabilities in DEs such as pull-in instability and the formation of wrinkles are studied with detailed precision using h and p refinements, pinpointing the robustness and the high performant

capability of the current framework.

2. Nonlinear continuum electromechanics

2.1. Kinematics: motion and deformation

Let us consider the motion of an electro-active body which in its initial or material configuration is defined by a domain V of boundary ∂V with outward unit normal \mathbf{N} . After the motion, the body occupies a spatial configuration defined by a domain v of boundary ∂v with outward unit normal \mathbf{n} , as shown in Figure 1. The motion of the electro-active polymer V is defined by a pseudo-time (t) dependent mapping field ϕ which links a material particle from material configuration \mathbf{X} to spatial configuration \mathbf{x} according to $\mathbf{x} = \phi(\mathbf{X}, t)$, where displacement boundary conditions can be defined as $\mathbf{x} = (\phi)_{\partial_u V}$ on the boundary $\partial_u V \subset \partial V$, where the notation $(\cdot)_{\partial_u V}$ is used to indicate the given value of a variable (\cdot) on the boundary $\partial_u V$. The two-point

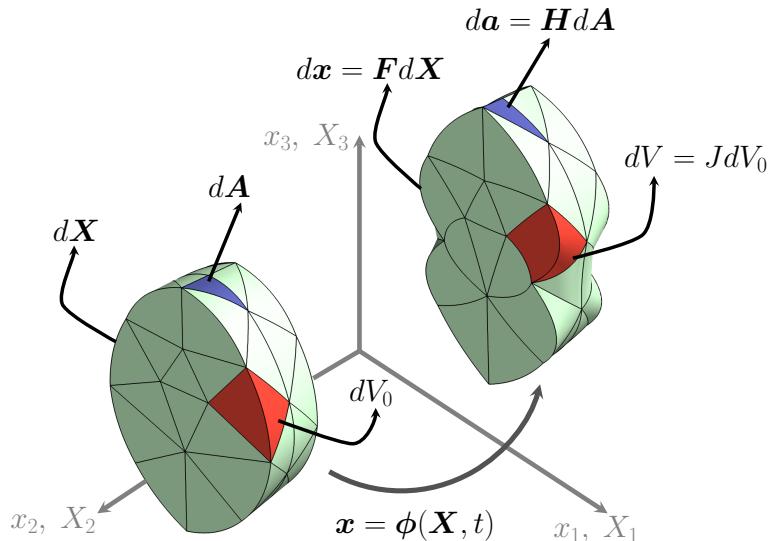


Figure 1: Deformation map of a continuum and illustration of the strain measures \mathbf{F} , \mathbf{H} and J .

deformation gradient tensor or fibre-map \mathbf{F} , which relates a fibre of differential length from the material configuration $d\mathbf{X}$ to the spatial configuration $d\mathbf{x}$, namely $d\mathbf{x} = \mathbf{F}d\mathbf{X}$, is defined as [39]

$$\mathbf{F} = \nabla_0 \mathbf{x} = \frac{\partial \phi(\mathbf{X}, t)}{\partial \mathbf{X}}, \quad (1)$$

where $\nabla_0(\cdot)$ is the material or Lagrangian gradient operator. In addition, $J = \det \mathbf{F}$ represents the Jacobian or volume-map of the deformation, which relates differential volume elements in the material configuration dV and the spatial configuration dv as $dv = JdV$. Finally, the element area vector is mapped from initial $d\mathbf{A}$ (colinear with \mathbf{N}) to final $d\mathbf{a}$ (colinear with \mathbf{n}) configuration by means of the two-point co-factor or adjoint tensor \mathbf{H} as $d\mathbf{a} = \mathbf{H}d\mathbf{A}$, which is related to the deformation gradient via the so-called Nanson's rule as $\mathbf{H} = J\mathbf{F}^{-T}$. With the help of the definition of the tensor cross product presented in [40] and utilised in the context of hyperelasticity

and electromechanics in [41, 22, 1], it is possible to re-write the area \mathbf{H} and volume J maps as

$$\mathbf{H} = \frac{1}{2} \mathbf{F} \times \mathbf{F}; \quad J = \frac{1}{3} \mathbf{H} : \mathbf{F}, \quad (2)$$

where for two-point second order tensors \mathbf{A} and \mathbf{B} , the tensor cross operation is computed as $[\mathbf{A} \times \mathbf{B}]_{iI} = \mathcal{E}_{ijk} \mathcal{E}_{IJK} A_{jJ} B_{kK}$, with \mathcal{E} the third order alternating tensor³. Based on the fundamental kinematic measures $\{\mathbf{F}, \mathbf{H}, J\}$ defined in (1) and (2), a set of symmetric kinematic measures namely \mathbf{C}, \mathbf{G} and C , can be defined as follows [22]

$$\mathbf{C} = \mathbf{F}^T \mathbf{F}; \quad \mathbf{G} = \frac{1}{2} \mathbf{C} \times \mathbf{C} = \mathbf{H}^T \mathbf{H}; \quad C = \frac{1}{3} \mathbf{G} : \mathbf{C} = J^2, \quad (3)$$

where \mathbf{C} is the right Cauchy-Green strain tensor and \mathbf{G} and C are its co-factor and determinant, respectively. Let us define $\delta \mathbf{u}$ and $\Delta \mathbf{u}$ as virtual and incremental variations of \mathbf{x} , respectively, where it will be assumed that $\delta \mathbf{u}$ and $\Delta \mathbf{u}$ satisfy compatible homogeneous displacement based boundary conditions that vanish on $\partial_u V$. Making use of the tensor cross product operation introduced above, the first and second directional derivatives of the symmetric kinematic measures $\{\mathbf{C}, \mathbf{G}, C\}$ with respect to virtual and incremental variation of the geometry can now be evaluated, starting with \mathbf{C}

$$DC[\delta \mathbf{u}] = (\nabla_0 \delta \mathbf{u})^T \mathbf{F} + \mathbf{F}^T \nabla_0 \delta \mathbf{u}; \quad (4a)$$

$$D^2 \mathbf{C}[\delta \mathbf{u}; \Delta \mathbf{u}] = (\nabla_0 \delta \mathbf{u})^T (\nabla_0 \Delta \mathbf{u}) + (\nabla_0 \Delta \mathbf{u})^T (\nabla_0 \delta \mathbf{u}), \quad (4b)$$

and similarly for the co-factor \mathbf{G}

$$DG[\delta \mathbf{u}] = \mathbf{C} \times DC[\delta \mathbf{u}]; \quad (5a)$$

$$D^2 \mathbf{G}[\delta \mathbf{u}; \Delta \mathbf{u}] = \mathbf{C} \times D^2 \mathbf{C}[\delta \mathbf{u}; \Delta \mathbf{u}] + DC[\Delta \mathbf{u}] \times DC[\delta \mathbf{u}], \quad (5b)$$

and finally for the determinant C

$$DC[\delta \mathbf{u}] = \mathbf{G} : DC[\delta \mathbf{u}] \quad (6a)$$

$$D^2 C[\delta \mathbf{u}, \Delta \mathbf{u}] = D^2 \mathbf{C}[\delta \mathbf{u}; \Delta \mathbf{u}] : \mathbf{G} + \mathbf{C} : (DC[\delta \mathbf{u}] \times DC[\Delta \mathbf{u}]), \quad (6b)$$

2.2. Translational and rotational equilibrium

The kinematics of the continuum presented in subsection 2.1 must be described through the conservation of linear momentum. In the absence of inertial effects, the global conservation of linear momentum leads to the integral translational equilibrium equations

$$\int_{\partial_t V} \mathbf{t}_0 dA + \int_V \mathbf{f}_0 dV = \mathbf{0}, \quad (7)$$

³In addition, throughout the paper, the symbol (\cdot) is used to indicate the scalar product or contraction of a single index $\mathbf{a} \cdot \mathbf{b} = a_i b_i$; the symbol $(:)$ is used to indicate double contraction of two indices $\mathbf{A} : \mathbf{B} = A_{ij} B_{ij}$; the symbol (\times) is used to indicate the cross product between vectors $[\mathbf{a} \times \mathbf{b}]_i = \xi_{ijk} a_j b_k$ and the symbol (\otimes) is used to indicate the outer or dyadic product $[\mathbf{a} \otimes \mathbf{b}]_{ij} = a_i b_j$.

where \mathbf{f}_0 represents the body force per unit undeformed volume V and \mathbf{t}_0 the traction force per unit undeformed area, applied on $\partial_t V \subset \partial V$ such that $\partial_t V \cup \partial_u V = \partial V$ and $\partial_t V \cap \partial_u V = \emptyset$. From (7), the local translational equilibrium equations and the associated boundary conditions can be written as

$$\text{DIV} \mathbf{P} + \mathbf{f}_0 = \mathbf{0} \quad \text{in } V; \quad (8a)$$

$$\mathbf{P} \mathbf{N} = \mathbf{t}_0 \quad \text{on } \partial_t V; \quad (8b)$$

$$\phi = (\phi)_{\partial_\phi V} \quad \text{on } \partial_u V, \quad (8c)$$

where \mathbf{P} represents the first Piola-Kirchoff stress tensor. Furthermore, conservation of rotational equilibrium leads to the well-known tensor condition $\mathbf{P} \mathbf{F}^T = \mathbf{F} \mathbf{P}^T$.

2.3. Electrostatics: Gauss's and Faraday's laws

In addition to the conservation of translation and rotational equilibrium presented in subsection 2.2, the electro-active polymer represented by the continuum described in subsection 2.1 is subjected in its material configuration V to an electric volume charge ρ_0^e per unit of undeformed volume and an electric surface charge ω_0^e per unit of undeformed area applied on $\partial_\omega V \subset \partial V$. Hence, the integral version of the Gauss's law can be written in a Lagrangian format as

$$\int_{\partial_\omega V} \omega_0^e dA + \int_V \rho_0^e dV = 0. \quad (9)$$

From (9), the local version of Gauss's law and the associated boundary conditions can be written as

$$\text{DIV} \mathbf{D}_0 - \rho_0^e = 0 \quad \text{in } V; \quad (10a)$$

$$\mathbf{D}_0 \cdot \mathbf{N} = -\omega_0^e \quad \text{on } \partial_\omega V, \quad (10b)$$

where \mathbf{D}_0 is the Lagrangian electric displacement vector. Alternatively, the spatial electric displacement vector \mathbf{D} can be obtained through the area push forward relationship $\mathbf{D}_0 = \mathbf{H}^T \mathbf{D}$, [5, 14]. Furthermore, in the absence of magnetic fields, the integral version of the static Faraday's law can be written in a Lagrangian form for a closed curve \mathcal{C} embedded in $V \cup \partial V$ as

$$\oint_{\mathcal{C}} \mathbf{E}_0 \cdot d\mathbf{X} = 0, \quad (11)$$

where \mathbf{E}_0 is the Lagrangian or material electric field vector. The local version of (11) and the associated boundary conditions can be written as

$$\mathbf{E}_0 = -\nabla_0 \varphi \quad \text{in } V; \quad (12a)$$

$$\varphi = (\varphi)_{\partial_\varphi V} \quad \text{on } \partial_\varphi V, \quad (12b)$$

where φ is the scalar electric potential. In (12), $\partial_\varphi V$ represents parts of the boundary ∂V where essential electric potential boundary conditions are applied such that $\partial_\omega V \cup \partial_\varphi V = \partial V$ and $\partial_\omega V \cap \partial_\varphi V = \emptyset$. The spatial electric field vector \mathbf{E} can be obtained by performing the push forward (standard fibre transformation) on material electric field i.e. $\mathbf{E}_0 = \mathbf{F}^T \mathbf{E}$.

2.4. The internal energy density: Objective convex multi-variable electro-elasticity

For the closure of the system of equations defined by (8), (10) and (12), two additional constitutive laws are needed relating deformation and electric displacements to stresses and electric fields in the continuum, satisfying appropriate constitutive inequalities, in particular ellipticity [5, 2]. In the case of reversible electro-elasticity, where thermal effects and electric polarisation induced hysteresis are disregarded, the internal energy density e per unit of undeformed volume can be solely defined in terms of the deformation and electric displacement, namely $e = e(\nabla_0 \mathbf{x}, \mathbf{D}_0)$. The requirement for objectivity (i.e. invariance with respect to rotations in the material configuration) implies that e must be independent of the rotational components of deformation. This can be facilitated by re-expressing the internal energy density in terms of a symmetric kinematic measure (such as the symmetric right Cauchy-Green tensor \mathbf{C}) as follows

$$e(\nabla_0 \mathbf{x}, \mathbf{D}_0) = e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0). \quad (13)$$

Denoting the virtual and incremental variations of the electric displacements with $\delta \mathbf{D}_0$ and $\Delta \mathbf{D}_0$, respectively, that satisfy compatible homogeneous displacement, the first law of thermodynamics yields

$$De_{\text{sym}}[\delta \mathbf{u}; \delta \mathbf{D}_0] = \mathbf{S} : \frac{1}{2} D\mathbf{C}[\delta \mathbf{u}] + \mathbf{E}_0 \cdot \delta \mathbf{D}_0, \quad (14)$$

where the second Piola-Kirchhoff stress tensor \mathbf{S} and the Lagrangian electric field \mathbf{E}_0 are computed as

$$\mathbf{S} = 2 \frac{\partial e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0)}{\partial \mathbf{C}}, \quad \mathbf{E}_0 = \frac{\partial e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0)}{\partial \mathbf{D}_0}, \quad (15)$$

wherein the first and second Piola-Kirchhoff tensors are related through $\mathbf{P} = \mathbf{F}\mathbf{S}$. Of particular importance to this development is multi-variable convexity in electromechanics, recently presented by Gil and Ortigosa [1, 2, 3], which automatically satisfies ellipticity. Using the convex multi-variable definition, the internal energy density is postulated as

$$e(\nabla_0 \mathbf{x}, \mathbf{D}_0) = W(\mathbf{F}, \mathbf{H}, J, \mathbf{D}_0, \mathbf{d}); \quad \mathbf{d} = \mathbf{F}\mathbf{D}_0, \quad (16)$$

where W represents a convex multi-variable functional in terms of the extended set of arguments $\{\mathbf{F}, \mathbf{H}, J, \mathbf{D}_0, \mathbf{d}\}$. For the requirement of objectivity, the convex multi-variable energy W can be re-expressed in terms of a set of symmetric kinematics defined in (3)

$$e(\nabla_0 \mathbf{x}, \mathbf{D}_0) = e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0) \stackrel{<}{=} W(\mathbf{F}, \mathbf{H}, J, \mathbf{D}_0, \mathbf{d}) \stackrel{>}{=} W_{\text{sym}}(\mathbf{C}, \mathbf{G}, C, \mathbf{D}_0), \quad (17)$$

where e_{sym} represents the internal energy in terms of the right Cauchy Green strain tensor \mathbf{C} and Lagrangian electric displacement \mathbf{D}_0 and W_{sym} represents an internal energy defined in terms of the extended symmetric mechanical kinematic set $\mathcal{V}_{\text{sym}}^m = \{\mathbf{C}, \mathbf{G}, C\}$ and the Lagrangian electric displacement \mathbf{D}_0 . It is worth noting that W_{sym} is not strictly convex with respect to the individual components of the set $\{\mathcal{V}_{\text{sym}}^m, \mathbf{D}_0\}$, but rather an objective re-expression of the convex multi-variable functional W . Furthermore, the inclusion of the term \mathbf{d} in W is useful in the context of studying material stability [1] which can be re-expressed in terms of the combination of kinematics in $\{\mathcal{V}_{\text{sym}}^m, \mathbf{D}_0\}$. Notice that, constructing a convex multi-variable energy (i.e. W)

is a necessary first step for a materially frame indifferent representation (i.e. W_{sym}), but the vice-versa is not necessarily true.

Definition of an objective internal energy density such as in (17) is an essential requirement for the particularisation of a multi-variable convex function to the case of small strains [38]. Furthermore, this definition of the internal energy e_{sym} (17) facilitates the introduction of a new set of work-conjugates which can now be defined as

$$\boldsymbol{\Sigma}_C = 2 \frac{\partial W_{\text{sym}}}{\partial \mathbf{C}}; \quad \boldsymbol{\Sigma}_G = 2 \frac{\partial W_{\text{sym}}}{\partial \mathbf{G}}; \quad \boldsymbol{\Sigma}_C = 2 \frac{\partial W_{\text{sym}}}{\partial C}; \quad \boldsymbol{\Sigma}_{D_0} = \frac{\partial W_{\text{sym}}}{\partial D_0}. \quad (18)$$

For notational convenience, the following sets, featuring in subsequent sections, are introduced

$$\mathcal{V}_{\text{sym}}^m = \{\mathbf{C}, \mathbf{G}, C\}; \quad \Sigma_{\mathcal{V}_{\text{sym}}}^m = \{\boldsymbol{\Sigma}_C, \boldsymbol{\Sigma}_G, \boldsymbol{\Sigma}_C\}; \quad (19a)$$

$$\mathcal{V}^e = \{D_0\}; \quad \Sigma_{\mathcal{V}}^e = \{\boldsymbol{\Sigma}_{D_0}\}; \quad (19b)$$

$$\mathcal{V} = \{\mathcal{V}_{\text{sym}}^m, \mathcal{V}^e\}; \quad \Sigma_{\mathcal{V}} = \{\Sigma_{\mathcal{V}_{\text{sym}}}^m, \Sigma_{\mathcal{V}}^e\} \quad (19c)$$

Following [22, 1, 2], a physically more insightful representation for the second Piola-Kirchhoff stress tensor \mathbf{S} and the Lagrangian electric field \mathbf{E}_0 can be obtained as

$$\mathbf{S} = \boldsymbol{\Sigma}_C + \boldsymbol{\Sigma}_G \times \mathbf{C} + \boldsymbol{\Sigma}_C \mathbf{G} \quad \mathbf{E}_0 = \boldsymbol{\Sigma}_{D_0}. \quad (20)$$

Subsequently, the internal energy $e_{\text{sym}} = e_{\text{sym}}(\mathbf{C}, D_0)$ can be further linearised leading to a tangent operator (facilitating a Newton-Raphson solution) which can be defined as follows

$$D^2 e_{\text{sym}}[\delta \mathbf{u}, \delta D_0; \Delta \mathbf{u}, \Delta D_0] = [\frac{1}{2} D \mathbf{C}[\delta \mathbf{u}] : \delta D_0] \begin{bmatrix} \mathbf{C} & \mathbf{Q}^T \\ \mathbf{Q} & \boldsymbol{\theta} \end{bmatrix} \begin{bmatrix} : \frac{1}{2} D \mathbf{C}[\Delta \mathbf{u}] \\ \Delta D_0 \end{bmatrix} + \mathbf{S} : \frac{1}{2} D^2 \mathbf{C}[\delta \mathbf{u}; \Delta \mathbf{u}], \quad (21)$$

with the fourth order elasticity tensor \mathbf{C} , the third order coupling tensor \mathbf{Q} and the second order dielectric tensor $\boldsymbol{\theta}$ defined as

$$\mathbf{C} = 4 \frac{\partial^2 e_{\text{sym}}(\mathbf{C}, D_0)}{\partial C \partial C}; \quad \mathbf{Q} = 2 \frac{\partial^2 e_{\text{sym}}(\mathbf{C}, D_0)}{\partial D_0 \partial C}; \quad \boldsymbol{\theta} = \frac{\partial^2 e_{\text{sym}}(\mathbf{C}, D_0)}{\partial D_0 \partial D_0}. \quad (22)$$

Analogous to (20), a more physically insightful representation of the tangent operator can be obtained as

$$D^2 e_{\text{sym}}[\delta \mathbf{u}; \Delta \mathbf{u}] = [\mathbb{M}_\delta]^T [\mathbb{H}_{W_{\text{sym}}}] [\mathbb{M}_\Delta] + (\boldsymbol{\Sigma}_G + \boldsymbol{\Sigma}_C \mathbf{C}) : (\frac{1}{2} D \mathbf{C}[\delta \mathbf{u}] \times \frac{1}{2} D \mathbf{C}[\Delta \mathbf{u}]) + \mathbf{S} : [(\nabla_0 \delta \mathbf{u})^T (\nabla_0 \Delta \mathbf{u})], \quad (23)$$

where

$$[\mathbb{M}_\delta]^T = [\frac{1}{2} D \mathbf{C}[\delta \mathbf{u}] : \frac{1}{2} D \mathbf{G}[\delta \mathbf{u}] : \frac{1}{2} D \mathbf{C}[\delta \mathbf{u}] \cdot \delta D_0]; \quad (24)$$

$$[\mathbb{M}_\Delta] = \begin{bmatrix} : \frac{1}{2} D \mathbf{C}[\Delta \mathbf{u}] \\ : \frac{1}{2} D \mathbf{G}[\Delta \mathbf{u}] \\ \frac{1}{2} D \mathbf{C}[\Delta \mathbf{u}] \\ \Delta D_0 \end{bmatrix}, \quad (25)$$

with the Hessian operator $[\mathbb{H}_{W_{\text{sym}}}]$ defined as

$$[\mathbb{H}_{W_{\text{sym}}}] = \begin{bmatrix} 4W_{\text{sym}}_{CC} & 4W_{\text{sym}}_{CG} & 4W_{\text{sym}}_{CC} & 2W_{\text{sym}}_{CD_0} \\ 4W_{\text{sym}}_{GC} & 4W_{\text{sym}}_{GG} & 4W_{\text{sym}}_{GC} & 2W_{\text{sym}}_{GD_0} \\ 4W_{\text{sym}}_{CC} & 4W_{\text{sym}}_{CG} & 4W_{\text{sym}}_{CC} & 2W_{\text{sym}}_{CD_0} \\ 2W_{\text{sym}}_{D_0C} & 2W_{\text{sym}}_{D_0G} & 2W_{\text{sym}}_{D_0C} & W_{\text{sym}}_{D_0D_0} \end{bmatrix}, \quad (26)$$

where the components of \mathcal{C} , \mathcal{Q} and θ can now be defined in terms of the set of work-conjugates Σ_V defined in (19) using Table 1. It is important to note that, as opposed to the convex multi-

\mathcal{C}	$\begin{aligned} & W_{\text{sym}}_{CC} + \mathcal{C} \times (W_{\text{sym}}_{GG} \times \mathcal{C}) + (W_{\text{sym}}_{CC} \mathcal{G} \otimes \mathcal{G}) \\ & + W_{\text{sym}}_{CG} \times \mathcal{C} + \mathcal{C} \times W_{\text{sym}}_{CG} \\ & + W_{\text{sym}}_{CC} \otimes \mathcal{G} + \mathcal{G} \otimes W_{\text{sym}}_{CC} \\ & + (\mathcal{C} \times W_{\text{sym}}_{CC}) \otimes \mathcal{G} + \mathcal{G} \otimes (W_{\text{sym}}_{CC} \times \mathcal{C}) \end{aligned}$
\mathcal{Q}^T	$W_{\text{sym}}_{CD_0} + \mathcal{C} \times W_{\text{sym}}_{GD_0} + \mathcal{G} \otimes W_{\text{sym}}_{CD_0}$
θ	$W_{D_0D_0}$

Table 1: Elasticity tensor \mathcal{C} , piezoelectric tensor \mathcal{Q} and dielectric θ tensor re-expressed in terms of the components of the Hessian operator.

variable Hessian operator expressed in terms of the fundamental kinematic set $\{\mathcal{F}, \mathcal{H}, J, \mathcal{D}_0, \mathcal{d}\}$ presented in [1], the constitutive term (first term) in (23) is not strictly positive definite as multi-variable convexity is not defined with respect to the set $\mathcal{V}_{\text{sym}}^m$ and, hence, ellipticity (i.e. rank-one convexity) of the internal energy based on this constitutive term alone cannot be established.

2.5. The Helmholtz-like energy density and Legendre transformation

In the case of pursuing a standard variational implementation via the finite element method, where the scalar electric potential is preferred as an unknown over the electric displacement field vector, it is typically preferred to work with the Helmholtz-like energy $\Phi = \Phi(\mathcal{C}, -\nabla_0 \varphi)$, defined as [1, 2]⁴

$$\Phi(\mathcal{C}, -\nabla_0 \varphi) = -\sup_{\mathcal{D}_0} \{-\nabla_0 \varphi \cdot \mathcal{D}_0 - e_{\text{sym}}(\mathcal{C}, \mathcal{D}_0)\}, \quad (27)$$

⁴The usage of Helmholtz-like energy is due to the fact that the Helmholtz energy is predominantly used in the context of thermal problems for the free energy expressed in terms temperature and the Legendre transformation therein to obtain the internal energy expressed in terms of entropy.

wherein the second Piola-Kirchhoff stress tensor \mathbf{S} and the Lagrangian electric displacement \mathbf{D}_0 are computed as

$$\mathbf{S} = 2 \frac{\partial \Phi(\mathbf{C}, \mathbf{E}_0)}{\partial \mathbf{C}} \Big|_{\mathbf{E}_0 = -\nabla_0 \varphi}; \quad \mathbf{D}_0 = -\frac{\partial \Phi(\mathbf{C}, \mathbf{E}_0)}{\partial \mathbf{E}_0} \Big|_{\mathbf{E}_0 = -\nabla_0 \varphi}. \quad (28)$$

Further linearisation of the Helmholtz-like energy leads to the three constitutive tensors analogous to (22)

$$\mathcal{C}_\Phi = \frac{\partial^2 \Phi(\mathbf{C}, \mathbf{E}_0)}{\partial \mathbf{C} \partial \mathbf{C}} \Big|_{\mathbf{E}_0 = -\nabla_0 \varphi}; \quad \mathcal{Q}_\Phi = \frac{\partial^2 \Phi(\mathbf{C}, \mathbf{E}_0)}{\partial \mathbf{C} \partial \mathbf{E}_0} \Big|_{\mathbf{E}_0 = -\nabla_0 \varphi}; \quad \boldsymbol{\theta}_\Phi = \frac{\partial^2 \Phi(\mathbf{C}, \mathbf{E}_0)}{\partial \mathbf{E}_0 \partial \mathbf{E}_0} \Big|_{\mathbf{E}_0 = -\nabla_0 \varphi}, \quad (29)$$

where the subscript $\mathbf{E}_0 = -\nabla_0 \varphi$ implies that energy is evaluated for the value of electric field obtained using (12). When the internal energy is a complex multi-variable function of the set \mathcal{V} , the nonlinearity of the convex multi-variable function e_{sym} can make it impossible to obtain an explicit representation of the constitutive tensors in (29). Whence, it is typically more suitable to perform this step numerically, by utilising a nonlinear iterative scheme. In the context of finite element analysis, by exploiting the relationship between the internal energy and the Helmholtz-like energy through the Legendre transform in (27), it is possible to compute, via a Newton-Raphson algorithm per quadrature point, these tensors as

$$[\boldsymbol{\theta}_\Phi]_{ij} = -[\boldsymbol{\theta}]_{ij}^{-1}; \quad [\mathcal{Q}_\Phi]_{ijk} = -[\boldsymbol{\theta}_\Phi]_{mi} [\mathcal{Q}]_{jkm}; \quad [\mathcal{C}_\Phi]_{ijkl} = [\mathcal{C}]_{ijkl} - [\mathcal{Q}]_{ijm} [\mathcal{Q}_\Phi]_{mkl}. \quad (30)$$

For a detailed representation of different electromechanical energies, the reader can refer to Gil and Ortigosa [1].

3. Variational formulation

3.1. Displacement-electric potential based variational formulation

A variational principle can be established by the total energy minimisation defined in terms of the internal energy of the system $e_{\text{sym}} = e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0)$. In this case, the total potential energy $\Pi_e(\mathbf{x}^*, \varphi^*, \mathbf{D}_0^*)$ can be written as

$$\Pi_e(\mathbf{x}^*, \varphi^*, \mathbf{D}_0^*) = \inf_{\mathbf{x}, \mathbf{D}_0} \sup_{\varphi} \left\{ \int_V e_{\text{sym}}(\mathbf{C}, \mathbf{D}_0) dV + \int_V \mathbf{D}_0 \cdot \nabla_0 \varphi dV - \Pi_{\text{ext}}(\mathbf{x}, \varphi) \right\}, \quad (31)$$

where $(\mathbf{x}^*, \varphi^*, \mathbf{D}_0^*)$ denotes the exact solution and $\Pi_{\text{ext}}(\mathbf{x}, \varphi)$ represents the external coupled electromechanical work additively decomposed into the purely mechanical $\Pi_{\text{ext}}^m(\mathbf{x})$ and electrical $\Pi_{\text{ext}}^e(\varphi)$ components

$$\begin{aligned} \Pi_{\text{ext}}^m(\mathbf{x}) &= \int_V \mathbf{f} \cdot \mathbf{x} dV + \int_{\partial_t V} \mathbf{t}_0 \cdot \mathbf{x} dA; \\ \Pi_{\text{ext}}^e(\varphi) &= - \int_V \rho_0^e \varphi dV - \int_{\partial_\omega V} \omega_0^e \varphi dA, \end{aligned}$$

where

$$\Pi_{\text{ext}}(\mathbf{x}, \varphi) = \Pi_{\text{ext}}^m(\mathbf{x}) + \Pi_{\text{ext}}^e(\varphi).$$

Application of the Legendre transform enables [1] above variational principle (31) to be reformulated as

$$\Pi_\Phi(\boldsymbol{x}^*, \varphi^*) = \inf_{\boldsymbol{x}} \sup_{\varphi} \left\{ \int_V \Phi(\boldsymbol{C}, -\nabla_0 \varphi) dV - \Pi_{\text{ext}}(\boldsymbol{x}, \varphi) \right\}. \quad (32)$$

It is now straightforward to carry out finite element discretisation of (32) in terms of $\{\boldsymbol{x}, \varphi\}$, for which we refer to [2]. In this context, (32) can be solved for in a standard monolithic way as described in Algorithm 1. In the current setting, equal order high order isoparametric finite elements are utilised for the interpolation of the primary variables i.e. $\{\boldsymbol{x}, \varphi\}$.

Algorithm 1 The nonlinear electromechanics solver

```

procedure MONOLITHIC SOLVER
    Input geometry, material properties and analysis parameters
    Initialise  $\mathbf{F} = \mathbf{0}$ ,  $\boldsymbol{x} = \mathbf{X}$  and  $\mathbf{R} = \mathbf{0}$ 
    for each increment  $n$  do
        Compute incremental nodal forces  $\Delta\mathbf{F}$ 
        Compute electromechanical nodal forces  $\mathbf{F} = \mathbf{F} + \Delta\mathbf{F}$ 
        Compute electromechanical residuals  $\mathbf{R} = \mathbf{R} - \Delta\mathbf{F}$ 
        while  $\|\mathbf{R}\|/\|\mathbf{F}\| > \text{tolerance}$  do
            Assemble  $\mathbf{K}$ 
            Solve  $\mathbf{Ku} = -\mathbf{R}$ 
            Update the geometry  $\boldsymbol{x} = \boldsymbol{x} + \mathbf{u}$ 
            for every quadrature point do
                Given  $\mathbf{E}_0$  compute  $\mathbf{D}_0$  implicitly via 22(c)
                Compute  $\mathcal{C}_\Phi$ ,  $\mathcal{Q}_\Phi$  and  $\boldsymbol{\theta}_\Phi$  using (30)
                Compute second Piola-Kirchhoff stress tensor  $\mathbf{S}$  using (20)
            end for
            Compute traction forces  $\mathbf{T}$ 
            Find  $\mathbf{R} = \mathbf{T} - \mathbf{F}$ 
        end while
    end for
end procedure

```

3.2. Incrementally linearised electromechanics

As presented in section 1, a particular class of coupling in electromechanics deals with small strains coupled with nonlinear electrostatics. To this end, the variational principle described in (32) can be particularised for the case of small strains. Extending, the strategy presented in [38, 42], let us consider the total potential energy (32) cast in an iterative (Newton-Raphson) form

$$\Pi_{\Phi_{\text{int}}}(\boldsymbol{x}_{k+1}^{*^{n+1}}, \varphi_{k+1}^{*^{n+1}}) = \int_V \Phi(\boldsymbol{C}_{k+1}^{n+1}, -(\nabla_0 \varphi)_{k+1}^{n+1}) dV, \quad (33)$$

where superscripts denote increments and subscripts denote iterations. A geometrically linearised formulation for electro-elasticity can be obtained by perturbing the potential energy (33). Ho-

wever note that in contrast to the approach presented in [38], particularisation of (33) for small strains through consistent linearisation with respect to the variations in electrostatic fields and mechanical fields cannot be performed simultaneously, as it leads to high order tensorial quantities without clear physical interpretations. However, as shown in Figure 2, for small deformations, the geometry could be solved for incrementally with only the electrostatic equations requiring a full Newton-Raphson procedure. This gives rise to the staggered scheme presented in the following subsection.

3.3. The staggered approach to incrementally linearised electromechanics

To elaborate the staggered approach emanating from the variational principle (31), let us reconsider (33) which in fully discretised form (using the finite element method) can be written as

$$\begin{bmatrix} \mathbf{K}_{\mathbf{uu}}^{n+1} & \mathbf{K}_{\mathbf{u}\phi}^{n+1} \\ \mathbf{K}_{\phi\mathbf{u}}^{n+1} & \mathbf{K}_{\phi\phi}^{n+1} \end{bmatrix} \begin{bmatrix} \mathbf{U}_{\mathbf{u}k+1}^{n+1} \\ \mathbf{U}_{\phi k+1}^{n+1} \end{bmatrix} = - \begin{bmatrix} \mathbf{R}_{\mathbf{u}k+1}^{n+1} \\ \mathbf{R}_{\phi k+1}^{n+1} \end{bmatrix}, \quad (34)$$

where $\mathbf{K}_{\mathbf{uu}}$ represents the fully discrete purely mechanical stiffness matrix, $\mathbf{K}_{\phi\phi}$ the fully discrete matrix associated with electrostatic variable(s) and $\mathbf{K}_{\mathbf{u}\phi}$ and $\mathbf{K}_{\phi\mathbf{u}}$ the fully discrete electromechanical coupled matrices. The above discrete form particularised to the staggered case of (31) can be established as shown in Algorithm 2. Algorithm 2 in particular implies a staggered system in which the geometry is updated incrementally, but the associated electrostatic variables are solved for iteratively. More specifically, as shown in Figure 2 the geometry, the mechanical and the coupling matrices are frozen during the iterative (implicit) solution of the electrostatic variables. An algorithmic representation of this staggered scheme is presented in Algorithm 2 where \mathbf{N} represents the external electrical nodal force vector. Note that an initial step for solving the mechanical variables is necessary in this case i.e. $\mathbf{K}_{\mathbf{uu}0}^0 \mathbf{U}_{\mathbf{u}0}^0 = -\mathbf{R}_{\mathbf{u}0}^0$ to account for Dirichlet driven problems. Since the above staggered approach requires the solution of a scalar field electric potential through solving the discretised Gauss's law, the saving in computational cost can be tremendous within its range of applicability (small strains). However as the voltage induced deformation increases a high number of increments might be needed to obtain the results of fully nonlinear monolithic approach.

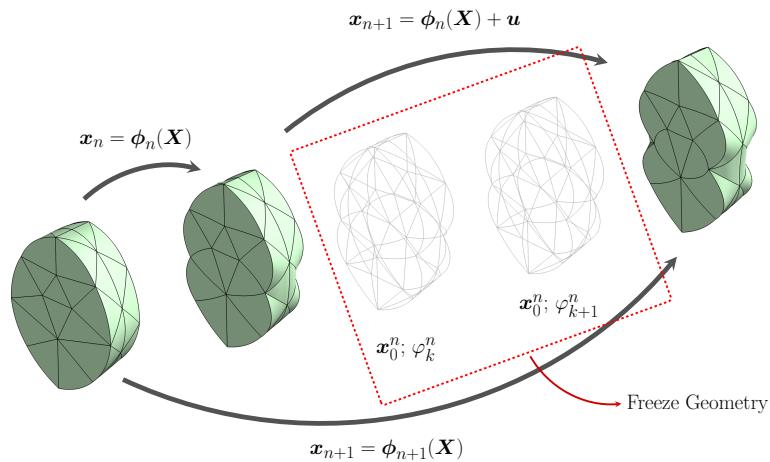


Figure 2: Schematic representation of the staggered incrementally linearised scheme.

Algorithm 2 The incrementally linearised electromechanics solver

procedure STAGGERED SOLVER

Assemble & solve the mechanical problem $\mathbf{K}_{\mathbf{uu}_0}^0 \mathbf{U}_{\mathbf{u}_0}^0 = -\mathbf{R}_{\mathbf{u}_0}^0$

for each increment n **do**

Compute electrostatic residual $\hat{\mathbf{R}}_{\phi_0}^{n+1} = \mathbf{R}_{\phi_0}^{n+1} - \mathbf{K}_{\phi\mathbf{u}_0^n} \mathbf{U}_{\mathbf{u}_0}^n$

while $\|\hat{\mathbf{R}}_{\phi_{k+1}}^{n+1}\| > tolerance$ **do**

Assemble & solve the electrostatic problem $\mathbf{K}_{\phi\phi_{k+1}}^{n+1} \mathbf{U}_{\phi_{k+1}}^{n+1} = -\hat{\mathbf{R}}_{\phi_{k+1}}^{n+1}$

Compute electrostatic traction forces $\mathbf{T}_{\phi_{k+1}}^{n+1}$

Update electrostatic residual $\hat{\mathbf{R}}_{\phi_{k+1}}^{n+1} = \mathbf{T}_{\phi_{k+1}}^{n+1} - \mathbf{N}_{\phi_{k+1}}^{n+1}$

Accumulate mechanical residual $\mathbf{R}_{\phi\mathbf{u}_{k+1}}^{n+1} = \mathbf{T}_{\mathbf{u}_{k+1}}^{n+1} - \mathbf{N}_{\mathbf{u}_{k+1}}^{n+1}$

end while

Compute the force vector $\mathbf{F}_{\mathbf{u}\phi_0^n} = \mathbf{K}_{\mathbf{u}\phi_0^n} \mathbf{U}_{\phi_{k+1}}^{n+1}$

Solve the corrected mechanical problem $\mathbf{K}_{\mathbf{uu}_0^n} \mathbf{U}_{\mathbf{u}_0}^{n+1} = -\mathbf{R}_{\mathbf{u}_0^n} - \mathbf{R}_{\phi\mathbf{u}_{k+1}}^{n+1} + \mathbf{F}_{\mathbf{u}\phi_0^n}$

Update the geometry $\mathbf{x}^{n+1} = \mathbf{x}^n + \mathbf{U}_{\mathbf{u}_0}^{n+1}$

Assemble $\mathbf{K}_{\mathbf{uu}_0^{n+1}}$, $\mathbf{K}_{\mathbf{u}\phi_0^{n+1}}$ and $\mathbf{K}_{\phi\mathbf{u}_0^{n+1}}$

Set $n + 1$ to n

end for

end procedure

4. Numerical examples

In this section a series of numerical examples for electromechanics are presented. These include a) (mesh refinement) h and (polynomial enrichment) p convergence studies for high order displacement-potential formulation for convex multi-variable internal energies presented in [section 3](#), b) the impact of accurate geometrical representations on the solution of large deformation electromechanical problems, c) comparison of monolithic (nonlinear) approach with the incrementally linearised staggered approach and the range of applicability of the latter approach in actuation, d) comparison of the current framework with the mixed Hu-Washizu variational principles presented in [\[2, 16\]](#) and, e) Impact of code transformation and data parallelism for an efficient implementation of convex multi-variable electro-elasticity in the context of high order finite elements. Apart from these benchmark studies, a series of examples pertaining to the massive deformation and the instabilities in DEs such as the formation of folds and wrinkles is studied in detail using the current framework. The finite element implementation involves equal order standard isoparametric discretisation of the electromechanical variables $\{\mathbf{x}, \varphi\}$, starting with at least quadratic basis functions. As a standard nomenclature in high order finite element analysis, polynomial interpolation over tetrahedra are denoted by p and polynomial enrichment over hexahedra are denoted by q . Unless otherwise specified, for all the examples, the curvilinear meshes are generated using the consistently linearised solid mechanics analogy presented in [\[38, 43\]](#), using a Mooney-Rivlin model with a Poisson's ratio of 0.45 and 10 load increments. In addition, all the analyses are carried out with the high performance domain-specific data parallel tensor contraction finite element framework **Fastor** based on the authors' previous work [\[44\]](#) and the massively parallel multi-frontal direct sparse solver **MUMPS** is used for the solution of system of linear equations. For the purpose of assessing the performance of the proposed monolithic approach the final example is solved using the algebraic multigrid (AMG) solver with the standard

Ruge-Stuben aggregations, c.f. [45, 46, 47, 48].

4.1. h & p convergence of the proposed high order framework

As a starting point, it is essential to examine the h and p convergence properties of the proposed high order displacement-potential formulation for large deformations large electric fields (fully coupled monolithic approach) described in section 2 and subsection 3.1. Hence, the objective of this example is to a) assess the convergence of different variables in $\{\boldsymbol{x}, \varphi, \mathcal{V}, \Sigma_{\mathcal{V}}\}$ using h and p refinements, b) examine optimality of the algorithm in terms of convergence properties on high order curved tetrahedral and hexahedral meshes around a dielectric elastomeric patch, obtained using the mesh deformation technique described in [38] and, c) showcase the scalability of the framework with high polynomial enrichment. The geometry of the dielectric patch is shown in Figure 3. A similar convergence study for an eleven field Hu-Washizu type mixed variational formulation in terms of the set of unknowns $\{\boldsymbol{x}, \boldsymbol{F}, \boldsymbol{H}, J, \varphi, \boldsymbol{d}, \boldsymbol{\Sigma}_F, \boldsymbol{\Sigma}_H, \boldsymbol{\Sigma}_J, \boldsymbol{\Sigma}_{D_0}, \boldsymbol{\Sigma}_d\}$ is presented in [2]. The constitutive model considered is based on a convex multi-variable energy functional expressed in terms of the invariants of the set \mathcal{V} as follows

$$W_{el,1}(\boldsymbol{C}, \boldsymbol{G}, C, \boldsymbol{D}_0) = \mu_1 \sum_{i=1}^n \alpha_i I_{\boldsymbol{C}}^i + \mu_2 \sum_{i=1}^n \beta_i I_{\boldsymbol{G}}^i - 2 \left(\mu_1 \sum_{i=1}^n i \alpha_i 3^{i-1} + 2 \mu_2 \sum_{i=1}^n i \beta_i 3^{i-1} \right) \ln \sqrt{C} \\ + \frac{\lambda}{2} (\sqrt{C} - 1)^2 + \frac{1}{2\varepsilon_1} II_{\boldsymbol{D}_0} + \frac{1}{2\varepsilon_2 \sqrt{C}} II_{\boldsymbol{d}}, \quad (35)$$

where $I_{(\bullet)}$ denotes the trace of the entity (\bullet) and $II_{(\bullet)}$ the squared of the \mathcal{L}^2 norm of the entity (\bullet) with $II_{\boldsymbol{d}} = \boldsymbol{D}_0 \cdot \boldsymbol{C}\boldsymbol{D}_0$. Furthermore, $\{\mu_1, \mu_2, \lambda, \varepsilon_1, \varepsilon_2\}$ represent positive material constants with α_i 's and β_i 's denoting scaling coefficients. The material parameters in (35) used for this example are presented in Table 2, where $n = 2$, $\alpha_1 = \beta_1 = 1$ and $\alpha_2 = \beta_2 = 0.2$ are chosen. The

μ_1 (Pa)	μ_2 (Pa)	λ (Pa)	ε_1 (N/V ²)	ε_2 (N/V ²)
1	1/2	1	4	4

Table 2: Material properties for example 4.1

problem is constructed so that smoothness of the solution is guaranteed. For that purpose, the following simple exact fields associated with the fields \boldsymbol{x} and φ are considered

$$\boldsymbol{x}^{\text{exact}} = \boldsymbol{X} + \begin{Bmatrix} \gamma_1 \sin X_1 \\ \gamma_2 \cos X_2 \\ \gamma_3 (\sin X_3 + \cos X_3) \end{Bmatrix}; \quad \varphi^{\text{exact}} = \varphi_0 \sin X_1, \quad (36)$$

where for this benchmark example, $\gamma_1 = 0.1$, $\gamma_2 = 0.2$, $\gamma_3 = 0.3$ and $\varphi_0 = 10^4 V$ are chosen. The deformation gradient tensor and the Lagrangian electric field can now be computed as

$$\boldsymbol{F}^{\text{exact}} = \begin{bmatrix} 1 + \gamma_1 \cos X_1 & 0 & 0 \\ 0 & 1 - \gamma_2 \sin X_2 & 0 \\ 0 & 0 & 1 + \gamma_3 (\cos X_3 - \sin X_3) \end{bmatrix}; \quad \boldsymbol{E}_0^{\text{exact}} = - \begin{Bmatrix} \varphi \cos X_1 \\ 0 \\ 0 \end{Bmatrix}. \quad (37)$$

The remaining exact fields in the set $\mathcal{V}^{\text{exact}} = \{\mathbf{C}^{\text{exact}}, \mathbf{G}^{\text{exact}}, C^{\text{exact}}, \mathbf{D}_0^{\text{exact}}\}$ can now be obtained for the smooth displacement and electric potential fields, from (37). Application of (3) on (37) yields the exact right Cauchy-Green tensor, its cofactor and determinant for the smooth fields (36) as

$$\mathbf{C}^{\text{exact}} = \mathbf{F}^{\text{exact}}{}^T \mathbf{F}^{\text{exact}}; \quad \mathbf{G}^{\text{exact}} = \frac{1}{2} \mathbf{C}^{\text{exact}} \times \mathbf{C}^{\text{exact}}; \quad C^{\text{exact}} = \frac{1}{3} \mathbf{G}^{\text{exact}} : \mathbf{C}^{\text{exact}}. \quad (38)$$

Similarly, the Lagrangian electric displacement vector can be computed by applying (15) on (35) as

$$\mathbf{D}_0^{\text{exact}} = \left(\frac{1}{\varepsilon_1} \mathbf{I} + \frac{1}{\varepsilon_2 \sqrt{C^{\text{exact}}}} \mathbf{C}^{\text{exact}} \right)^{-1} \mathbf{E}_0^{\text{exact}}. \quad (39)$$

Once all the elements of the set $\mathcal{V}^{\text{exact}}$ have been determined, it is possible to obtain the set of exact work conjugates $\Sigma_{\mathcal{V}}^{\text{exact}} = \{\Sigma_{\mathbf{C}}^{\text{exact}}, \Sigma_{\mathbf{G}}^{\text{exact}}, \Sigma_C^{\text{exact}}, \Sigma_{\mathbf{D}_0}^{\text{exact}}\}$ via (18). These enable to compute the second Piola-Kirchhoff stress tensor $\mathbf{S}^{\text{exact}}$ from (20). Finally, the associated volumetric force and electric charge in mechanical and electrical equilibrium with the exact first Piola-Kirchhoff stress tensor $\mathbf{P}^{\text{exact}} = \mathbf{F}^{\text{exact}} \mathbf{S}^{\text{exact}}$ and exact Lagrangian electric displacement field $\mathbf{D}_0^{\text{exact}}$ are determined from (8) and (10), respectively as

$$\mathbf{f}_0(\mathbf{x}^{\text{exact}}, \varphi^{\text{exact}}) = -\text{DIV} \mathbf{P}^{\text{exact}}; \quad \rho_0(\mathbf{x}^{\text{exact}}, \varphi^{\text{exact}}) = \text{DIV} \mathbf{D}_0^{\text{exact}}. \quad (40)$$

For the convergence studies, three different high order curvilinear unstructured tetrahedral meshes and one high order curvilinear structured hexahedral mesh for the dielectric patch in Figure 3 are considered with 532 elements, 9220 elements, 26807 elements for the first three and 5000 elements for the last one, respectively, as shown in Figure 5. As will be described shortly, these discretisations are chosen such that the maximum segment fitting in a curved element (denoted by h) is successively refined for h -convergence. The placements of high order nodes on curved boundaries of the meshes are given particular importance, in that they have been computed through an arc-length based projection from high order planar meshes with Warburton nodal distribution [38]. The curved volume mesh is then obtained by applying the consistently linearised elastic analogy [38, 43, 49, 50]. It is worth mentioning that the quality of the curvilinear meshes directly impact the h and p convergence property (optimality) of the finite element interpolation scheme. To this end, Table 3 and Table 4 report the three fundamental quality measures (where Q_1 quantifies distortion of edges, Q_2 quantifies distortion of faces and Q_3 , also known as scaled Jacobian, quantifies volumetric distortion of the element itself) of the curved tetrahedral and hexahedral meshes, respectively; c.f. [38]. Optimal symmetric quadrature rules for tetrahedra reported in [51] are utilised for numerical integration of high order tetrahedral elements. For hexahedral elements, a simple tensor product based quadrature rule is utilised.

p	Coarse (532 elements)			Medium (9220 elements)			Fine (26807 elements)					
	DoFs	Q_1	Q_2	Q_3	DoFs	Q_1	Q_2	Q_3	DoFs	Q_1	Q_2	Q_3
$p = 2$	1027×4	0.888	0.875	0.790	17934×4	0.962	0.924	0.885	52590×4	0.953	0.909	0.866
$p = 3$	3077×4	0.881	0.843	0.746	54018×4	0.936	0.869	0.799	158352×4	0.947	0.893	0.838
$p = 4$	6861×4	0.874	0.813	0.738	120473×4	0.926	0.850	0.771	352838×4	0.935	0.871	0.800
$p = 5$	12911×4	0.851	0.807	0.711	226516×4	0.904	0.831	0.721	662848×4	0.924	0.843	0.765
$p = 6$	21759×4	0.846	0.771	0.652	381373×4	0.891	0.828	0.713	1115196×4	0.901	0.837	0.757

Table 3: Distortion quality of high order curvilinear tetrahedral meshes

p	DoFs	Q_1	Q_2	Q_3
$p = 2$	49943×4	0.999	0.984	0.914
$p = 3$	147364×4	0.993	0.972	0.889
$p = 4$	341885×4	0.987	0.961	0.853
$p = 5$	659106×4	0.981	0.934	0.820
$p = 6$	1129027×4	0.971	0.900	0.797

Table 4: Distortion quality of high curvilinear hexahedral mesh

For studying h -convergence properties, only tetrahedral meshes are considered. The convergence rate of different primary and derived variables $\{\mathbf{x}, \varphi, \mathcal{V}, \Sigma_{\mathcal{V}}\}$ are then studied by comparing the interpolated solution and analytical solution for a fixed p and successive h -refinement, similar to the strategy followed in [38, 2]. To monitor the convergence rate, the $\mathcal{L}^2(V)$ norm of the error is computed for all quantities of interest. For curvilinear meshes the $\mathcal{L}^2(V)$ norms of the variables

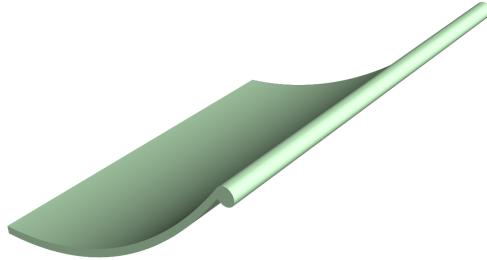


Figure 3: CAD geometry of the dielectric patch

are reported as a function of the diameter (the largest segment that can fit within a curved region) of the element computed through a straightforward sampling strategy illustrated in Figure 4 for curved tetrahedral and hexahedral elements. After traversing the whole computational mesh, the minimum value of all the diameters is chosen as h [52, 53]. The tetrahedral meshes reported in Table 3 are chosen such that they correspond to successive refinement in this quantity.

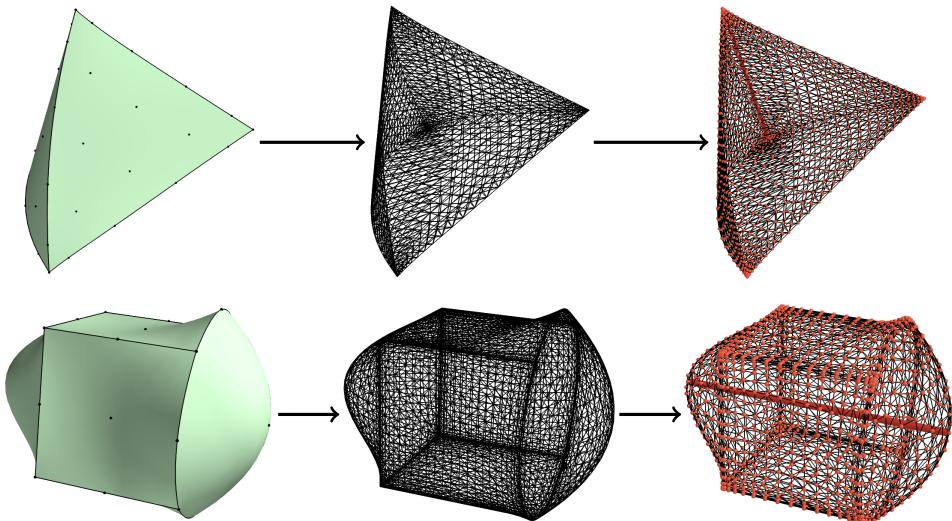


Figure 4: Illustration of computing the diameter (largest segment) in curvilinear meshes using sampling and tessellation strategy, for $p=5$ tetrahedra and $q=2$ hexahedra

Figure 6 shows h -convergence of the variables $\{\mathbf{x}, \varphi, \mathcal{V}, \Sigma_{\mathcal{V}}\}$ for four levels of p -refinement (i.e. $p = 2, 3, 4, 5$). Note that unlike in [2], while the displacements and electric potential are primary

(solved) variables, the rest of variables are secondary (derived). As can be observed in [Figure 6](#) for

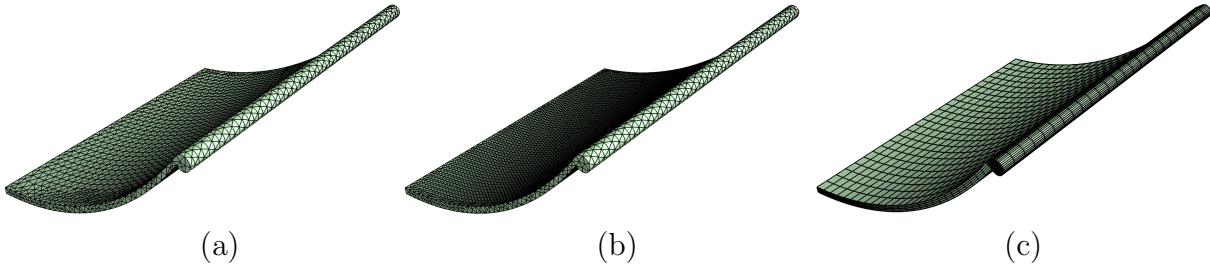


Figure 5: Three representative curvilinear meshes used for convergence studies; a) medium tetrahedral mesh (9220 elements) b) fine tetrahedral mesh (26807 elements), and c) fine hexahedral mesh (5000 elements)

both low and high order polynomial interpolations, the expected rate of convergence is achieved for both primary ($p+1$ convergence rate) and derived (p convergence rate) variables. Notice that for this study, the multi-precision floating point library (MPFR) is utilised for arbitrary floating point accuracy. Apart from pinpointing the optimality and accuracy of the high order electromechanical solver on complex curved meshes, the h -convergence plots also prove that the framework scales reliably with high p and millions of degrees of freedom. Next, the p -convergence of tetrahedral elements for a fixed mesh (the finest mesh) is considered and compared to the p -convergence properties of hexahedral elements. Notice from [Table 3](#) and [Table 4](#) that the hexahedral mesh is chosen such that for every p , the number of nodes in the tetrahedral and hexahedral meshes are within a 5% difference. Once again, the degree of interpolation is successively increased leading to p convergence of the $\mathcal{L}^2(V)$ norm of the same afore-mentioned quantities. [Figure 7](#) shows p convergence of the $\mathcal{L}^2(V)$ for arguments of the set $\{\mathbf{x}, \varphi, \mathcal{V}, \Sigma_{\mathcal{V}}\}$. Once again, optimal rates of convergence for both curvilinear tetrahedral and hexahedral elements are obtained.

4.2. Effect of accurate boundary representation in nonlinear electro-elasticity

Having studied the convergence properties of high order curvilinear finite elements for modelling convex multi-variable electromechanics in the previous section, the objective of this section is to examine if accurate geometrical representation through high order curvilinear finite elements, does play a role in capturing the physics of soft elastomeric materials better compared to the standard high order finite elements with planer faces/edges. The examples in this section are also motivated by the fact that, the vast majority of finite element simulations of electromechanical devices are carried out using either low order or high order finite elements with planer faces [[54](#), [55](#), [56](#), [2](#), [16](#), [23](#), [34](#)], with possible geometrical simplifications and at times even de-featuring [[57](#)].

To this end, the notable example of electromechanical plate with a hole is chosen for examination [[54](#)]. For this example, once again, one tetrahedral mesh and one hexahedral mesh is chosen and the polynomial degree is successively enriched, while keeping the mesh size h fixed. Additionally, for both examples, the material model chosen is based on the following convex multi-variable strain energy representation

$$W_{el,2} = \mu_1 I_C + \mu_2 I_G - 2(\mu_1 + \mu_2) \ln \sqrt{C} + \frac{\lambda}{2} (\sqrt{C} - 1)^2 + \frac{1}{2\varepsilon_1} II_{D_0} + \frac{1}{2\sqrt{C}\varepsilon_2} II_d, \quad (41)$$

with material constants as given in [Table 5](#). The Poisson's ratio corresponding to parameters μ_1 , μ_2 and λ , is $\nu = 0.357$. The mesh quality information for the plate with the hole is listed in [Table 6](#)

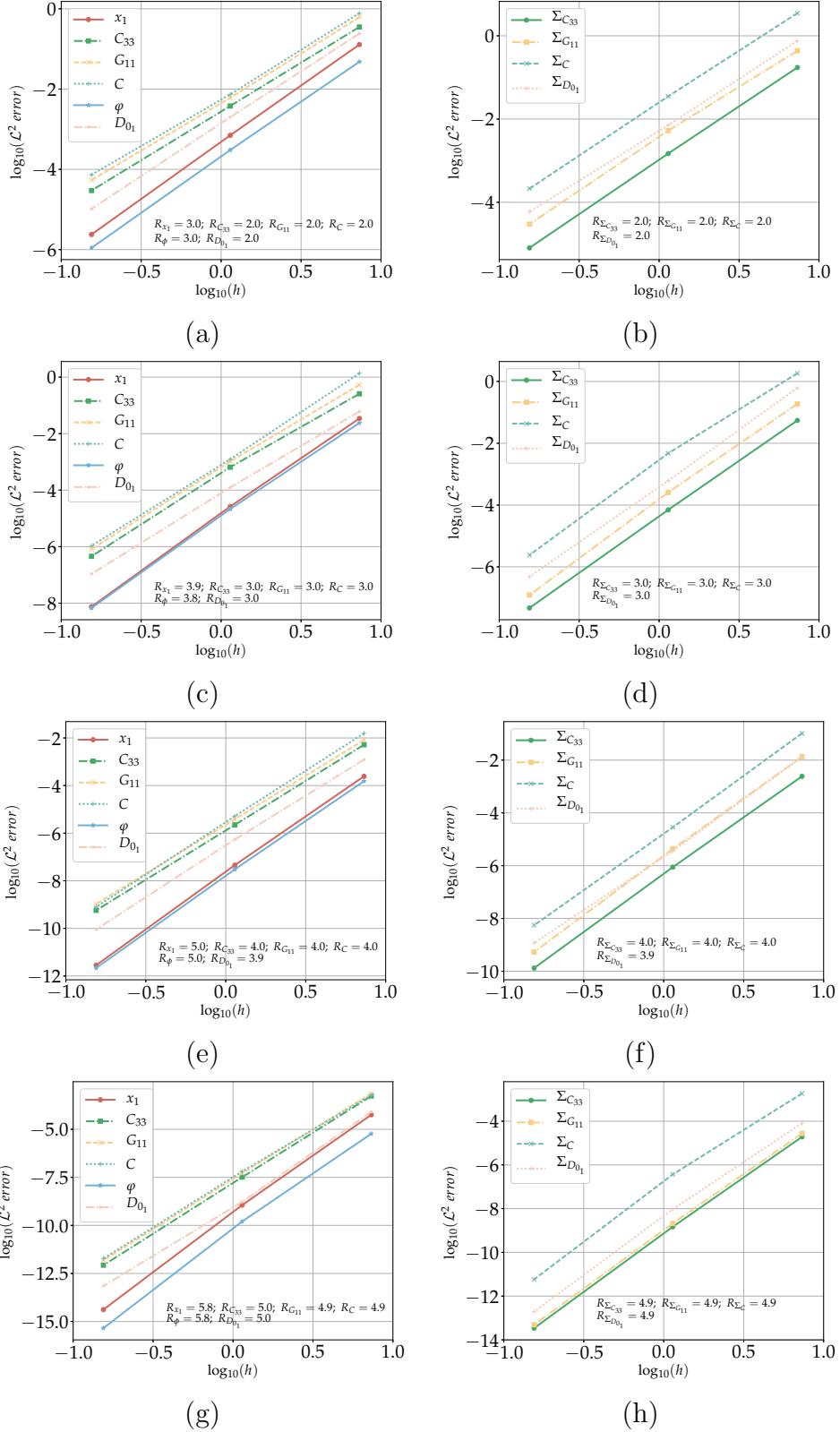


Figure 6: h convergence of $\mathcal{L}^2(V)$ norm of the error for different kinematic and kinetic variables, (a) & (b) $p = 2$, (c) & (d) $p = 3$, (e) & (f) $p = 4$ and (g) & (h) $p = 5$. R_ζ indicates the rate of convergence of quantity ζ .

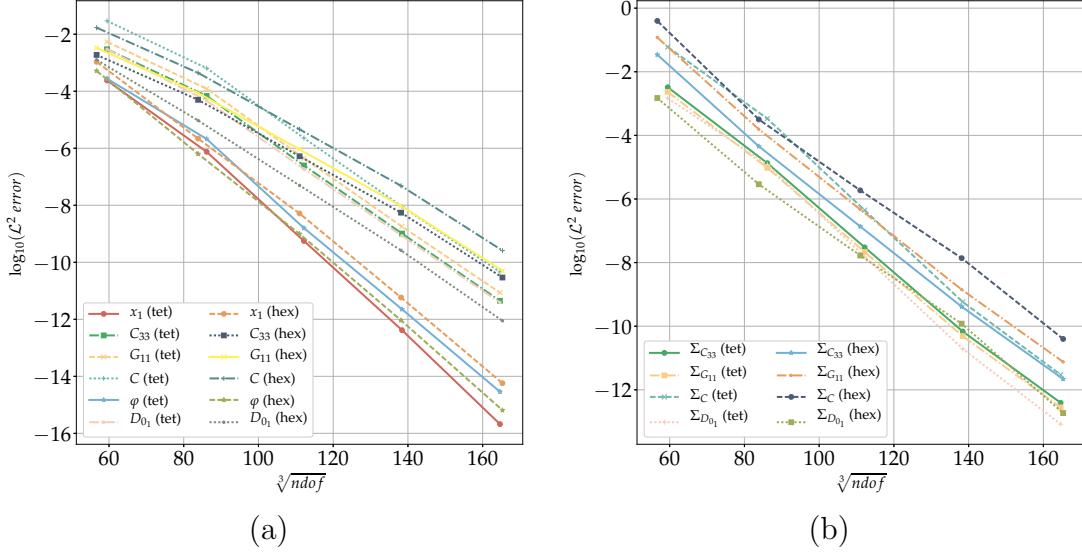


Figure 7: p convergence of $\mathcal{L}^2(V)$ norm of the error for different kinematic and kinetic variables for mesh with 9220 elements.

μ_1 (Pa)	μ_2 (Pa)	λ (Pa)	ε_1 (N/V^2)	ε_2 (N/V^2)
$1e5$	$1e5$	$5e5$	ε_0	$5\varepsilon_0$

Table 5: Material parameters for (41) with the vacuum permittivity $\varepsilon_0 = 8.85418781 \times 10^{-12}$ (N/V^2)

and the geometry and the tetrahedral and hexahedral meshes are shown in Figure 8 and Figure 9, respectively. A constant electric voltage of 2×10^8 V/m is applied through the whole thickness

p	Tetrahedral			Hexahedral				
	DoFs	Q_1	Q_2	Q_3	DoFs	Q_1	Q_2	Q_3
$p/q = 2$	340×4	0.982	0.965	0.948	363×4	0.980	0.961	0.942
$p/q = 3$	938×4	0.990	0.980	0.971	1012×4	0.982	0.966	0.951
$p/q = 4$	2021×4	0.986	0.973	0.960	2165×4	0.981	0.963	0.947
$p/q = 5$	3681×4	0.990	0.980	0.970	3966×4	0.979	0.959	0.941
$p/q = 6$	6097×4	0.990	0.979	0.969	6559×4	0.977	0.955	0.935

Table 6: Distortion quality of high order curvilinear tetrahedral and hexahedral meshes

and symmetric mechanical Dirichlet boundary conditions are imposed on the outer boundaries of the plate. This induces stretching of the electromechanical plate and as a result of the outer boundaries being fixed the plate is thickened in the region away from the hole and shrunk in thickness in the vicinity of the hole. Figure 10 shows the quadratic convergence of Newton-Raphson for the last load increment for all polynomial degrees on curvilinear tetrahedral and hexahedral meshes, respectively. It is worth mentioning that, since the planar and curvilinear meshes do not possess the same volume, a systematic study of the difference in error norms of quantities is not feasible. However, since both planar and curved meshes share the same ($p/q = 1$) vertices, a comparison of certain quantities of interest at these vertices can be carried out. Notice that since the node is shared between neighbouring elements appropriate stress recovery is required. This will lead to some oscillatory results in stresses (c.f. subsection 4.4 for further investigation in this regard). In the current setting, the order of quadrature rule to integrate stresses have been purposefully increased to $2(p + 1)$ to remove these oscillations as far as feasible. A representation of stress concentration for the plate with circular hole is shown

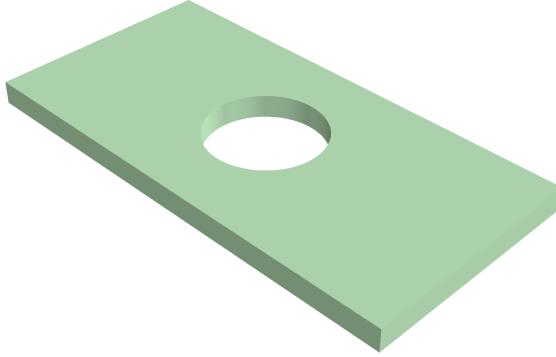


Figure 8: CAD representation of plate with circular hole with dimensions $10 \times 20 \times 2 \text{ m}^3$. The circular hole with a radius of 5m is centred at $[0, 0, 0]^T$

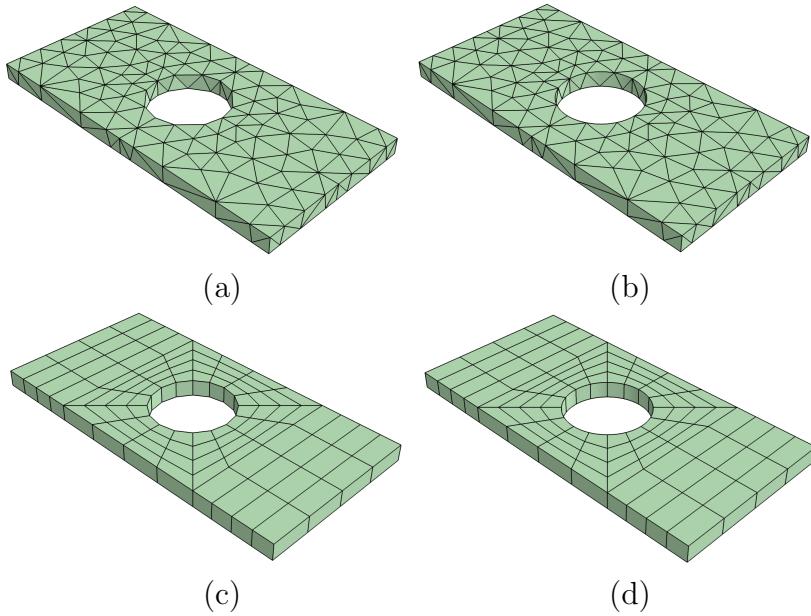


Figure 9: Curvilinear meshes used for comparison a) $p = 3$ planar tetrahedra, b) $p = 3$ curvilinear tetrahedra, c) $p = 3$ planar hexahedra and, d) $p = 3$ curvilinear hexahedra

figuratively in Figure 11 only for $p/q = 5$ meshes, for the final deformed configuration. It is evident from the figures that as opposed to the meshes with planar faces/edges, curvilinear meshes perform much better in reporting a smoother representation of the stress near the circular region. A similar conclusion is drawn across all polynomial degrees for tetrahedral and hexahedral elements (not reported, for the purpose of brevity). For instance, the maximum hydrostatic pressure at the tip of the circular hole within the plate located at $[5, 20, 2]^T$ is $p_{hyd} = -4.982 \times 10^4 \text{ Pa}$ for the hexahedral mesh with planar faces/edges, whereas only $p_{hyd} = -0.104 \text{ Pa}$ for curvilinear hexahedral mesh. This clearly confirms that the appearance of non-physical stress concentrations can be significantly reduced through curvilinear meshes accurately representing the geometry. Similar behaviour has also been reported by the computational fluid dynamics community, where an inaccurate geometric representation leads to non-physical entropy production [58, 57]. In line with the theme of this work, the aim of this study has been to assess and report the performance of curvilinear meshes in comparison to high order meshes with planar facets purely based on

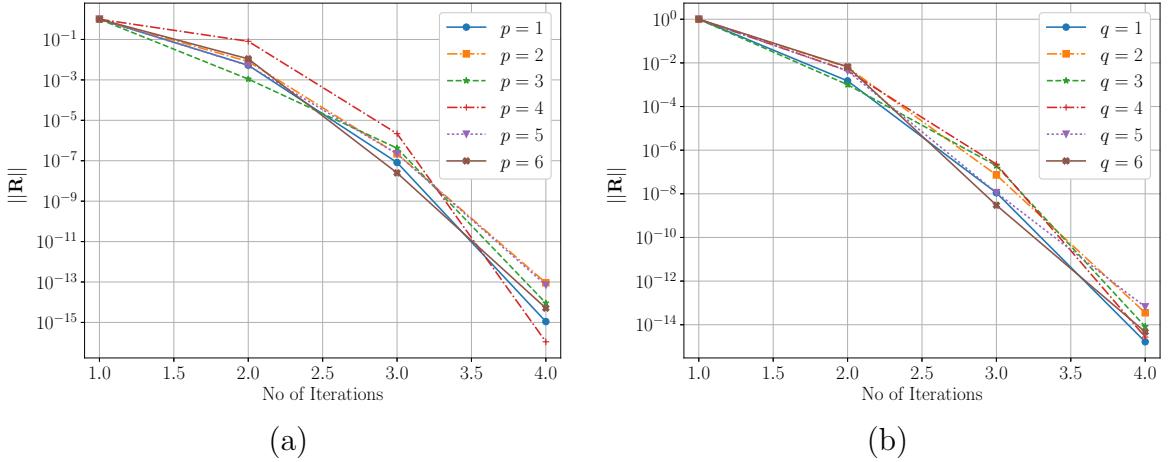


Figure 10: Quadratic convergence of Newton-Raphson on curvilinear a) tetrahedral mesh and, b) hexahedral mesh

geometrical enhancements.

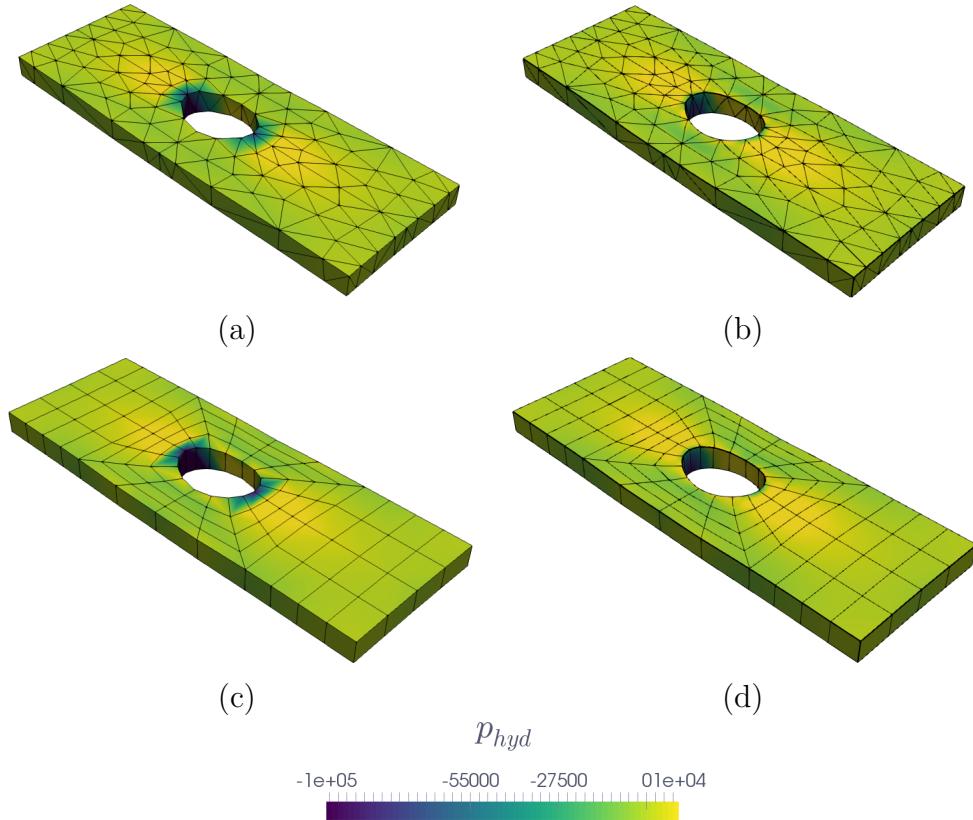


Figure 11: Evolution of voltage induced hydrostatic pressure in a plate with circular hole, a) planar $p = 5$ mesh, b) curved $p = 5$ mesh, c) planar $q = 5$ mesh and, d) curved $q = 5$ mesh

4.3. Linearised electrostriction through the staggered approach

The objective of this example is to study the performance and the range of applicability of the staggered approach for small strain actuation and energy harvesting problems. To this end,

two actuation problems are considered, based on two material models. For the first problem, the material model used is that of (41) with material constant given by (5). This problem is in particular a simple one due to the fact that electric field is a linear function of electric displacement and hence chosen for the purpose of benchmarking the staggered scheme. For the second problem, the following stabilised convex multi-variable energy functional is considered [3]

$$W_{el,2} = \mu_1 I_C + \mu_2 I_G - (2\mu_1 + 4\mu_2 + 12\mu_e) \ln \sqrt{C} + \frac{\lambda}{2} (\sqrt{C} - 1)^2 + \frac{1}{2\varepsilon_1} II_{D_0} + \frac{1}{2\varepsilon_2 \sqrt{C}} II_d \quad (42)$$

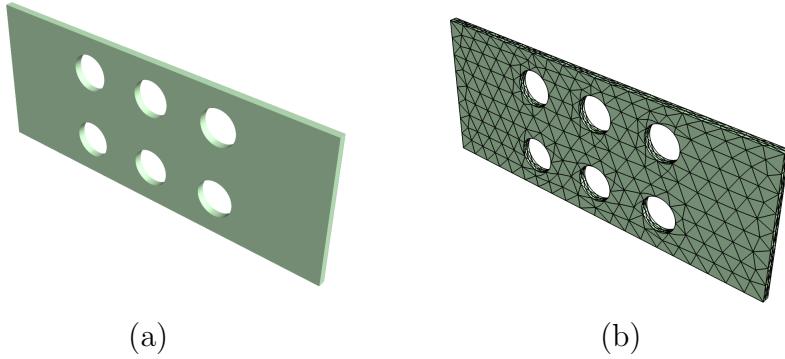
$$+ \mu_e \left(I_C^2 + \frac{2}{\mu_e \varepsilon_e} I_C II_d + \frac{1}{\mu_e^2 \varepsilon_e^2} II_d^2 \right),$$

with material properties listed in [Table 7](#). In this model the electric field and electric displacement

μ_1 (Pa)	μ_2 (Pa)	μ_e (Pa)	λ (Pa)	ε_1^{-1} (V^2/N)	ε_2 (N/V^2)	ε_2 (N/V^2)
$2.3e4$	$1.15e5$	800	$8e5$	0	$4.5\varepsilon_0$	$1050\varepsilon_0$

[Table 7](#): Material parameters for (42) with the vacuum permittivity $\varepsilon_0 = 8.85418781 \times 10^{-12}$ (N/V^2)

are coupled implicitly and hence this model tests the true performance of the staggered approach. The computational meshes considered are shown in [Figure 12](#) together with the CAD geometry. The mesh has 3819 elements and for an interpolation degree $p = 3$ there are 47317 points in the mesh corresponding to a total of 189268 degrees of freedom. To study the performance of the



[Figure 12](#): Electromechanical plate with holes a) CAD geometry and, b) $p = 4$ curvilinear mesh

staggered scheme the error incurred in staggered scheme is quantified relative to (fully nonlinear) monolithic approach as

$$\|e\|_{\mathcal{L}^2}^{\text{staggered}} = \left[\frac{\|\boldsymbol{\eta}_{\text{monolithic}} - \boldsymbol{\eta}_{\text{staggered}}\|}{\|\boldsymbol{\eta}_{\text{monolithic}}\|} \right]^{\frac{1}{2}}. \quad (43)$$

Typically, the range of applicability of the staggered electrostrictive scheme can be determined by successively increasing the electric voltage and observing the error produced by the scheme [25]. In the current setting, the following non-dimensional quantity is chosen to quantify the

increase in electric voltage

$$\mathcal{W} = \Phi / (H \sqrt{\frac{\mu}{\varepsilon}}), \quad (44)$$

where Φ denotes the applied potential and H represents the plate thickness. For the ideal dielectric model μ and ε can be related to material parameters as

$$\mu = 2(\mu_1 + \mu_2); \quad \frac{1}{\varepsilon} = \frac{1}{\varepsilon_1} + \frac{1}{\varepsilon_2}. \quad (45)$$

Similarly, for the stabilised material described in (42), μ and ε can be found as

$$\mu = 2(\mu_1 + 2\mu_2 + 6\mu_e); \quad \frac{1}{\varepsilon} = \frac{1}{\varepsilon_1} + \frac{1}{\varepsilon_2} + \frac{12}{\varepsilon_e}. \quad (46)$$

The plate is clamped at the two ends along the length and the electric voltage is applied through the thickness of the plate such that \mathcal{W} increases from 0.1 to 1.0. Figure 13 shows the relative \mathcal{L}^2 norm of the staggered approach as function of number of increments for different values of \mathcal{W} . As can be observed, certainly as the \mathcal{W} and hence the voltage induced deformation increases, the error incurred in staggered approach increases. The convergence of the staggered approach is approximately linear for both models. Interestingly, even under the application of extremely high voltage, the results of the staggered scheme seem to converge to the results of the monolithic approach. However, opting for such high number of increments may not be computationally practical. For engineering accuracy and large scale problems, the staggered scheme pays off as it only involves an iterative solution of the Gauss's law for the scalar electric potential.

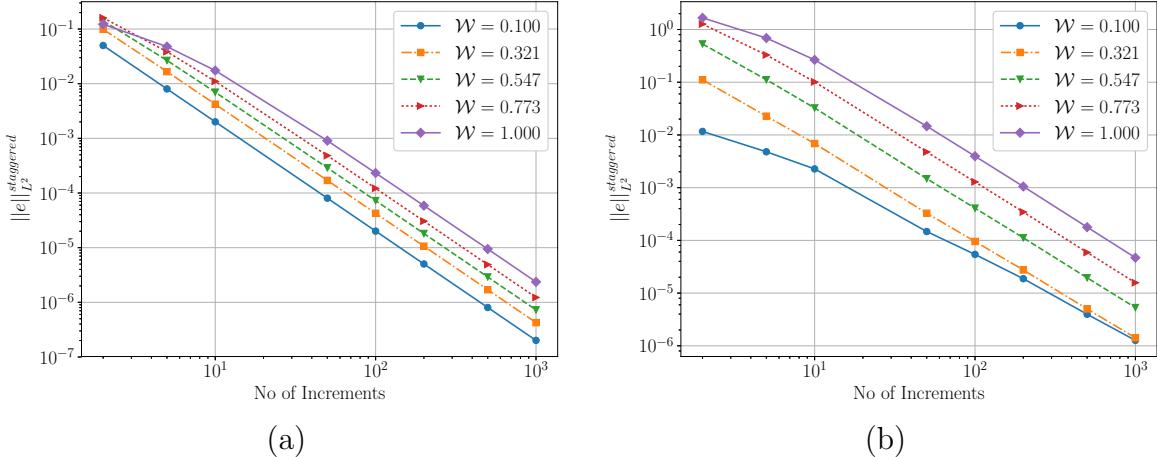


Figure 13: Convergence of the staggered scheme with respect to the monolithic solver as a function of increase in applied voltage quantified by \mathcal{W} as defined in (44)

Finally, Figure 14 shows the amount of voltage induced strains in the plate. Notice that despite large displacements, the strains are still small in the context of dielectric elastomers.

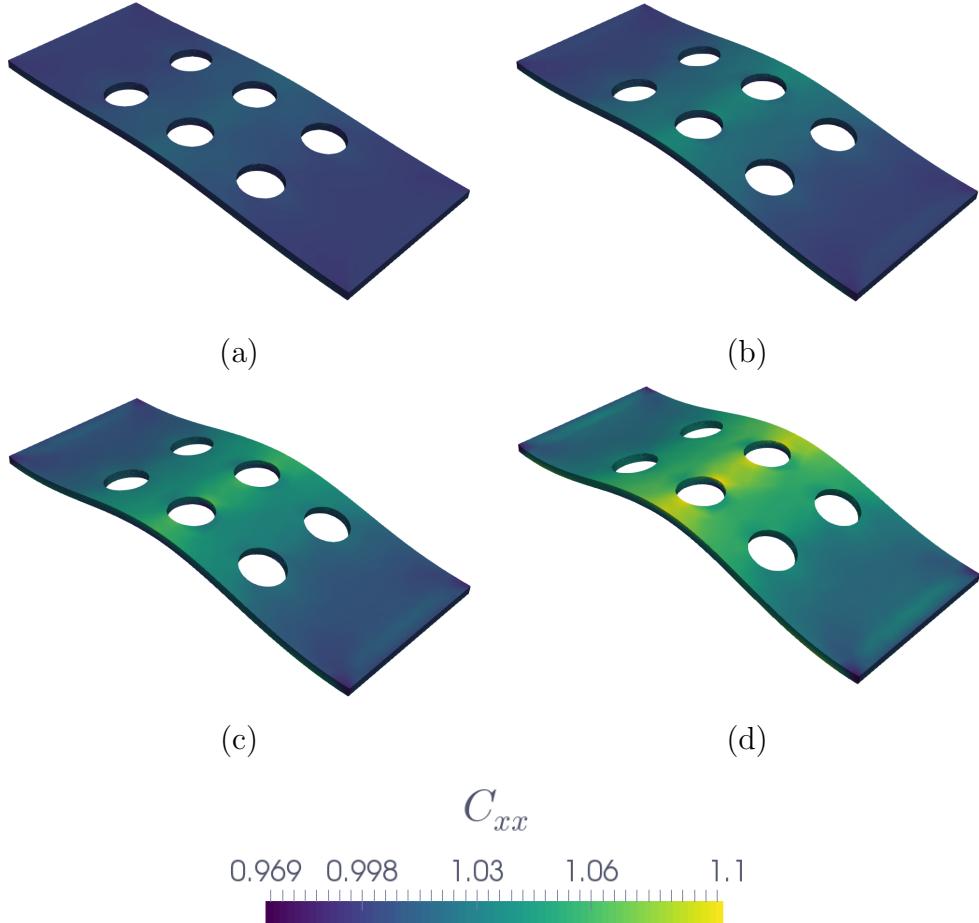


Figure 14: Voltage induced strains C_{xx} in electromechanical plate

4.4. Comparison of higher order displacement potential based formulation with 11 field mixed formulation for electro-elasticity

The objective of the examples presented in this section is to compare the performance of the displacement-potential formulation (**DPF**) discretised using high order tetrahedral and hexahedral finite elements with the 11 field mixed Hu-Washizu variational formulation (**MWF**) (based on the set $\{\mathbf{x}, \mathbf{F}, \mathbf{H}, J, \varphi, \mathbf{d}, \Sigma_{\mathbf{F}}, \Sigma_{\mathbf{H}}, \Sigma_J, \Sigma_{D_0}, \Sigma_d\}$) for electro-elasticity presented in [1, 2, 16], on curvilinear meshes. In particular, **MWF** is an extension of the mixed formulations to polyconvex elasticity presented in [41, 22, 62]. While for the **DPF** formulation, an equal order interpolation for all the variables (\mathbf{x}, φ) is used (for tetrahedra and hexahedra), the **MWF** formulation utilises a quadratic interpolation for \mathbf{x} and φ , linear discontinuous interpolation for $\mathbf{F}, \mathbf{H}, J, \mathbf{d}, \Sigma_{\mathbf{F}}, \Sigma_{\mathbf{H}}, \Sigma_{D_0}, \Sigma_d$ and a piecewise constant interpolation for J and Σ_J and is primarily developed for 10-noded tetrahedra [2]. These formulations are compared against with an eye on two fundamental issues pertaining to the modelling of large deformation large electric field electromechanics namely, shear-locking and volumetric locking. To this end, three dielectric cantilever patches with aspect ratio $\{10, 100, 1000\}$ are considered, as shown in Figure 15. The problem is analysed under compressible and nearly incompressible scenarios with Poisson's ratio 0.45 and 0.499 respectively. A series of curvilinear meshes are then produced for $p = \{2, 3, 4, 5\}$ and $q = \{2, 3, 4, 5\}$ for each aspect ratio. In addition, three h -refinement levels are considered

using a coarse, a medium and a fine description for each aspect ratio and each polynomial degree. An attempt is made to keep the total number of nodes in the computational mesh for every polynomial degree within the same range (within 5% difference). The necessary information regarding the quality of the curvilinear meshes are listed in [Table 8](#) and [Table 9](#), for tetrahedral meshes and hexahedral meshes, respectively. Note that, for a given p and h -refinement level, the size of the computational mesh is the same for all aspect ratios, in that, to obtain different aspect ratios, the mesh is appropriately scaled in the thickness direction. This certainly leads to extremely stretched elements but does not change the distortion of quality of the mesh as far as curvilinear mesh quality measures are concerned.

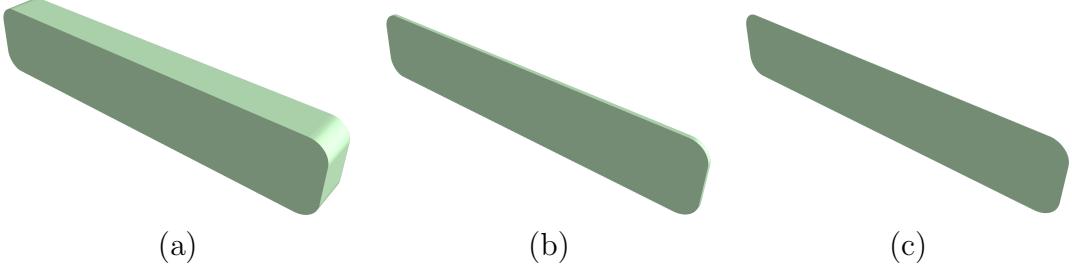


Figure 15: Geometry of the patch for three different aspect ratios a) 10, b) 100 and, c) 1000

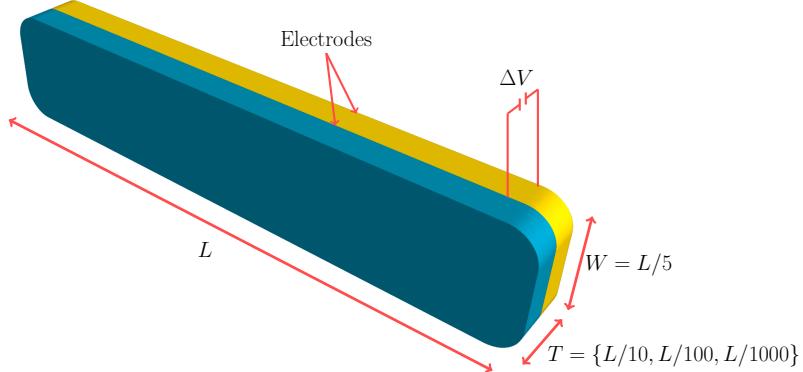


Figure 16: Description of the boundary conditions

p	Coarse				Medium				Fine			
	DoFs (N. Elements)	Q_1	Q_2	Q_3	DoFs (N. Elements)	Q_1	Q_2	Q_3	DoFs (N. Elements)	Q_1	Q_2	Q_3
$p = 2$	2842×4 (1652)	0.971	0.943	0.909	4686×4 (2730)	0.965	0.936	0.911	9916×4 (5910)	0.980	0.959	0.938
$p = 3$	2859×4 (396)	0.911	0.784	0.586	4656×4 (710)	0.925	0.837	0.734	9838×4 (1866)	0.952	0.898	0.836
$p = 4$	2835×4 (126)	0.852	0.690	0.500	4693×4 (326)	0.872	0.733	0.611	9819×4 (698)	0.945	0.886	0.819
$p = 5$	2856×4 (82)	0.873	0.758	0.651	4651×4 (172)	0.952	0.908	0.866	9889×4 (396)	0.944	0.890	0.840

Table 8: Distortion quality of high order curvilinear tetrahedral meshes

For comparison an ideal dielectric elastomer is considered with the following convex multi-variable internal energy description

$$W_{el,3} = \mu_1 I_C + \mu_2 I_G - 2(\mu_1 + \mu_2) \ln \sqrt{C} + \frac{\lambda}{2} (\sqrt{C} - 1)^2 + \frac{1}{2\varepsilon_1} II_d, \quad (47)$$

with the material parameters as given in [Table 10](#).

p	Coarse				Medium				Fine			
	DoFs (N. Elements)	Q_1	Q_2	Q_3	DoFs (N. Elements)	Q_1	Q_2	Q_3	DoFs (N. Elements)	Q_1	Q_2	Q_3
$q = 2$	2885×4 (244)	0.972	0.987	0.957	4665×4 (472)	0.986	0.970	0.953	9805×4 (928)	0.990	0.980	0.969
$q = 3$	2853×4 (76)	0.959	0.913	0.863	4605×4 (148)	0.979	0.957	0.935	9805×4 (292)	0.979	0.958	0.935
$q = 4$	2873×4 (32)	0.818	0.618	0.361	4650×4 (68)	0.950	0.897	0.842	9793×4 (124)	0.950	0.898	0.842
$q = 5$	2756×4 (16)	0.822	0.625	0.369	4700×4 (32)	0.834	0.656	0.401	9801×4 (64)	0.853	0.723	0.781

Table 9: Distortion quality of high order curvilinear hexahedral meshes

Parameters	μ_1 (Pa)	μ_2 (Pa)	λ (Pa)	ε_1 (N/V ²)
Values (for $\nu = 0.45$)	1e5	0	9e5	$4\varepsilon_0$
Values (for $\nu = 0.499$)	1e5	0	4.99e7	$4\varepsilon_0$

Table 10: Material parameters for (47) with the vacuum permittivity $\varepsilon_0 = 8.85418781 \times 10^{-12}$ (N/V²)

In order to be able to apply appropriate boundary conditions, it is made sure that all computational meshes have at least two elements across the thickness. For every aspect ratio, the patch is clamped at $[0, X, Y]^T$ and a constant electric voltage is applied across the half thickness $T/2$ (where T represents the patch thickness), as shown in Figure 16. It is certainly not feasible to apply the same amount of electric voltage while varying the patch aspect ratio and hence Table 11 lists the Dirichlet boundary conditions considered for the three aspect ratios.

Aspect Ratio	10	100	1000
Electric Voltage (V/m)	1e7	7.5e4	1.8e2

Table 11: Applied electric voltage as Dirichlet boundary condition for three aspect ratios

In order to compare different kinematic and kinetic quantities of interest for different formulations, two different physical points in the patches are considered namely, A - the node located at $[L, W/2, T/2]^T$, and B - the point interior to the computational mesh $[0.98L, 0.49W, 0.49T]^T$ (c.f. Figure 16).

Table 12 and Table 13 compare the results obtained for different kinematics and kinetics quantities with displacement potential formulation using high order tetrahedral and hexahedral elements with the 11 field mixed formulations, for aspect ratio 10 and Poisson's ratio 0.45 and 0.499, respectively. For this comparison, the problem is solved in 50 load increments and the comparison is performed at increment 12, which corresponds to the onset of nonlinear deformation. The first observation from the tables is that the **DPF** based discretisations have an asymptotically upper bound convergence while the **MWF** based results have an asymptotically lower bound convergence. This can be observed studying the convergence of a given quantity for a fixed polynomial degree and successive h -refinements. This convergence pattern is more pronounced in the case of **MWF**. Consequently, for a fixed polynomial degree, it can be observed that **MWF** has a consistent 2%-5% higher displacements and strains estimate rate.

Regarding the stresses, as can be seen from the σ_{xz} results, the differences between the formulations are significant. While the results for the **MWF** show a convergence pattern, the results of the **DPF** do not seem to converge as coherently, with some clear pressure oscillations (as p/q and h -refinements are carried out). The σ_{zz} and hydrostatic pressure still do not seem to be affected, primarily due to the fact that these components have a much higher electro-mechanical stress contribution coming from the electric displacement (which is the applied Dirichlet boundary condition).

The conclusion drawn from the above set of results is that, for thick electromechanical beams, both high order **DPF** and **MWF** implementations capture the physical behaviour of the system

accurately without showing any signs of shear or volumetric lockings. Importantly, this signifies that high order discretisations can capture nearly incompressible scenarios reliably [59, 60, 61]. The oscillations in stresses is an expected phenomenon for **DPF** implementations [2, 62, 41] which seems to persist even with polynomial enrichment. Unlike the problems reported in [subsection 4.1](#), the convergence rate of stresses here can not be ascertained, since the polynomial enrichment here, comes with the trade-off of coarsening the computational mesh to keep the problem size the same. In this regard, **MWF** implementation maintains a consistent accuracy by virtue of explicitly solving for kinetic variables.

Note that so far, only the aspect ratio of the patch has been considered and no remarks have been made regarding the aspect ratio of the elements in the mesh. It is important to mention that high order elements are at a severe disadvantage here compared to their low-order or **MWF** counterparts, in that, polynomial enrichment leading to coarsening of the mesh, also inherently implies dramatic stretching for high order elements. It is certainly not feasible to control the aspect ratio of the patch and the aspect ratio of the elements at the same time, while also having to keep the size of the computational mesh the same across all polynomial enrichments. This also signifies that, for $p = q$, the computational meshes do not necessarily exhibit the same elemental aspect ratios. It is still astounding to observe that, high order elements perform fairly well regardless.

Quantity	Refinement	$p = 2$	$p = 3$	$p = 4$	$p = 5$	$q = 2$	$q = 3$	$q = 4$	$q = 5$	11Mixed
$u_x^A(m)$	Coarse	0.15	0.1506	0.1502	0.1505	0.1505	0.1506	0.1506	0.1506	0.1505
	Medium	0.1505	0.1503	0.1505	0.1506	0.1506	0.1507	0.1506	0.1506	0.1542
	Fine	0.1505	0.1508	0.1505	0.1506	0.1506	0.1507	0.1506	0.1506	0.1535
$u_z^A(m)$	Coarse	0.9572	0.9555	0.9565	0.9551	0.958	0.9597	0.9598	0.9603	1.049
	Medium	0.9566	0.9576	0.9603	0.9596	0.9583	0.96	0.9604	0.9603	1.0022
	Fine	0.9571	0.96	0.9597	0.9594	0.9585	0.96	0.9604	0.9603	0.9948
F_{xz}^A	Coarse	-0.0176	-0.0186	-0.018	-0.0184	-0.0178	-0.0184	-0.0183	-0.0183	-0.0199
	Medium	-0.0183	-0.0181	-0.018	-0.0183	-0.0178	-0.0185	-0.0183	-0.0183	-0.0191
	Fine	-0.0183	-0.0186	-0.0178	-0.0184	-0.0178	-0.0185	-0.0183	-0.0183	-0.0191
H_{zx}^A	Coarse	-0.0184	-0.0184	-0.0182	-0.0186	-0.0181	-0.0184	-0.0183	-0.0183	-0.0198
	Medium	-0.0184	-0.0181	-0.0183	-0.0182	-0.018	-0.0185	-0.0184	-0.0183	-0.019
	Fine	-0.0184	-0.0181	-0.0186	-0.0184	-0.018	-0.0185	-0.0184	-0.0183	-0.0192
$\sigma_{xz}^B(\text{Pa})$	Coarse	68.817	-34.4955	8.4401	2.5424	20.9594	-16.3045	-4.2272	-8.6977	-24.0193
	Medium	1.5908	-10.4219	16.6441	-17.0173	11.9318	-4.436	-4.8418	-8.9388	-15.3804
	Fine	-3.9871	-63.5881	19.221	-15.5529	7.1833	-4.433	-4.8497	-8.9573	-4.7419
$\sigma_{zz}^B(\text{Pa})$	Coarse	228.8202	242.9458	167.4105	225.4981	227.3906	208.8909	220.5269	233.9105	262.6845
	Medium	226.3231	304.8385	190.6815	232.3852	228.7893	202.9177	228.5124	234.1146	262.0383
	Fine	226.7235	228.4325	192.3997	227.9372	229.1137	202.8566	228.503	234.1104	233.9687
$p_{hyd}^B(\text{Pa})$	Coarse	258.7676	262.7815	228.257	256.8133	249.9223	258.8183	253.0553	260.2188	283.1432
	Medium	256.2272	239.7233	239.8217	256.7475	249.6441	259.3416	254.3613	260.1902	279.219
	Fine	257.2179	258.3805	255.9474	256.809	249.4617	259.3333	254.3594	260.1998	229.3169

Table 12: Comparison of high order displacement potential and 11 field mixed formulations for different kinematic and kinetic measures, for aspect ratio 10 and Poisson's ratio 0.45

Turning the attention to thinner beams, the results for the patch with the aspect ratio 100 are analogous to those with the aspect ratio 10 and offer no further insight, and are hence omitted for the purpose of brevity. The same analysis is then carried out with extremely thin beams with aspect ratio 1000 featuring highly stretched elements. The corresponding loading from [Table 11](#) is applied. [Table 14](#) and [Table 15](#) compare the results obtained for different kinematics and kinetics quantities with displacement potential formulation using high order tetrahedral and hexahedral elements with the 11 field mixed formulations, for aspect ratio 1000 and Poisson's ratio 0.45 and 0.499, respectively.

As can be observed from the tables, the $p = 2$ elements exhibit severe shear locking in this case. Comparing the results of thick and thin beams, it is evident that this phenomenon occurs solely due to the high aspect ratio (slenderness of the beam) [63, 64]. Interestingly, although not as dramatic, the **MWF** implementation also exhibits shear locking. On the other hand, hexahedral

Quantity	Refinement	$p = 2$	$p = 3$	$p = 4$	$p = 5$	$q = 2$	$q = 3$	$q = 4$	$q = 5$	11Mixed
$u_x^A(m)$	Coarse	0.0022	-0.0329	-0.0563	-0.0564	-0.0603	-0.0618	-0.0607	-0.0618	0.0004
	Medium	0.0022	-0.0489	-0.0559	-0.0574	-0.0642	-0.0615	-0.0615	-0.0605	-0.0194
	Fine	0.0022	-0.0434	-0.0577	-0.0582	-0.0614	-0.0616	-0.0615	-0.6165	-0.0312
$u_z^A(m)$	Coarse	0.0062	2.1839	2.9353	3.0931	3.1536	3.2548	3.2635	3.2618	0.5988
	Medium	0.0024	2.496	2.9301	2.9984	3.2623	3.2082	2.9336	3.2518	1.8414
	Fine	0.0104	2.8819	2.933	2.9997	3.2017	3.2099	3.1395	3.2548	2.3124
F_{xz}^A	Coarse	-0.0007	-0.0441	-0.0505	-0.0515	-0.0626	-0.0691	-0.0603	-0.0621	-0.0004
	Medium	-0.0002	-0.0565	-0.0564	-0.0526	-0.0643	-0.063	-0.0592	-0.0044	-0.0001
	Fine	-0.0002	-0.0501	-0.0561	-0.0582	-0.063	-0.063	-0.0613	-0.0593	-0.0004
H_{zx}^A	Coarse	-0.0007	-0.0441	-0.0205	-0.0515	-0.0627	-0.0691	-0.0604	-0.0621	-0.0448
	Medium	-0.0002	-0.0565	-0.0565	-0.0527	-0.0643	-0.063	-0.0592	-0.0044	-0.0115
	Fine	-0.0002	-0.0501	-0.0561	-0.0583	-0.063	-0.063	-0.0613	-0.0593	-0.038
$\sigma_{xz}^B(\text{Pa})$	Coarse	-0.058	-1.0561	0.2593	0.191	0.7204	0.7001	0.7089	0.7071	-0.2034
	Medium	-0.0656	-1.022	5.1248	-39.3918	0.8225	0.5512	0.3729	-9.6306	0.6378
	Fine	-0.1293	-0.8051	10.014	28.8039	0.8635	0.5506	0.364	-206.3348	0.2668
$\sigma_{zz}^B(\text{Pa})$	Coarse	47.1482	-395.4936	-33.2631	538.5343	-2.8369	3.3726	3.3814	3.3796	2.1594
	Medium	0.6649	49.3672	-414.2999	2113.4412	-4.1608	6.1501	1.802	2217.1176	4.6069
	Fine	-0.6984	13.3289	330.097	2189.0055	-5.4436	6.173	1.4294	3478.8777	14.6644
$p_{hyd}^B(\text{Pa})$	Coarse	38.1504	-407.0878	-41.0376	528.9243	0.4472	6.5811	6.5899	6.5881	-5.0029
	Medium	-7.9517	40.8982	-420.7844	985.5186	-0.8781	9.456	5.1897	739.779	6.6814
	Fine	-9.1691	5.7871	350.556	710.8945	-2.1342	9.4748	4.8289	1165.274	15.8006

Table 15: Comparison of high order displacement potential and 11 field mixed formulations for different kinematic and kinetic measures, for aspect ratio 1000 and Poisson’s ratio 0.499

at the expense of a huge oscillation in derived variables such as stresses. To separate the issue of volumetric and shear locking further a different set of experiments are now considered using only **DPF** implementation by applying an extremely high electric voltage on the patches and lowering the Poisson’s ratio to $\nu = 0.35$. The loading scenario for different aspect ratios is listed in Table 16.

Aspect Ratio	10	100	1000
Electric Voltage (V/m)	$2.6e7$	$1e5$	$4e2$

Table 16: Applied electric voltage as Dirichlet boundary condition for three aspect ratios

Figure 17 shows the hydrostatic pressure for the patch with aspect ratio 10 (thick) solved with the medium mesh. Once again, it can be seen that even for large deformations and finite strains, the **DPF** implementation with tetrahedra does not suffer from any locking, although some oscillation in the pressure could be observed, across all polynomial degrees. Similarly, Figure 18 shows the evolution of the co-factor component H_{13} on the coarse mesh. For this aspect ratio, which could be considered a thin beam, **DPF** implementation with tetrahedra once again produces accurate results.

As evident by now, the challenging problem is the extremely thin beams with aspect ratio 1000 shown in Figure 18. Notice that both $p = 2$ and $p = 3$ elements exhibit severe shear locking for this aspect ratio, confirming the fact that this locking is purely due to stretching of the elements. Only $p = 4$ elements (and beyond) are able to capture this deformation accurately. Interestingly, as shown in Figure 20 the $q = 2$ elements also lock at this level of deformation. Certainly, the problem of shear locking could in part be remedied by using h -refinement. As shown in Figure 21, even with the fine mesh the $q = 2$ elements suffer from shear locking.

The examples presented here show a clear advantage for the application of high order elements in electromechanics. Modelling of dielectric elastomeric beams with high aspect ratio is indeed a challenging problem. Mixed formulations based on convex multi-variable electro-elastic strain energies have an unrivalled performance in capturing the pressure and other derived quantities of interest, but still could suffer from shear locking at extremely high aspect ratios. Note that the performance of high order elements could still be superior if the element’s aspect ratio was preserved during polynomial enrichment (which was not the case for here). The analyses perfor-

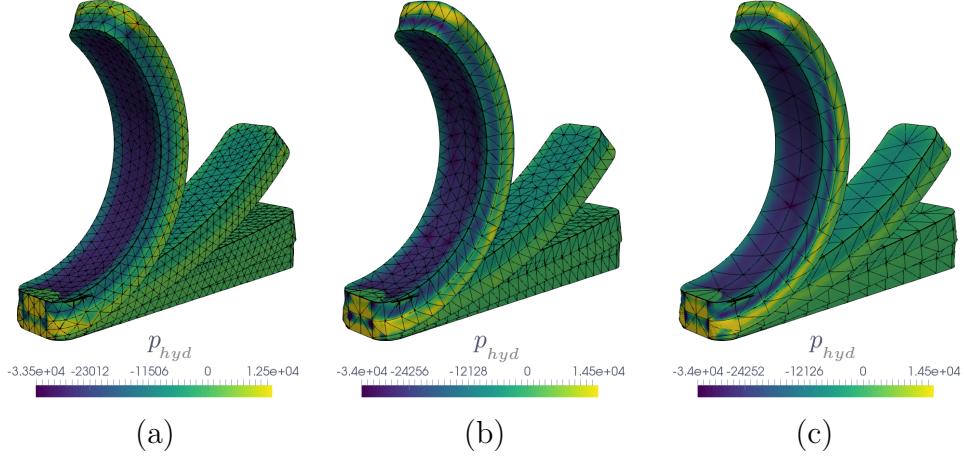


Figure 17: Voltage induced hydrostatic pressure p_{hyd} for the medium mesh using a) $p = 2$, b) $p = 3$ and, c) $p = 4$

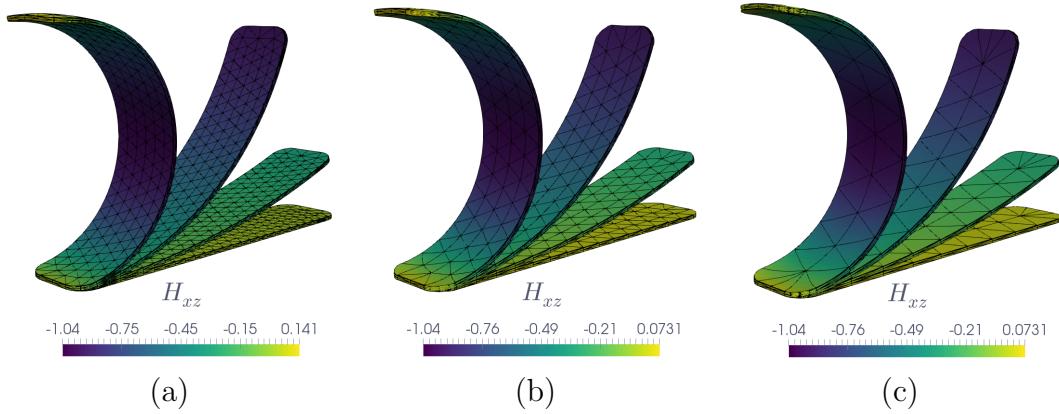


Figure 18: Voltage induced large strain (H_{xz}) for the coarse mesh using a) $p = 2$, b) $p = 3$ and, c) $p = 4$

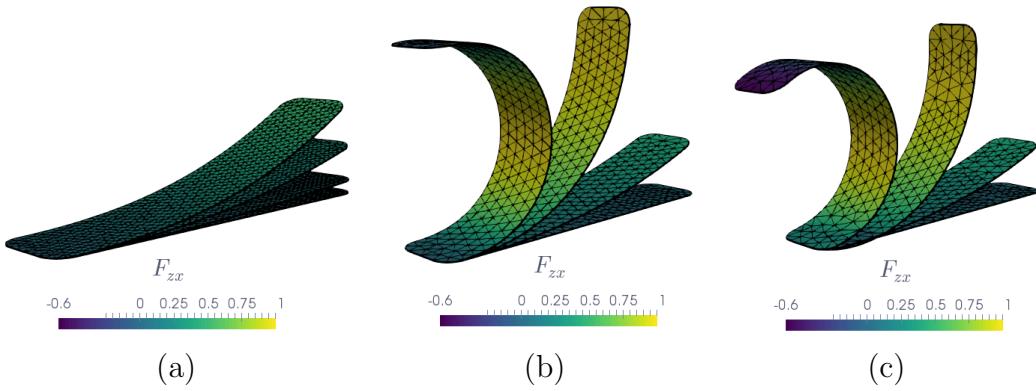


Figure 19: Voltage induced large strain (F_{zx}) for the fine mesh using a) $p = 2$, b) $p = 3$ and, c) $p = 4$

med in this section was carried out with the clear goal of keeping the size of the computational domain the same across all polynomial enrichments.

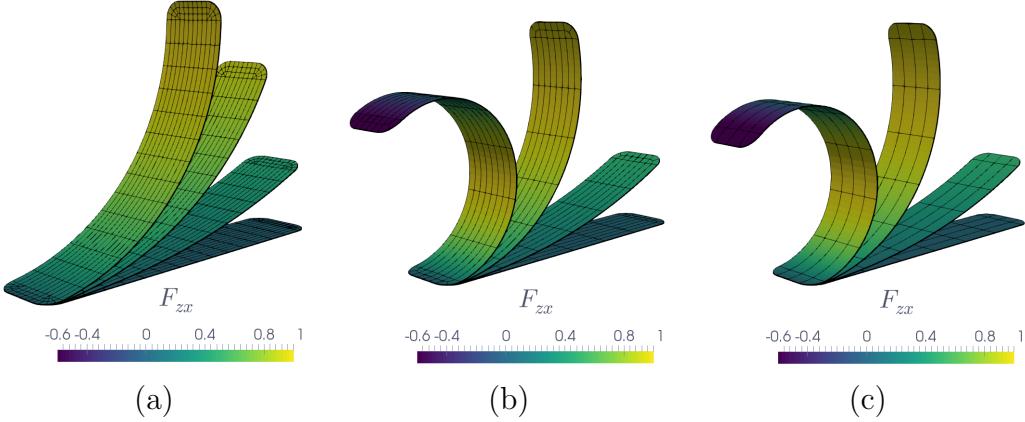


Figure 20: Voltage induced large strain (F_{zx}) for the fine mesh using a) $q = 2$, b) $q = 3$ and, c) $q = 4$

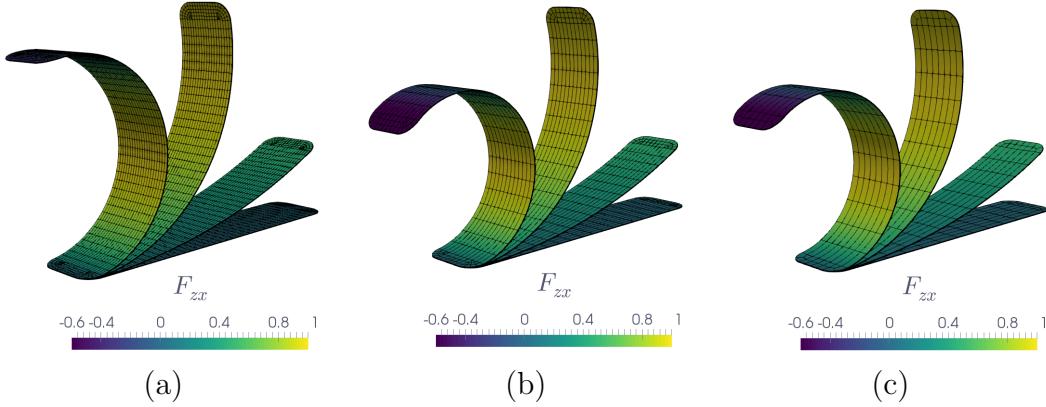


Figure 21: Voltage induced large strain (F_{zx}) for the fine mesh using a) $q = 2$, b) $q = 3$ and, c) $q = 4$

4.5. Kernel-based data parallel code generation for electromechanics on curvilinear meshes

Having studied the h and p convergence properties of the framework and its performance with respect to mixed variational principles, the objective of this section is to present some implementation details of the framework. The present high order curvilinear finite element framework is tied and closely developed on top of **Fastor** library [44]. **Fastor** (<https://github.com/romeric/Fastor>) is an open source (MIT licensed) C++ based generic tensor contraction library that can perform heavy numerical computations such as mathematical/algorithmic transformations and graph optimisations at compile time using a powerful in-built expression template engine. The transformed algorithms are then used by **Fastor** to generate data parallel (SIMD-vectorised) code for modern CPU and potentially GPU architectures. Recently, it has been shown in [44] that **Fastor** is able to generate carefully crafted data parallel code (SSE-SSE4.2, AVX-AVX2-AVX512, FMA) for the local assembly of low and high order finite elements for a series of convex multi-variable electro-elastic models, exhibiting many fold performance improvement over hand-written C code. As described earlier, since convex multi-variable electro-elasticity models require Legendre transformation through a Newton-Raphson scheme per quadrature point to find the appropriate constitutive tensors, local assembly of the finite elements become a critical hotspot for performance optimisations (see [38, 44] for performance benchmarks).

To this effect, a problem is manufactured where the performance of the current framework (using **Fastor**) is studied and compared to the equivalent hand-optimised C code. This problem

pertains only performance studies of numerical integration of work-conjugates (20), Hessian (22) and the subsequent quadrature point Newton-Raphson required for the Legendre transform (30) and (26), during local assembly (also see Algorithm 1). A set of similar benchmark problems has been presented in the authors' previous work on Fastor [44], where explicit code snippets for the aforementioned set of operations are presented. The interested reader is advised to consult Poya et. al. [44] and Fastor's repository, for a series of such performance studies, as for the purpose of brevity, complete implementation details of the problem are not presented here.

This performance study once again considers both curved hexahedral and tetrahedral meshes. The geometry considered for the problem is that of an electromechanical component shown in Figure 22 together with curved meshes. For the purpose of benchmark, refined meshes are chosen such that distortion qualities are all almost unity.

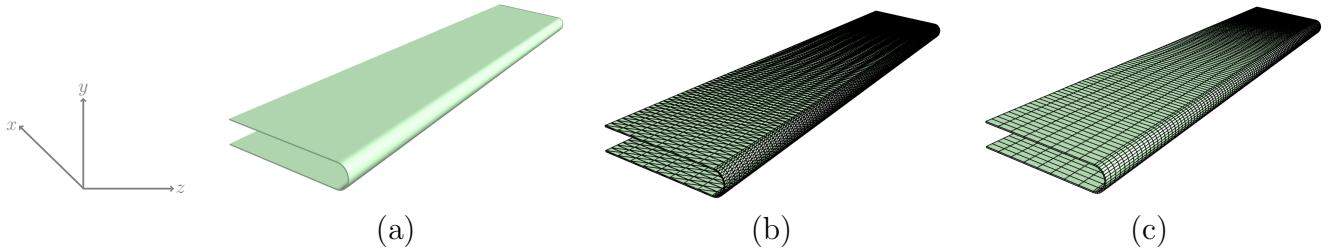


Figure 22: a) CAD geometry, b) curved tetrahedral mesh and, c) curved hexahedral mesh for electromechanical component with 1m thickness, 20m height (y -axis), 100m width (x -axis) and 500m length (z -axis)

Note that the tetrahedral and hexahedral meshes do not have the same number of nodes. The relevant information regarding the meshes is listed in Table 17. The geometry essentially consists of two plates connected through an extruded arc/fillet (in the z direction). No boundary conditions have been applied on the arc part of the geometry. The electromechanical component is mechanically fixed at one longitudinal end and two different electric voltages are applied across the thickness of the top and bottom plates, namely 4.5×10^7 V for the top plate and 8.1×10^7 V for the bottom plate. The problem exhibits massive snap-back behaviour as the different applied voltages make the component bend initially but after the point of electromechanical instability the difference in displacements in both plates causes the component to deform in the opposite direction.

Mesh (N. Elements)	$p/q = 2$	$p/q = 3$	$p/q = 4$	$p/q = 5$	$p/q = 6$	$p/q = 7$	$p/q = 8$
Tetrahedral (43200)	367196	1223568	2689668	5082200	8366140	12921312	18777812
Hexahedral (7200)	351276	1048144	2322900	4348344	7297276	11342496	16656804

Table 17: Degrees of freedom associated with each polynomial degree for tetrahedral and curvilinear meshes

To study the performance of numerical integration, the stabilised convex multi-variable model (42) is chosen with material properties listed in Table 7. The problem is then analysed under the aforementioned boundary conditions and the performance of numerical integration of work-conjugates Σ_V (20), Hessian and constitutive tensors of the internal energy \mathcal{C} , \mathcal{Q} , θ (22) and the quadrature point Newton-Raphson required for the Legendre transform (30) to obtain the constitutive tensors of Helmholtz-like energy \mathcal{C}_Φ , \mathcal{Q}_Φ , θ_Φ (26) are monitored under three different compilation flags. These correspond to the same code compiled with -O3 -march=native but with auto-vectorisation turned off (i.e. -fno-tree-vectorize for gcc, -fno-vectorize for

LLVM’s clang and `-no-vec` for Intel’s ICC), same code compiled with `-O3 -march=native` where the compiler is permitted to auto-vectorise the code and Fastor transformed code compiled with `-O3 -march=native`. All performance measurements have been recorded on a single core of Intel(R) Xeon(R) CPU E5-2650 v2 @2.60GHz with AVX instruction sets, 20MB private L1-cache and 32GB memory, running Ubuntu 16.04 and GCC-7, LLVM’s Clang 4.0 and ICC 17.0.3 are used to compile all the codes. To keep the comparison fair, careful attention is paid to the compilers generated assembly code.

Figure 23 shows the relative speed of the code (with and without auto-vectorisation) in comparison to Fastor’s generated code with all the three aforementioned compilers. Since with every polynomial enrichment the quadrature order has to be increased accordingly, as a result the computational cost of numerical integration increases almost cubically for all the cases (for three-dimensional problems). Nevertheless, Fastor’s data parallelism falling into the category of latency hiding techniques, hides much of this computational complexity. For both hexahedra and tetrahedra, Fastor’s SIMD vectorised code shows consistently nearly 2X speed-up over compilers auto-vectorised code on an AVX capable processor and 3-3.5X speed up over non-vectorised (but well-optimised) code. It can also be seen from the figures, that unlike Fastor, auto-vectorisation is compiler specific and in some cases the code may not benefit from it, specifically with Intel’s ICC in this case, Figure 23(e,f). On the other hand, Fastor is extremely less sensitive to the compiler (and optimisations) used. This is potentially the ultimate speed-up achievable for non-trivial codes using data parallelism on AVX architectures over non-vectorised codes (ideal speed-up is 4X for double precision). For complex algorithms, such as numerical integration of convex multi-variable electro-elastic models, compile time code transformation and graph search optimisation are necessary to be able to achieve this performance. One again, these aspects of Fastor library are presented and benchmarked in detail in [44].

Finally, Figure 24 shows the evolution of voltage induced hydrostatic pressure p_{hyd} at different loading stages, for $q = 3$ hexahedral mesh. Notice the massive snap-back behaviour of system and butterfly shape it occupies at the final deformed configuration.

4.6. Virtual prototyping of massive deformations and instabilities in dielectric elastomers through high performance numerical simulations

The objective of this section is to numerically study the behaviour of highly stretchable dielectric elastomers undergoing massive deformations through virtual prototyping of a series of electromechanical components. One of the main goals of these set of examples is to gain insights into the onset of instabilities such as pull-in instability [24, 25] and the subsequent formation of wrinkles in dielectric elastomers. The occurrence of these phenomena have been experimentally verified by Plante and Dubowsky [24], where a large applied electric potential has led to massive deformation of a dielectric elastomer sheet with eventual partial wrinkling of the sheet. Zhao et. al. [25], Mao [26] and others [27, 28, 65] have also reported instabilities in DEs through numerical studies, albeit in relatively simplified settings. Here, an attempt is made to model dielectric elastomeric components which could potentially be applied as compliant actuators in soft robotics, medical devices and similar applications. From a computational point of view, four ingredients are needed to simulate such massive deformations in DEs namely, a) a convex multi-variable expansion of the electromechanical internal energy ensuring ellipticity [1, 2, 3, 16] b) accurate representation of the dielectric elastomeric components, c) high order finite element analyses to capture extreme deformations and finally, d) high performance kernels for rapid prototyping and modelling DEs. Having established all the aforementioned necessities, a diligent

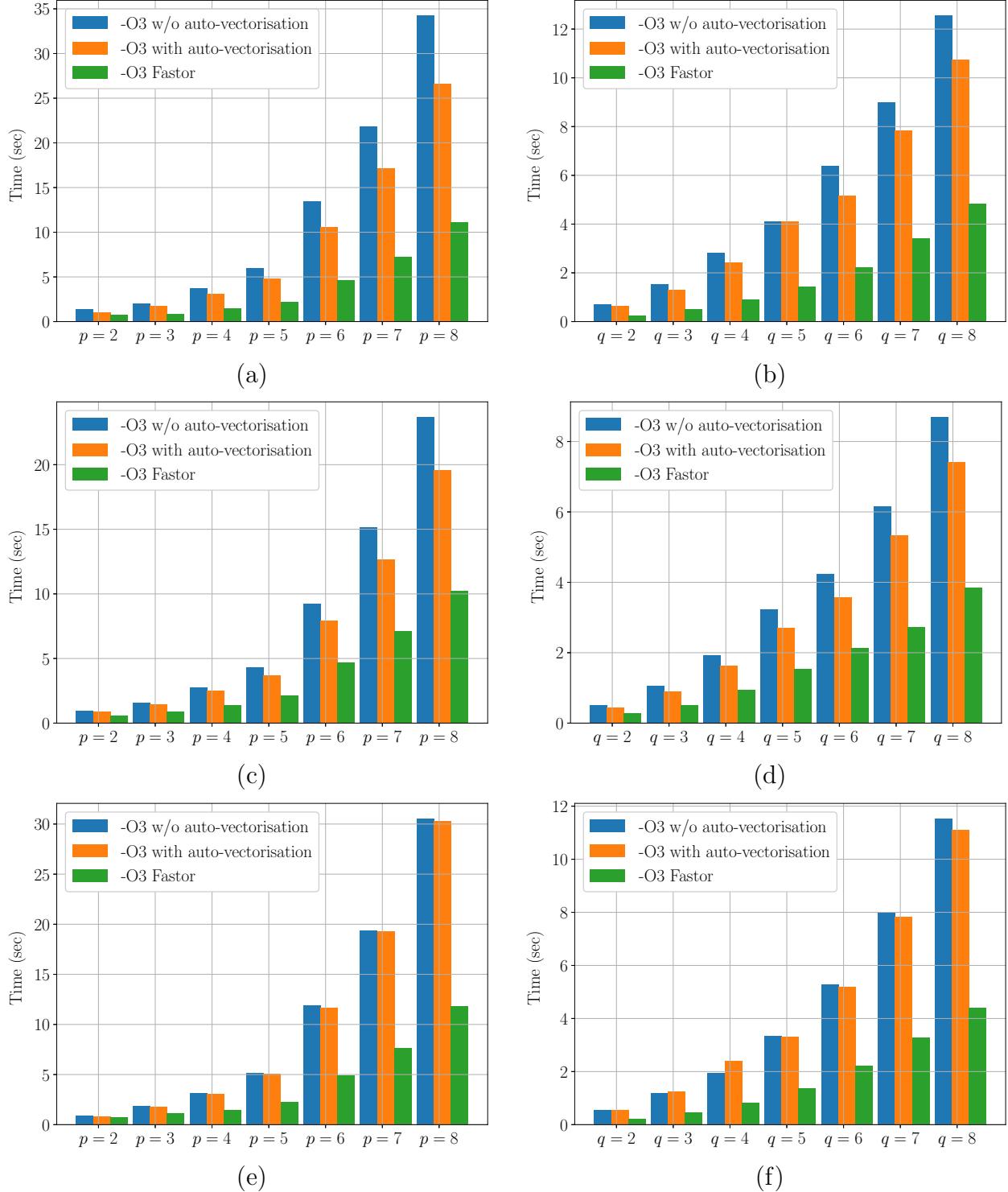


Figure 23: Speed-up achieved by compile time algorithmic transformation and data parallelism for numerical integration of work-conjugates and Hessian of convex multi-variable electro-elastic model (42) with high order curved tetrahedral and hexahedral meshes with compilers a,b) GCC, c,d) LLVM's Clang and, e,f) Intel's ICC

endeavour is put in the upcoming examples to present application-oriented numerical modelling of dielectric elastomers. Unless otherwise specified, all the examples in this section are modelled

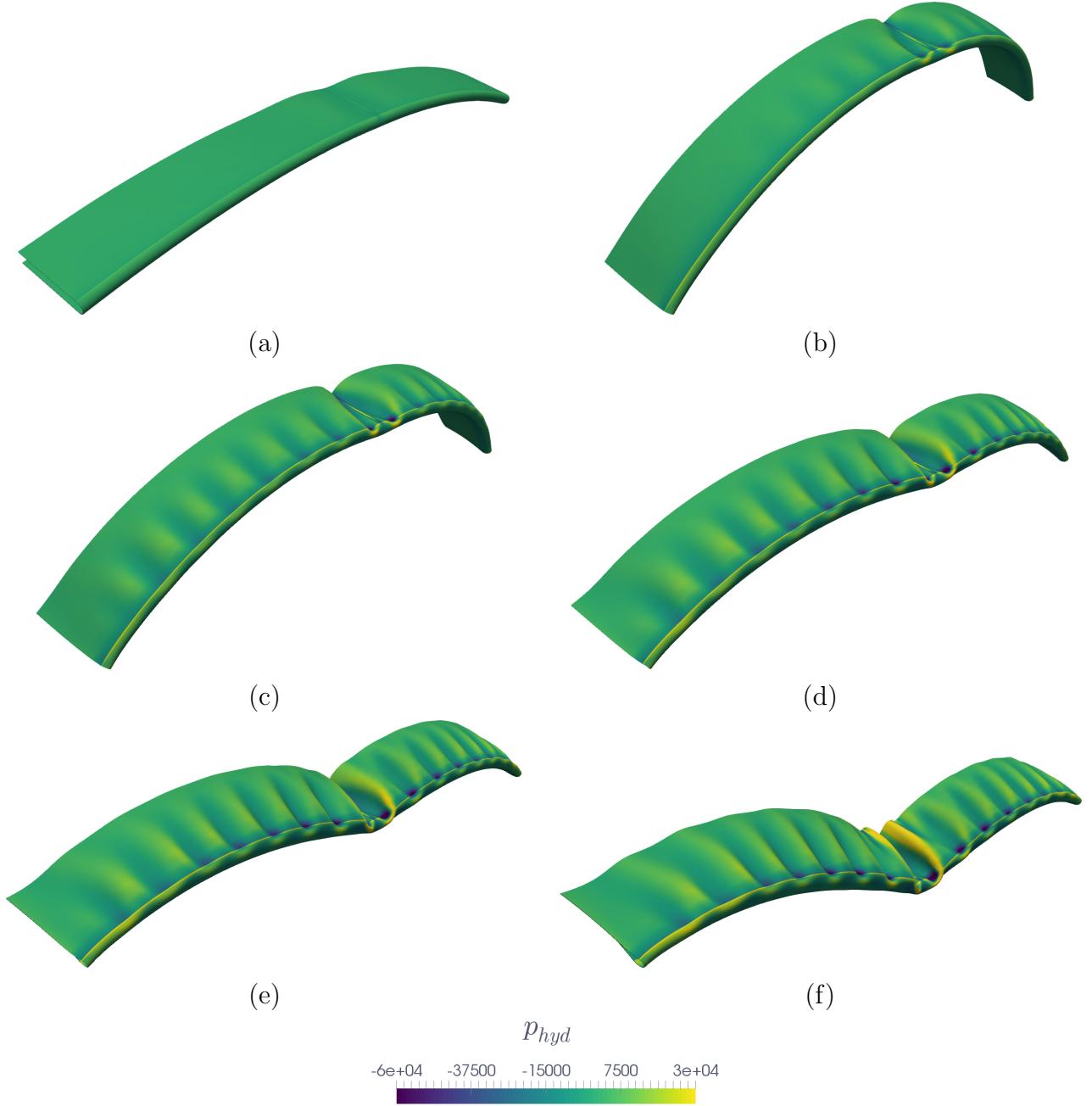
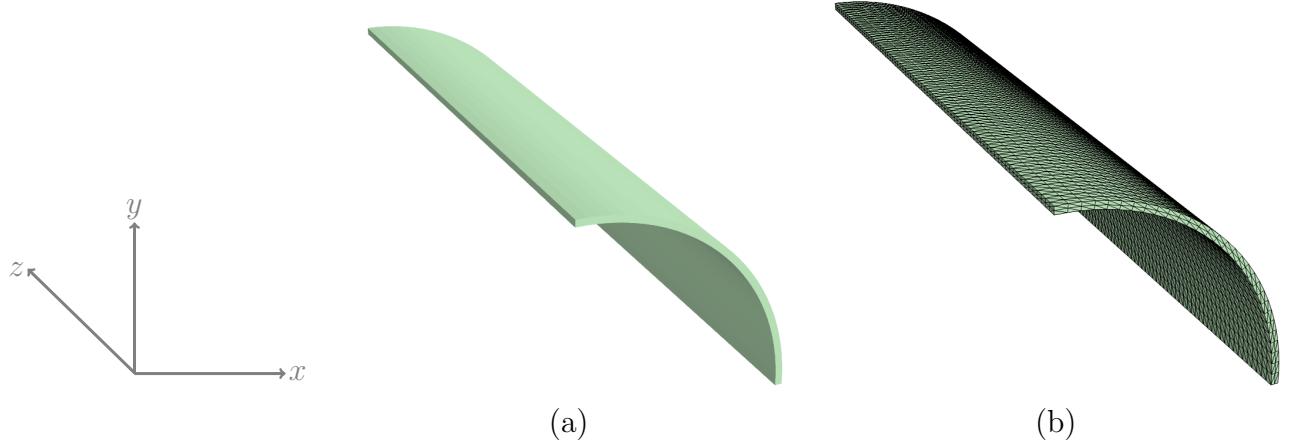


Figure 24: Evolution of voltage induced hydrostatic pressure p_{hyd} in electromechanical component with accumulated load factor Λ being a) $\Lambda = 0.416$ b) $\Lambda = 0.806$, c) $\Lambda = 0.889$, d) $\Lambda = 0.944$, e) $\Lambda = 0.972$ and, f) $\Lambda = 1.0$. Note that in the figures shown, the electromechanical component is mirrored along the z -axis

using the ideal dielectric elastomer material presented in (41) with material parameters listed in [Table 5](#). For majority of these examples, the meshes are extremely fine in order to capture the formation of wrinkles (despite polynomial enrichment) and, as a consequence, the distortion quality of curvilinear meshes are almost unity.

4.6.1. A massively deformable dielectric cylinder

The first example considered is that of a dielectric shell-like cylindrical structure shown in [Figure 25\(a\)](#) with 10m outer radius, 0.333m thickness and 25m length, centred at $[0, 0, 0]^T$. The model is made up of 27000 ($p = 3$) curvilinear tetrahedral elements with 179122 points in the computational mesh corresponding to a total of 716488 degrees of freedom (shown in [Figure 25\(b\)](#)). An ideal dielectric elastomer is used as material model and the material properties are shown in



[Figure 25](#): Geometry and a curved tetrahedral mesh of the dielectric shell-like cylindrical structure with 10m radius 0.333m thickness and 25m length, centred at $[0, 0, 0]^T$

[Table 18](#). The following set of boundary conditions are applied.

μ_1 (Pa)	μ_2 (Pa)	λ (Pa)	ε_1^{-1} (V^2/N)	ε_2 (N/V^2)
$1e5$	0	$4e5$	0	$4\varepsilon_0$

[Table 18](#): Material parameters with the vacuum permittivity $\varepsilon_0 = 8.85418781 \times 10^{-12}$ (N/V^2)

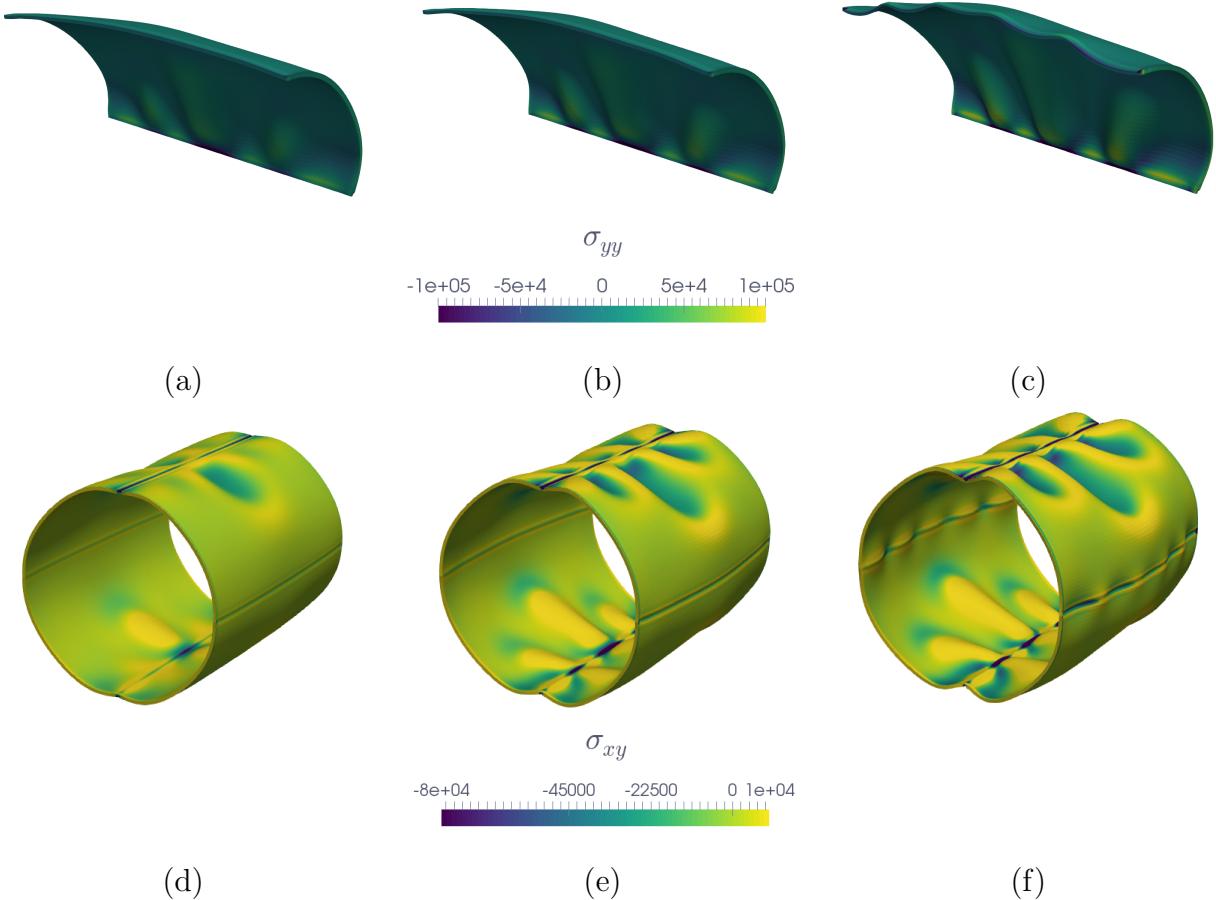
$$\begin{aligned}
u_y &= 0 \text{ [m]} & \text{at } \mathbf{X} &= [0, y, z]^T; \\
u_x &= 0 \text{ [m]} & \text{at } \mathbf{X} &= [x, 0, z]^T; \\
\mathbf{u} &= \mathbf{0} \text{ [m]} & \text{at } \mathbf{X} &= [10, y, z]^T; \\
\varphi &= 0 \text{ [V]} & \text{at } & R_{in}; \\
\omega_0 &= 1.4 \times 10^{-4} \Lambda \text{ [Q/m}^2\text{]} & \text{at } & R_{out},
\end{aligned}$$

where $R_{in} = 9.6667\text{m}$ is the inner radius, $R_{out} = 10\text{m}$ is the outer radius and Λ

$$\Lambda \in [0, 1] \subset \mathbb{R}$$

is simply a load factor controlling the application of surface charge such that, $\Lambda = 0$. represents start of the load increment and $\Lambda = 1$. represents end of the load increment (application of full load). The usage of load factor as a pseudo-time step to quantify the application of load is a standard practice in nonlinear quasi-static analyses [\[66, 67, 68\]](#). The above non-uniform set of boundary conditions causes the cylinder to deform asymmetrically. Furthermore, the disparity

in electric charge between the electrodes across the thickness causes the thickness to shrink. As shown in [Figure 26](#), when the voltage is small, the elastomer deforms slightly, and the charge increases with the electric voltage approximately linearly. As the voltage increases, the elastomer's surface area expands significantly, and a small increase in the voltage adds a large amount of displacement on the cylinder. Consequently, after the voltage reaches a maximum value the voltage needed to maintain the charge drops [\[24\]](#). This corresponds to the pull-in instability [\[25\]](#) causing regions of the cylinder to snap-back to fully new configuration possessing a different thickness. As more charge is applied at this state, the difference between the thick and thin regions causes the surface of the cylinder to fold and form wrinkles. This can be clearly observed in [Figure 26\(c,f\)](#). Notice that at the extreme loading state the elements closer to symmetry surfaces become considerably thinner and a complete zone of compression is formed at this region. Since



[Figure 26](#): Evolution of charge induced stresses in dielectric cylindrical structure with accumulated load factor Λ being a) σ_{yy} at $\Lambda = 0.678$ b) σ_{yy} at $\Lambda = 0.785$, c) σ_{yy} at $\Lambda = 1.0$ d) σ_{xy} at $\Lambda = 0.678$ e) σ_{xy} at $\Lambda = 0.785$, f) σ_{xy} at $\Lambda = 1.0$

the formation of wrinkles causes the cylinder to maintain different circumferential expansion at different regions along the length, one measure to quantify actuation property is to report the electric voltage as a function of circumferential expansion. One such graph is shown in [Figure 27](#) for different positions along the length of the cylinder. The figure confirms the experimental findings in [\[24\]](#) that the formation of wrinkles is beyond the point of pull-in instability, when the surface is a mixture of thinner and thicker regions.

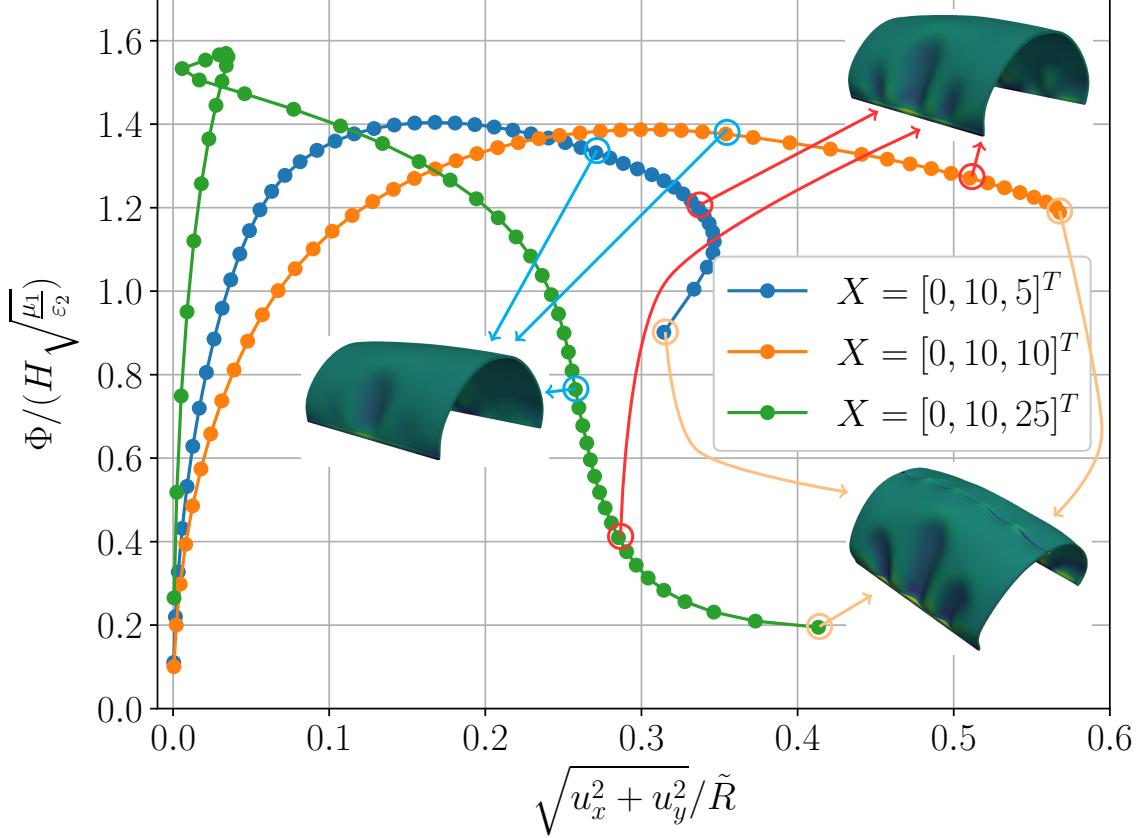


Figure 27: Non-dimensional quantity $\Phi/(H\sqrt{\frac{\mu_1}{\varepsilon_2}})$ quantifying electric voltage as a function of circumferential expansion, showing points of snap-through, pull-in instability and formation of coarse wrinkles in dielectric elastomer. H stands for thickness and $\tilde{R} = 9.8333$ is the averaged radius accounting for thickness stretch

Next, two $q = 3$ hexahedral meshes are chosen for the cylinder and two sets of boundary conditions are applied on each one respectively. The meshes correspond to 1.5M and 3.2M degrees of freedom, approximately. The first set of boundary conditions is similar to the previous analysis, however instead of surface charge now an electric potential of $5.2 \times 10^7 V$ is applied at the inner radius. The second set of boundary conditions corresponds to

$$\begin{aligned}
 \mathbf{u} &= \mathbf{0} \text{ [m]} & \text{at } \mathbf{X} &= [0, y, z]^T; \\
 \mathbf{u} &= \mathbf{0} \text{ [m]} & \text{at } \mathbf{X} &= [x, 0, z]^T; \\
 \varphi &= 0 \text{ [V]} & \text{at } R_{in}; \\
 \varphi &= 5.2 \times 10^7 \Lambda \text{ [V]} & \text{at } R_{out},
 \end{aligned}$$

where Λ is now the load factor quantifying the application of electric potential for the nonlinear quasi-static analysis. Due to symmetry, only 1/8th of geometries are analysed. Figure 28 shows the evolution of the F_{xz} component of the deformation gradient tensor with voltage history. The first observation is that since the boundary condition is now through applied voltage and not electric charge as in the previous analysis, the voltage will always constantly increase and

hence there is no negative or downward slope in voltage vs strain curves, up until the point of electric breakdown. The second observation from [Figure 28](#) is that, unlike in the case of applied charge where the surfaces near symmetry boundary conditions become thinner, here these surfaces undergo twisting to accommodate for the constant applied electric voltage. The formation of coarse wrinkles at $[0, x, y]^T$ is not present in this case.

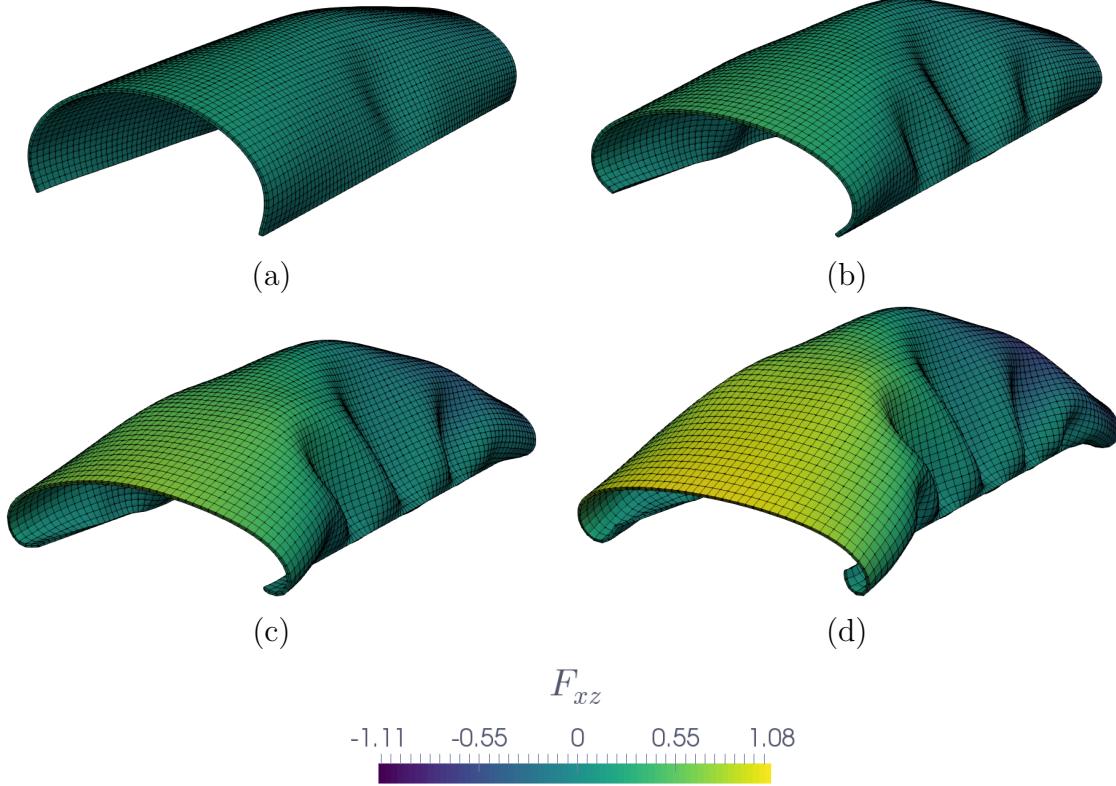


Figure 28: Evolution of voltage induced strain F_{xz} in dielectric cylindrical structure with accumulated load factor Λ being a) $\Lambda = 0.538$ b) $\Lambda = 0.769$, c) $\Lambda = 0.923$ and, d) $\Lambda = 1.0$

Finally, unlike the two previous analyses, the third example with the second set of boundary conditions for the hexahedral mesh mentioned above is completely uniform in terms of loading. As can be seen in [Figure 29](#) this set of boundary conditions imposes a homogeneous circumferential expansion. However, once the electric voltage is high enough, coarse wave-like wrinkles starting from the central point along the length of the cylinder are formed propagating towards the two ends of the cylinder.

The above three examples of cylindrical shell-like DEs pinpoint the different actuation properties that can be activated different sets of boundary conditions. In addition, they verify the capability of the current framework to cope with extremely fine meshes and high polynomial enrichment to capture massive deformations and wrinkling in dielectric elastomers.

4.6.2. Capsule

The next example considered is that of a dielectric capsule, shown in [Figure 30\(a,b\)](#) centered at $[0, 0, 0]^T$ with an in-plane radius of 10m, thickness of 0.5m and a bulging factor of 2m in the out of plane direction. The mesh considered (shown in [Figure 30\(c\)](#)) has 10112 elements and for an interpolation degree $q = 3$ there are 365584 points in the computational mesh which corresponds

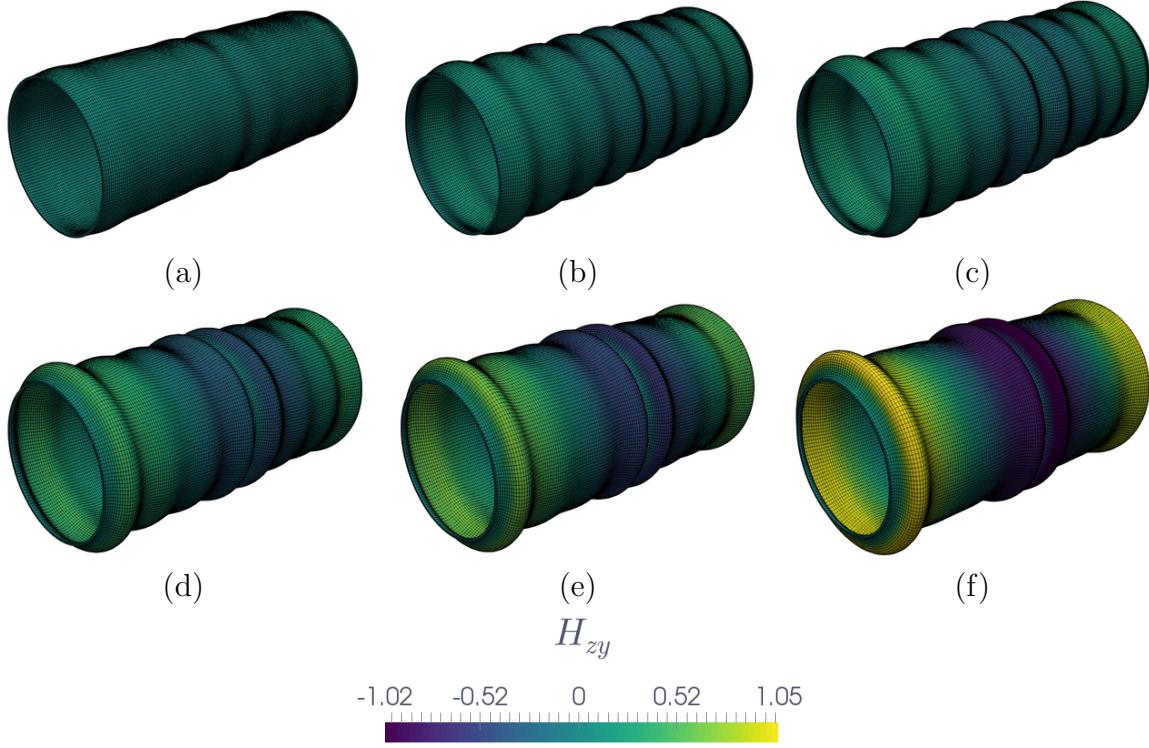


Figure 29: Evolution of voltage induced strain (co-factor of the deformation gradient) H_{zy} in dielectric cylindrical structure with accumulated load factor Λ being a) $\Lambda = 0.490$ b) $\Lambda = 0.686$, c) $\Lambda = 0.784$, d) $\Lambda = 0.882$, e) $\Lambda = 0.941$ and, f) $\Lambda = 1.0$

to 1462336 degrees of freedom. Due to symmetry, only 1/4th of the geometry is analysed. A set of symmetric mechanical Dirichlet boundary conditions are applied on the mesh with an applied electric voltage across the thickness, as described in the following.

$$\begin{aligned}
 \mathbf{u} = \mathbf{0} \text{ [m]} & \quad \text{at } \mathbf{X} = [0, y, z]^T; \\
 u_z = 0 \text{ [m]} & \quad \text{at } \mathbf{X} = [x, 0, z]^T; \\
 \varphi = 0 \text{ [V]} & \quad \text{at } R_{in}; \\
 \varphi = 2.5 \times 10^7 \Lambda \text{ [V]} & \quad \text{at } R_{out}.
 \end{aligned}$$

This problem is in particular interesting in shape and pattern forming through actuation. [Figure 31](#) shows the hydrostatic pressure at different loading stages. As can be observed the capsule undergoes massive deformation to adopt a squared shape from an initial circular configuration, as the electric voltage is applied. The capsule deforms significantly when the electric voltage is high to adopt the new shape. After the point of snap-through as more voltage is applied, a compression zone is formed at the centre. As the electric voltage is increased, this compression zone rapidly moves from the centre towards the edges of the deformed capsule signalling the propagation of compressed band through a highly inflated capsule. The occurrence of this phenomenon is indeed analogous to the ones reported in the previous section, in that, after the point of pull-in instability the thicker regions expand at the expense of the thinner region resulting in zones of compression and propagation of pressure as a wave through the capsule.

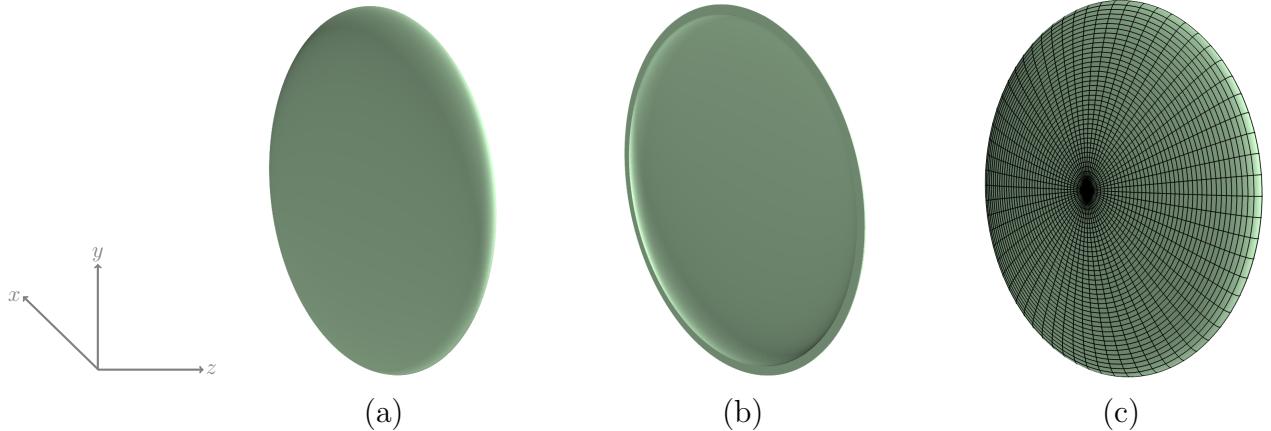


Figure 30: Dielectric capsule centred at $[0, 0, 0]^T$ with an in-plane radius of 10m, thickness of 0.5m and a bulging factor of 2m a,b) CAD geometry and c) $q = 3$ structured hexahedral mesh

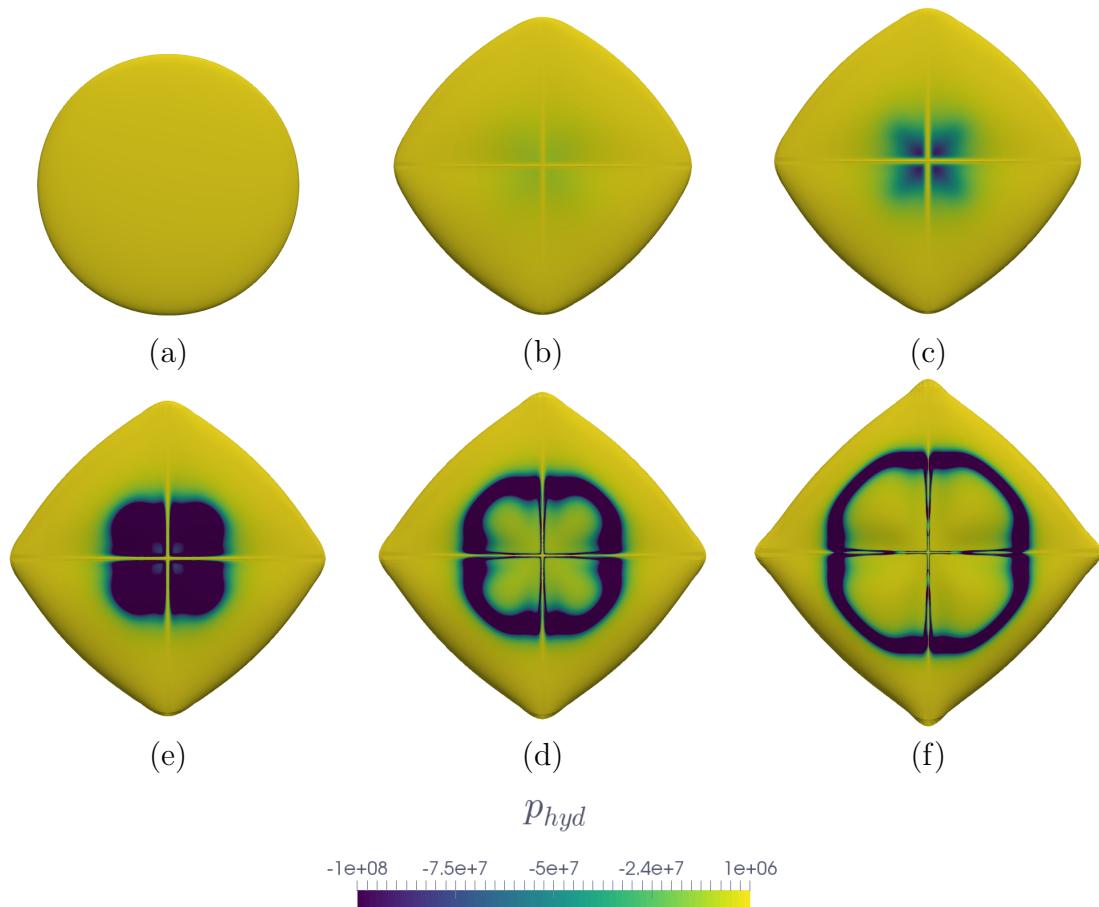
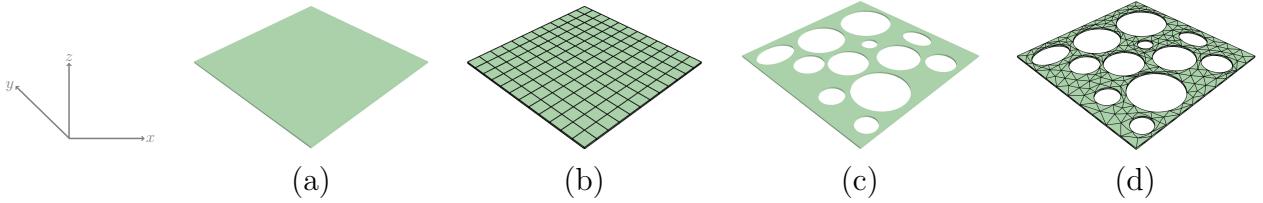


Figure 31: Evolution of voltage induced hydrostatic pressure in dielectric capsule with the accumulated load factor Λ being a) $\Lambda = 0.411$, b) $\Lambda = 0.767$, c) $\Lambda = 0.823$, d) $\Lambda = 0.882$, e) $\Lambda = 0.946$ and, f) $\Lambda = 1.0$

4.6.3. Capturing folds and coarse wrinkles in dielectric plates with and without inclusions

The objective of this study is to show the effect of inclusions and the capability of the current high order curvilinear framework in capturing folds and coarse wrinkles with very coarse meshes

through polynomial enrichments. To this effect, two dielectric plates are considered one without holes or inclusions and one with multiple circular inclusions distributed non-homogeneously through the plate, as shown in [Figure 32](#). The mesh considered for the plate without inclusion is an extremely coarse $12 \times 12 \times 1$ structured hexahedral mesh (with only one element across the thickness). For a polynomial degree of $q = 6$ there are 37303 points in the computational mesh corresponding to a total of 149212 degrees of freedom. For the plate with inclusions, the mesh considered is an unstructured tetrahedral with 2147 elements, as shown in [Figure 32\(d\)](#). For an interpolation degree of $p = 4$, there are 26943 points in the computational mesh corresponding to a total of 107772 degrees of freedom.



[Figure 32](#): Geometries and meshes for the dielectric plates of size $100 \times 100 \times 1 \text{ m}^3$, a,b) without inclusions and, c,d) with inclusions

The following set of boundary conditions are applied on both plates

$$\begin{aligned}
\mathbf{u} &= \mathbf{0} [\text{m}] & \text{at } \mathbf{X} &= [0, y, 0]^T; \\
\mathbf{u} &= \mathbf{0} [\text{m}] & \text{at } \mathbf{X} &= [100, y, 0]^T; \\
\mathbf{u} &= \mathbf{0} [\text{m}] & \text{at } \mathbf{X} &= [x, 0, 0]^T; \\
\mathbf{u} &= \mathbf{0} [\text{m}] & \text{at } \mathbf{X} &= [x, 100, 0]^T; \\
\varphi &= 0 [\text{V}] & \text{at } \mathbf{X} &= [x, y, 0]^T; \\
\varphi &= 4.98 \times 10^7 \Lambda [\text{V}] & \text{at } \mathbf{X} &= [x, y, 1]^T.
\end{aligned}$$

[Figure 33](#) shows the evolution of deformation in the plate (without inclusions) at multiple loading stages and [Figure 34](#) shows the final configuration of the plate. The formation of folds and coarse wrinkles can be clearly seen from the figures. Notice how the high order curvilinear elements are able to capture this intrinsic property of dielectric elastomers with a very coarse discretisation.

Similarly, [Figure 35](#) shows the evolution of strain component F_{xz} in the plate with inclusions at multiple loading stages and [Figure 36](#) shows the final configuration of the plate. The formation of folds can be clearly seen from the figures. However, unlike in the case of plate with no inclusions, the deformation pattern is completely non-uniform and the plate does not deform as much. Once again, notice how the high order curvilinear elements are able to capture foldings in dielectric elastomers with an increased level of detail despite a coarse discretisation.

While folding and the formation of single layer coarse wrinkles can be captured accurately using the high order curvilinear elements, as presented in the next section, multi-layer wrinkling and the propagation and nucleation of extremely fine wrinkles would still require a refined computational mesh.

4.6.4. Dielectric thin film undergoing massive wrinkling

The objective of this final example is to study the voltage induced instability in a thin dielectric elastomeric film undergoing massive formation, propagation and nucleation of wrinkles.

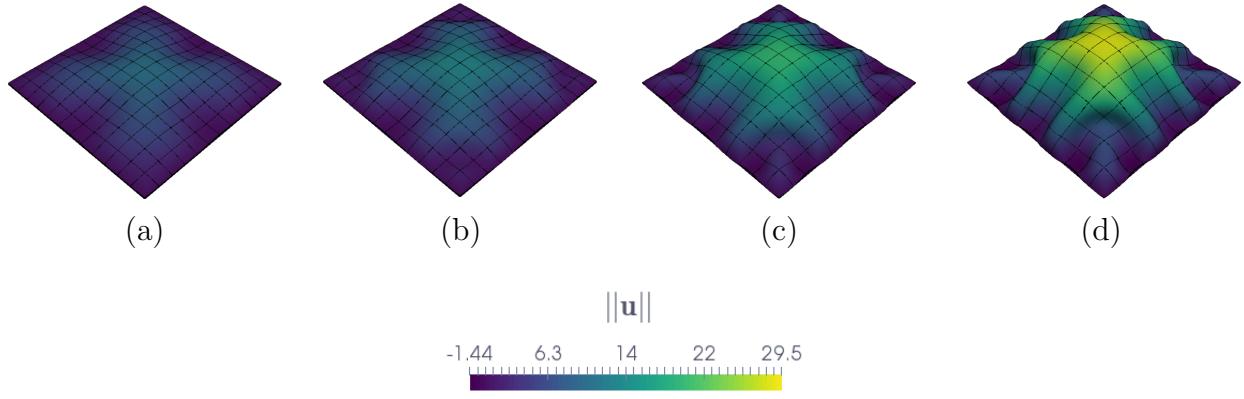


Figure 33: Evolution of norm of displacements in dielectric plate with the accumulated load factor Λ being a) $\Lambda = 0.482$, b) $\Lambda = 0.602$, c) $\Lambda = 0.843$ and, d) $\Lambda = 1.0$

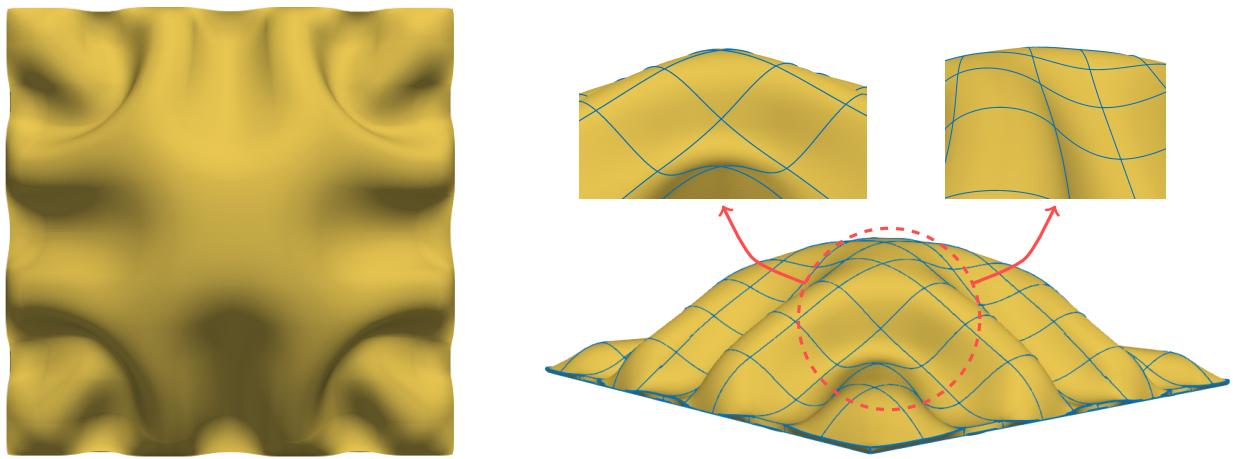


Figure 34: Formation of folds in dielectric plate (at the final deformed configuration) captured on a $12 \times 12 \times 1$ hexahedral mesh with $q = 6$ polynomial interpolation

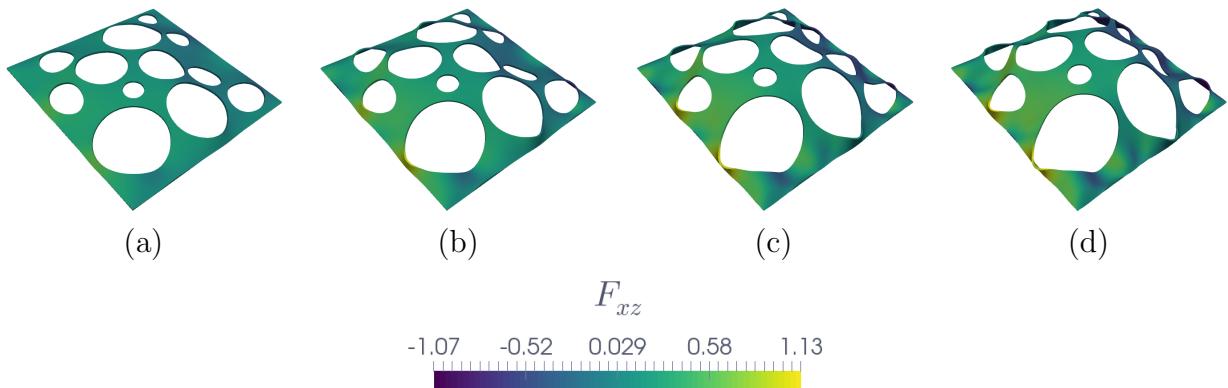


Figure 35: Evolution of strains F_{xz} in dielectric plate with inclusions with the accumulated load factor Λ being a) $\Lambda = 0.482$, b) $\Lambda = 0.602$, c) $\Lambda = 0.843$ and, d) $\Lambda = 1.0$

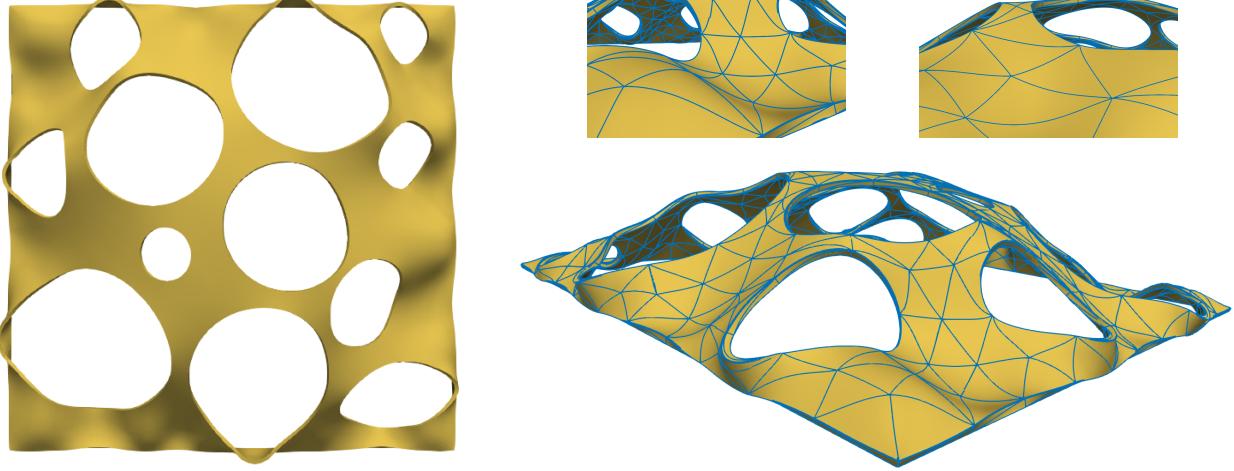


Figure 36: Formation of folds in dielectric plate with inclusions (at the final deformed configuration) captured on an unstructured tetrahedral mesh with $p = 4$ polynomial interpolation

Instabilities in the form of wrinkling have also been studied in [26, 27, 28]. For the analysis, a hexahedral mesh with two polynomial enrichments is considered namely $q = 2$ and $q = 6$ with 268140 and 1.4M degrees of freedom, respectively. The latter allows for a highly detailed propagation of wrinkles through the film in terms of geometry and solution accuracy. The CAD geometry and mesh of the film are shown in Figure 37 with 100m radius and 1m thickness centred at $[0, 0, 0]^T$.

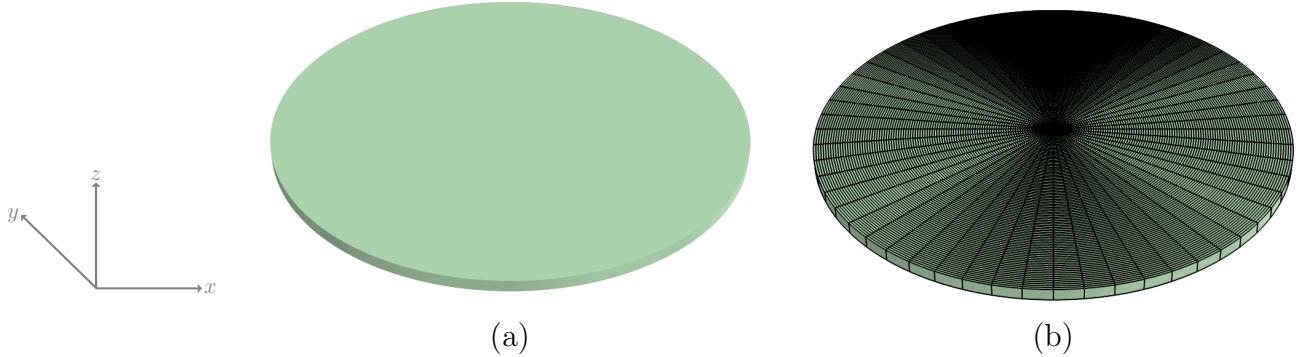


Figure 37: Dielectric thin film with 100m radius and 1m thickness centred at $[0, 0, 0]^T$, a) CAD geometry and, b) $q = 2$ mesh

The following set of boundary conditions is applied on the film

$$\begin{array}{ll} \mathbf{u} = \mathbf{0} \text{ [m]} & \text{at } \mathbf{X} = [x, y, 0]^T \text{ s.t. } x^2 + y^2 = 100^2; \\ \varphi = 0 \text{ [V]} & \text{at } \mathbf{X} = [x, y, 0]^T; \\ \varphi = 5 \times 10^7 \Lambda \text{ [V]} & \text{at } \mathbf{X} = [x, y, 1]^T. \end{array}$$

This above set of boundary conditions essentially implies fixing the mechanical variables at the outer boundary of the base of the film and applying an electric voltage across the thickness of the film. This forces the film to bend perpendicularly and expand in area. The electric voltage

is applied adaptively on the film through 1000 load increments. Figure 38 shows the evolution of stress components σ_{xz} and σ_{yz} in dielectric film for $q = 2$ mesh. As can be observed under low electric voltage, the region near the boundary of the circular film starts bulging up. This allows the film to expand in area and shrink in thickness in this region. As the voltage is increased the thick regions surrounding the thin ones start to expand and bulge facilitating the formation and further propagation of wrinkles. Under further increase in voltage all the regions start to possess similar thickness and the film straightens again allowing for nucleation of many wrinkles. The process of wrinkling starts once again, when the voltage is further increased. The process keeps repeating itself as voltage keeps increasing up until the point of complete electric breakdown. Note that the formation of deep channels near the boundary require polynomial enhancement despite a very fine mesh to capture the massive voltage induced bending caused by wrinkles.

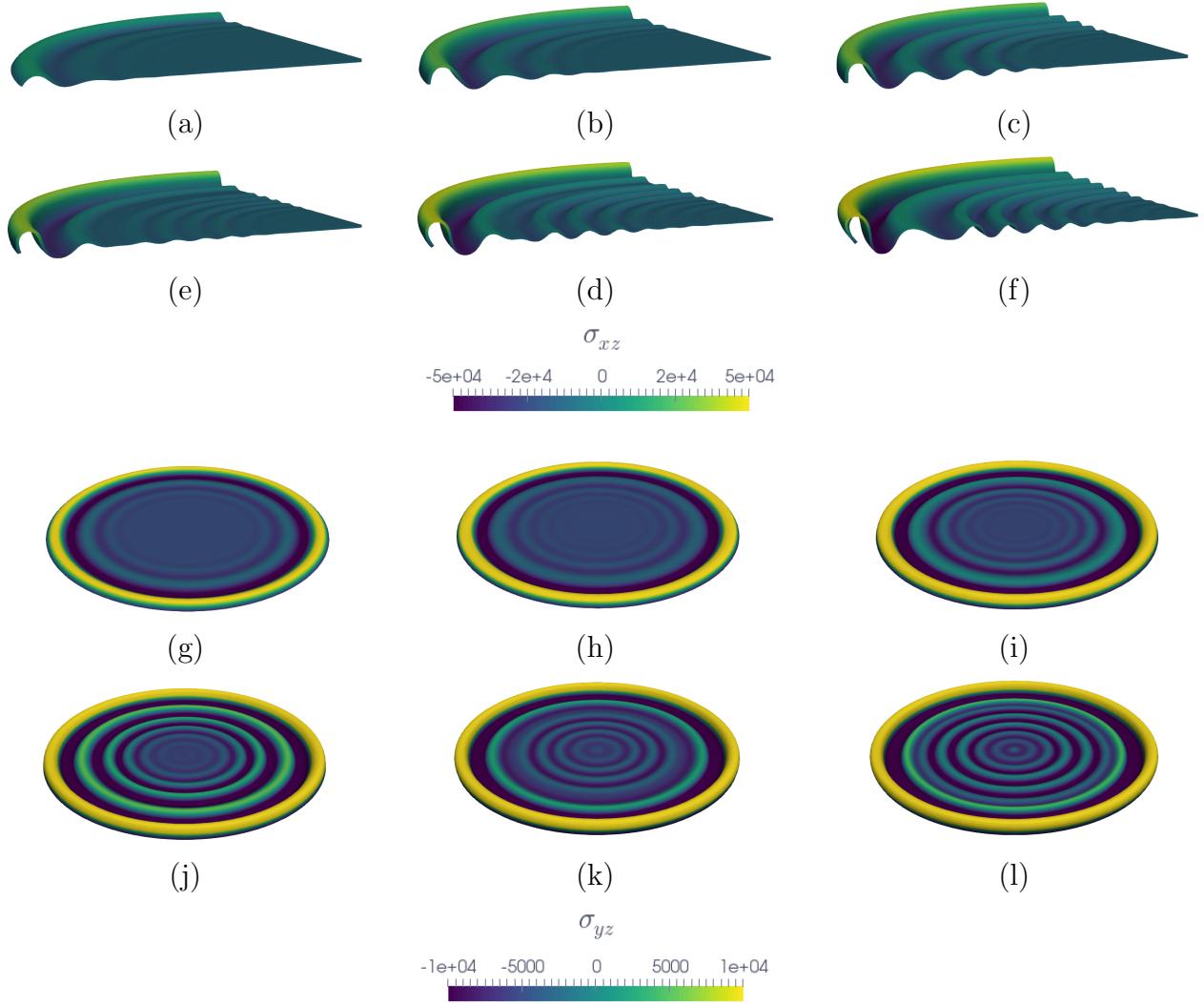


Figure 38: Evolution of voltage induced stress components σ_{xz} (a,b,c,d,e,f) and σ_{yz} (g,h,i,j,k,l) in dielectric film with the accumulated load factor Λ being a,g) $\Lambda = 0.580$, b,h) $\Lambda = 0.725$, c,i) $\Lambda = 0.798$, d,j) $\Lambda = 0.870$, e,k) $\Lambda = 0.943$ and, f,l) $\Lambda = 1.0$

The same problem is then analysed using $q = 6$ elements, however now after the formation a certain number of wrinkles in the film, the load is released. Figure 39 shows the deformation

history of the film together with the evolution of hydrostatic pressure. As the load is released, the deep channels around the boundary start to move inwards, leaving a thick bended layer behind. As the whole load cycle is completed the film starts flattening and forming a plate with multiple tiny wrinkles on the surface.

Finally, the aspect ratio of the film is decreased from 1/100 to 1/20 i.e. the radius is decreased to $20m$ and the same problem with release load cycle is analysed using the $q = 6$ mesh. The deformation pattern is the same as in the previous analysis, however, since now there is little room for the formation of multiple thick and thin regions, wrinkling is limited. As shown in [Figure 40](#) a deep channel is formed near the boundary that starts propagating inwards till it reaches the centre. The film starts occupying a balloon shape at this stage. As the voltage is increased, the centre of the film starts bending in the opposite direction, allowing the formation of another layer of wrinkles to propagate outwards. Once the wrinkle is propagated through the film by reaching the boundaries, another layer of wrinkles starting from the centre starts propagating outwards. The process is repeated and in the process, multiple wrinkles start nucleating before hitting the boundary. Under the completion of the release cycle, the film starts flattening leaving multiple extremely fine wrinkles behind on the surface.

Overall, the problem of analysing the formation of wrinkles in dielectric elastomers is a challenging one, in that it requires a very fine mesh. With coarse meshes it is possible to capture folds using high order curvilinear elements, however multi-layer wrinkling might not even occur using coarse meshes. Even high polynomial enrichments do not allow for a wave like deformation of a single element in a wrinkled state, and hence h -refinement is absolutely necessary. Where polynomial enrichment really helps is in massive bending of single elements during the formation of deep channels. The examples discussed afore, pinpoint the robustness and the high performant capability of the framework in capturing massive deformations and wrinkling in dielectric elastomers with a remarkable level of detail.

5. Concluding remarks

A high order finite element implementation of the convex multi-variable electro-elasticity for large deformations large electric fields simulations and its particularisation to the case of small strains through a staggered scheme is presented. Accurate geometrical representation through a high performance curvilinear finite element framework based on a posteriori mesh deformation technique is developed to accurately discretise the underlying displacement-potential variational formulation. The performance of the method under near incompressibility and bending actuation scenarios is analysed with extremely thin and highly stretched components and compared to the performance of mixed variational principles. Although convex multi-variable constitutive models are elliptic and hence, materially stable for the entire range of deformations and electric fields, other forms of physical instabilities are not precluded in these models. In particular, physical instabilities present in dielectric elastomers such as pull-in instability, snap-through and the formation, propagation and nucleation of wrinkles and folds are numerically studied with a detailed precision in this paper, verifying experimental findings. In this context, the combination of h and p refinement proves to be essential to capture the inherent instabilities in dielectric elastomers. While the formation of folds and coarse wrinkles can be accurately captured by high order curvilinear elements using extremely coarse meshes, nucleation of wrinkles and multi-layer wrinkling require mesh refinement in addition to polynomial enrichment. For the case of small strains, the staggered scheme is shown to capture the electrostrictive behaviour of EAPs fairly well with

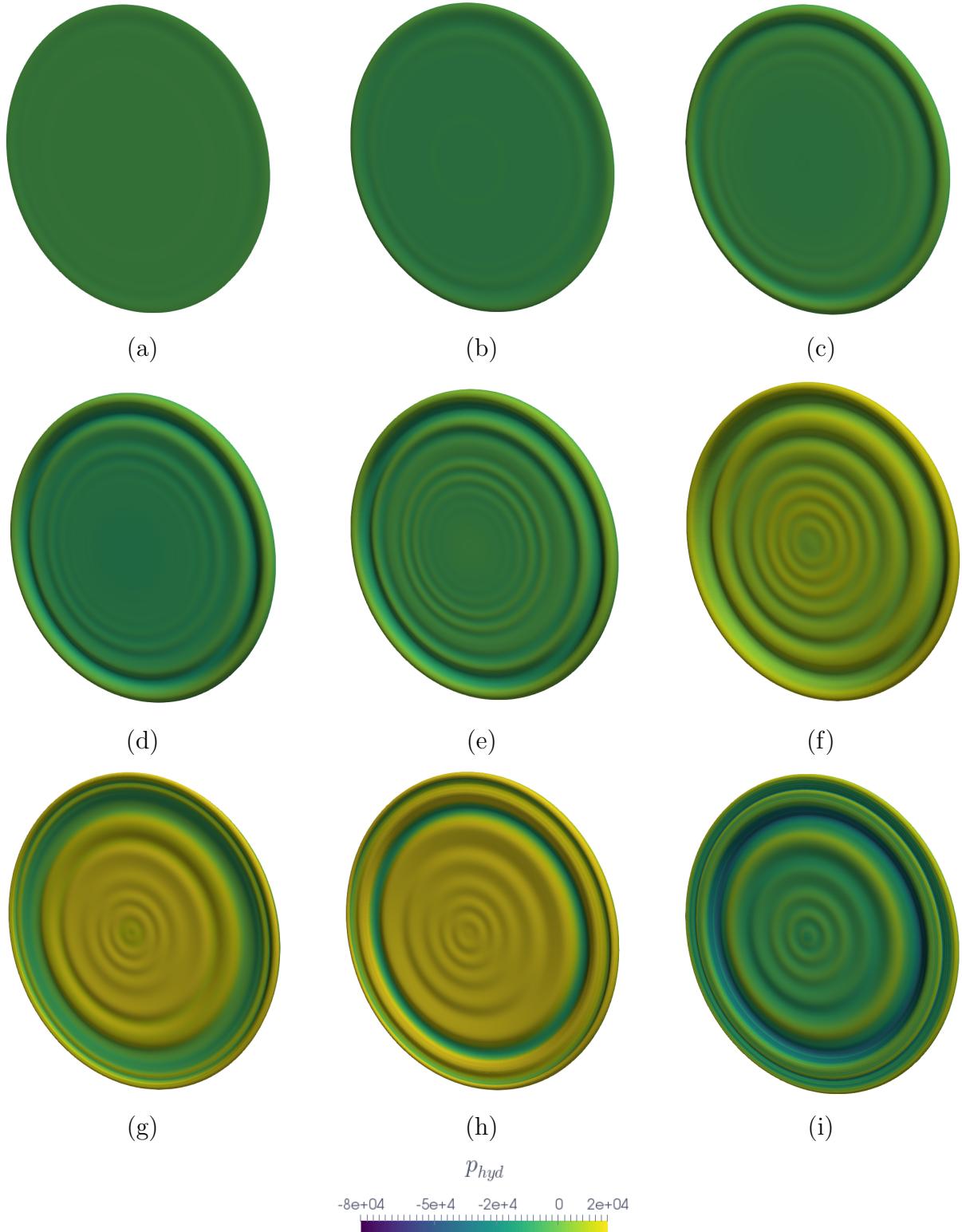


Figure 39: Evolution of voltage induced hydrostatic pressure in dielectric film with the accumulated load factor Λ being a) $\Lambda = 0.221$, b) $\Lambda = 0.332$, c) $\Lambda = 0.443$, d) $\Lambda = 0.554$, e) $\Lambda = 0.665$, f) $\Lambda = 0.887$, g) $\Lambda = 0.943$, h) $\Lambda = 0.971$, and, i) $\Lambda = 1.0$

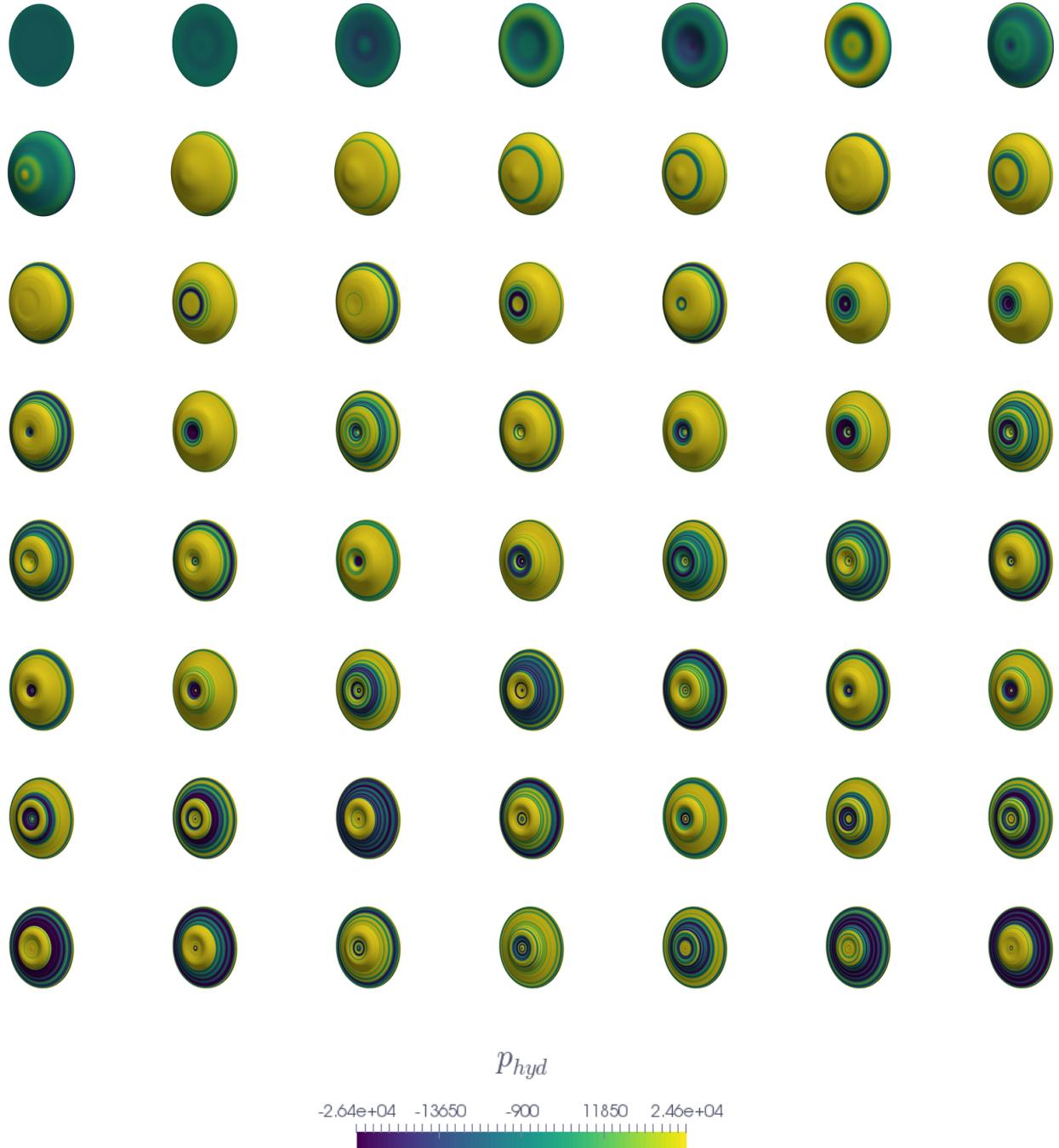


Figure 40: Evolution of voltage induced hydrostatic pressure in dielectric film with the accumulated load factor Λ being (left to right - top to bottom) $\Lambda = 0.0, \Lambda = 0.124, \Lambda = 0.249, \Lambda = 0.374, \Lambda = 0.499, \Lambda = 0.624, \Lambda = 0.749, \Lambda = 0.874, \Lambda = 0.886, \Lambda = 0.899, \Lambda = 0.911, \Lambda = 0.918, \Lambda = 0.921, \Lambda = 0.924, \Lambda = 0.928, \Lambda = 0.93, \Lambda = 0.934, \Lambda = 0.936, \Lambda = 0.94, \Lambda = 0.942, \Lambda = 0.949, \Lambda = 0.952, \Lambda = 0.955, \Lambda = 0.958, \Lambda = 0.96, \Lambda = 0.962, \Lambda = 0.964, \Lambda = 0.965, \Lambda = 0.966, \Lambda = 0.968, \Lambda = 0.969, \Lambda = 0.97, \Lambda = 0.971, \Lambda = 0.972, \Lambda = 0.974, \Lambda = 0.975, \Lambda = 0.976, \Lambda = 0.978, \Lambda = 0.979, \Lambda = 0.98, \Lambda = 0.981, \Lambda = 0.982, \Lambda = 0.984, \Lambda = 0.985, \Lambda = 0.986, \Lambda = 0.988, \Lambda = 0.989, \Lambda = 0.99, \Lambda = 0.991, \Lambda = 0.992, \Lambda = 0.994, \Lambda = 0.995, \Lambda = 0.996, \Lambda = 0.998, \Lambda = 0.999$ and, $\Lambda = 1.0$

a threshold point in applied voltage beyond which the fully coupled nonlinear solver becomes

computational more favourable. Importantly, the staggered approach can also be applied to large displacements small strains problems. The staggered approach is competitive due to the fact that for its range of applicability, only the scalar field Gauss's law needs to be solved for iteratively.

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