Mixed finite elements for port-Hamiltonian models of von Kármán beams

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Abstract: A port-Hamiltonian formulation of von Kármán beams is presented. The variables selection lead to a non linear interconnection operator, while the constitutive laws are linear. The model can be readily discretized by exploiting a coenergy formulation and a mixed finite element method. The mixed formulation does not demand the H^2 regularity requirement typical of standard Galerkin discretization of thin structures. A numerical test is performed to assess the convergence rate of the solution.

Keywords: Port-Hamiltonian systems, von Kármán beams, Mixed Finite Elements

1. INTRODUCTION

Linear elastic structures have been largely investigated into the port-Hamiltonian (pH) framework as well as the heat equation (consult for instance Macchelli and Melchiorri (2004) for the Timoshenko beam, Aoues et al. (2017) for the Euler-Bernoulli beam, Brugnoli et al. (2019b,a) for thick and thin plates). Recently, more complicated models arising from fluid dynamics have also been considered (Cardoso-Ribeiro et al., 2019, 2020b; Rashad et al., 2021a,b; Califano et al., 2021; Altmann and Schulze, 2017).

The Hamiltonian foundation of non-linear elasticity dates back to the late 80s (Simo et al., 1988). A pH formulation of flexible beams undergoing large deformations can be found in Macchelli et al. (2007, 2009). One fundamental non linear theory is the von Kármán one, that applies for moderately large deformation of beams, plates and shells. This theory has puzzled mathematicians and physicists as the derivation of the model rely on some not well justified assumptions. Mathematical conditions under which this theory is actually applicable and will give reasonable results when solved are discussed in Ciarlet (1980, 1990). In particular, for this model the stresses-strains constitutive relation has to be linear (isotropic) and the Kirchhoff-Love displacement field assumption (i.e. the normals to the middle surface remain normal to the deformed middle surface) hold. Existence and uniqueness of solutions for the full dynamical problem were established in Lagnese and Leugering (1991) for one-dimensional beams and in Puel and Tucsnak (1996) for plates. Since the full dynamical von Kárman problem is conservative (see e.g. Bilbao et al. (2015)), a pH realization of this system exists.

The development of new models within the pH framework has been accompanied with an increased interest in numerical discretization methods, capable of retaining the main features of the distributed system in its finite-dimensional counterpart (see Rashad et al. (2020) for a comprehensive review on this topic). Recently, it has become evident that there is a strict link between discretization of port-Hamiltonian systems and mixed finite elements (Cardoso-Ribeiro et al., 2020a). An example of this connection is given in Kirby and Kieu (2015), where a velocity-stress formulation for the wave dynamics is shown to be Hamiltonian and its mixed discretization preserves such a structure.

In this contribution, the von Kármán beam model is formulated as a pH system. The selection of energy variables will be such to make the Hamiltonian quadratic in these variables. As a consequence of this choice, the non linearities of the model are included in the interconnection operator, whereas the constitutive relations remain linear. The obtained model can be discretized using mixed finite elements. To this aim, a weak formulation that does not demand for H^2 regularity for the vertical displacement (as in classical Galerkin discretization of beams and plates, cf. Gustafsson et al. (2018)) is obtained. This means that the vertical displacement can be discretized using H^1 conforming elements (i.e. Continuous Galerkin elements), rather then using the more computationally demanding H^2 conforming elements, like the Hermite polynomials. A numerical test is carried out to evaluate the convergence rate of the discrete solution with respect to an analytical one.

The paper is organized as follows. In Section 2 the classical model for von Kármán beams is recalled. Then, a pH realization of the classical model is detailed in Sec. 3. The mixed finite element discretization strategy is discussed in Sec. 4. In Sec. 5 a numerical convergence test is performed to assess the rate of convergence of the solution.

2. VON KÁRMÁN BEAMS

The classical von Kármán beam model is presented in (Reddy, 2010, Chapter 4). Under the hypothesis of isotropic material, the extensional-bending stiffness is zero when the x-axis is taken along the geometric centroidal axis. This means that a pure traction (bending) deformation produces no bending (traction) stress. With this assumption, the problem, defined on an interval $\Omega = [0, L]$, takes the following form

$$\rho A\ddot{u} = \partial_x n_{xx},
\rho A\ddot{w} = -\partial_{xx}^2 m_{xx} + \partial_x (n_{xx}\partial_x w),$$
(1)

together with the stresses and strains expressions

$$n_{xx} = EA\varepsilon_{xx}, \qquad m_{xx} = EI\kappa_{xx},$$

 $\varepsilon_{xx} = \partial_x u + 1/2(\partial_x w)^2, \quad \kappa_{xx} = \partial_{xx}^2 w.$ (2)

Variable u is the horizontal displacement, w is the vertical displacement (cf. Fig. 1), n_{xx} is the axial stress resultant and m_{xx} is the bending stress resultant. The coefficients ρ, A, E, I are the beam's mass density, the cross section, the Young modulus and the second moment of area.

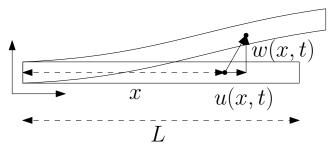


Fig. 1. Schematic representation of a von Kármán beam

The total energy of the model (Hamiltonian functional)

$$H = \frac{1}{2} \int_{\Omega} \left\{ \rho A(\dot{u}^2 + \dot{w}^2) + n_{xx} \varepsilon_{xx} + m_{xx} \kappa_{xx} \right\} d\Omega, \quad (3)$$

consists of the kinetic energy and both membrane and bending deformation energies. This model proves conservative, see Bilbao et al. (2015). Indeed, this implies that a port-Hamiltonian realization of the system exists. We shall demonstrate how to construct a port-Hamiltonian realization, equivalent to (1).

3. THE EQUIVALENT PORT-HAMILTONIAN REALIZATION

To find a suitable port-Hamiltonian system, we first select a set of new energy variables to make the Hamiltonian functional quadratic

$$\alpha_{u} = \rho A \dot{u},$$
 Horizontal momentum,
 $\alpha_{\varepsilon} = \varepsilon_{xx},$ Axial strain,
 $\alpha_{w} = \rho A \dot{w},$ Vertical momentum,
 $\alpha_{\kappa} = \kappa_{xx},$ Curvature. (4)

The energy is quadratic in these variables
$$H = \frac{1}{2} \int_{\Omega} \left\{ \frac{\alpha_u^2 + \alpha_w^2}{\rho A} + E A \alpha_{\varepsilon}^2 + E I \alpha_{\kappa}^2 \right\} \, \mathrm{d}\Omega. \tag{5}$$

By computing the variational derivative of the Hamiltonian, one obtains the so-called coenergy variables:

$$e_{u} := \delta_{\alpha_{u}} H = \dot{u}, \qquad e_{\varepsilon} := \delta_{\alpha_{\varepsilon}} H = n_{xx}, e_{w} := \delta_{\alpha_{w}} H = \dot{w}, \qquad e_{\kappa} := \delta_{\alpha_{\kappa}} H = m_{xx}.$$
 (6)

Before stating the final formulation, consider the unbounded operator operator $C(w)(\cdot): L^2(\Omega) \to L^2(\Omega)$, that acts as follows

$$C(w)(\cdot) = \partial_x(\cdot \partial_x w). \tag{7}$$

The domain of the operator is given by

$$D(\mathcal{C}) = \{ v \mid v \, \partial_x w \in H^1(\Omega) \}. \tag{8}$$

Proposition 1. The formal adjoint of the $C(w)(\cdot)$ is given

$$C(w)^*(\cdot) = -\partial_x(\cdot)\partial_x(w). \tag{9}$$

Proof 1. Consider a smooth scalar fields with compact support $\psi \in C_0^{\infty}(\Omega)$ and $\xi \in C_0^{\infty}(\Omega)$. The formal adjoint of $\mathcal{C}(w)(\cdot)$ satisfies the relation

$$\langle \psi, \mathcal{C}(w)(\xi) \rangle_{\Omega} = \langle \mathcal{C}(w)^*(\psi), \xi \rangle_{\Omega},$$
 (10)

where $\langle f, g \rangle_{\Omega} = \int_{\Omega} f g \ d\Omega$. The proof follows from the

$$\langle \psi, \mathcal{C}(w)(\xi) \rangle_{\Omega} = \langle \psi, \partial_x(\xi \, \partial_x w) \rangle_{\Omega},$$

$$= \langle -\partial_x \psi, \xi \partial_x w \rangle_{\Omega},$$

$$= \langle -\partial_x \psi \, \partial_x w, \xi \rangle_{\Omega}.$$
(11)

This means that

$$C(w)^*(\cdot) = -\partial_x(\cdot)\partial_x w, \tag{12}$$

leading to the final result.

The pH realization is then given by the following system

$$\frac{\partial}{\partial t} \begin{pmatrix} \alpha_u \\ \alpha_{\varepsilon} \\ \alpha_w \\ \alpha_{\kappa} \end{pmatrix} = \begin{bmatrix} 0 & \partial_x & 0 & 0 \\ \partial_x & 0 & (\partial_x w) \partial_x & 0 \\ 0 & \partial_x (\cdot \partial_x w) & 0 & -\partial_{xx}^2 \\ 0 & 0 & \partial_{xx}^2 & 0 \end{bmatrix} \begin{pmatrix} e_u \\ e_{\varepsilon} \\ e_w \\ e_{\kappa} \end{pmatrix}.$$
(13)

The second line of system (13) represents the time derivative of the axial strain ε_{xx} . To close the system, variable w has to be accessible. For this reason, its dynamics has to be included. The augmented system reads

$$\frac{\partial}{\partial t} \begin{pmatrix} \alpha_u \\ \alpha_\varepsilon \\ \alpha_w \\ \alpha_\kappa \\ w \end{pmatrix} = \underbrace{\begin{bmatrix} 0 & \partial_x & 0 & 0 & 0 & 0 \\ \partial_x & 0 & (\partial_x w) \partial_x & 0 & 0 \\ 0 & \partial_x (\cdot \partial_x w) & 0 & -\partial_{xx}^2 & -1 \\ 0 & 0 & \partial_{xx}^2 & 0 & 0 \\ 0 & 0 & 1 & 0 \end{bmatrix}}_{\mathcal{J}} \begin{pmatrix} e_u \\ e_\varepsilon \\ e_w \\ e_\kappa \\ \delta_w H \end{pmatrix}$$
(14)

The operator \mathcal{J} is formally skew-adjoint. If only the kinetic and deformation energies are considered, it holds $\delta_w H = 0$. In general this term allows accommodation of other potentials, such as gravitational. Suitable boundary variables are then obtained considering the power balance

$$\dot{H} = \langle e_u, e_{\varepsilon} \rangle_{\partial \Omega} + \langle e_w, e_{\varepsilon} \partial_x w - \partial_x e_{\kappa} \rangle_{\partial \Omega} + \langle \partial_x e_w, e_{\kappa} \rangle_{\partial \Omega}.$$
(15)

For the boundary terms, the following notation has been used

$$\langle f, g \rangle_{\partial\Omega} = fg|_0^L = f(L)g(L) - f(0)g(0).$$

The boundary conditions, summarized in Table 1, are consistent with the ones assumed in Puel and Tucsnak (1996) for deriving a global existence result for this model.

BCs	Traction	Bending	
Dirichlet BCs.	$e_u _0^L$	$e_w _0^L$	$\partial_x e_w _0^L$
Neumann BCs.	$ e_{\varepsilon} _{0}^{L}$	$e_{\varepsilon}\partial_x w - \partial_x e_{\kappa} _0^L$	$e_{\kappa} _{0}^{L}$

Table 1. Boundary conditions for von Karman beams.

4. MIXED FINITE ELEMENT DISCRETIZATION

To perform the numerical discretization, the constitutive relations are first incorporated in the dynamics. The link between the energy variables (4) and the coenergy variables (6) is given by the linear transformation

$$\begin{pmatrix}
\alpha_u \\
\alpha_\varepsilon \\
\alpha_w \\
\alpha_\kappa
\end{pmatrix} = \begin{bmatrix}
\rho A & 0 & 0 & 0 \\
0 & C_a & 0 & 0 \\
0 & 0 & \rho A & 0 \\
0 & 0 & 0 & C_b
\end{bmatrix} \begin{pmatrix}
e_u \\
e_\varepsilon \\
e_w \\
e_\kappa
\end{pmatrix},$$
(16)

where $C_a = (EA)^{-1}$ and $C_b = (EI)^{-1}$ are the axial and bending compliances respectively. The physical parameters are assumed to be constant for simplicity. A pure coenergy formulation can then be employed once (16) is plugged into (14)

$$\begin{pmatrix} \rho A \dot{e}_{u} \\ C_{a} \dot{e}_{\varepsilon} \\ \rho A \dot{e}_{w} \\ \dot{w} \end{pmatrix} = \begin{bmatrix} 0 & \partial_{x} & 0 & 0 & 0 \\ \partial_{x} & 0 & \partial_{x} w \, \partial_{x} & 0 & 0 \\ 0 & \partial_{x} (\cdot \partial_{x} w) & 0 & -\partial_{xx}^{2} & -1 \\ 0 & 0 & \partial_{xx}^{2} & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \end{bmatrix} \begin{pmatrix} e_{u} \\ e_{\varepsilon} \\ e_{w} \\ e_{\kappa} \\ \delta_{w} H \end{pmatrix}.$$

$$(17)$$

To derive the discrete system, first (17) is put into weak form. To this aim the test functions $(\psi_u, \psi_{\varepsilon}, \psi_w, \psi_{\kappa}, \psi)$ are introduced. For sake of simplicity, no dependency between the displacements and the energy is considered, i.e. $\delta_w H = 0$:

$$\langle \psi_{u}, \, \rho A \, \dot{e}_{u} \rangle_{\Omega} = \langle \psi_{u}, \, \partial_{x} e_{\varepsilon} \rangle_{\Omega} \, .$$

$$\langle \psi_{\varepsilon}, \, C_{a} \, \dot{e}_{\varepsilon} \rangle_{\Omega} = \langle \psi_{\varepsilon}, \, \partial_{x} e_{u} \rangle_{\Omega} + \langle \psi_{\varepsilon}, \, (\partial_{x} w) \, \partial_{x} e_{w} \rangle_{\Omega} \, ,$$

$$\langle \psi_{w}, \, \rho A \dot{e}_{w} \rangle_{\Omega} = \langle \psi_{w}, \, \partial_{x} (e_{\varepsilon} \partial_{x} w) \rangle_{\Omega} - \langle \psi_{w}, \, \partial_{xx}^{2} e_{\kappa} \rangle_{\Omega} \, ,$$

$$\langle \psi_{\kappa}, \, C_{b} \, \dot{e}_{\kappa} \rangle_{\Omega} = \langle \psi_{\kappa}, \, \partial_{xx}^{2} e_{w} \rangle_{\Omega} \, ,$$

$$\langle \psi, \, \dot{w} \rangle_{\Omega} = \langle \psi, \, e_{w} \rangle_{\Omega} \, .$$

$$(18)$$

Then the integration by parts is performed on the first, third and fourth line. This choice is such to retain the skew-symmetric structure at the discrete level and to lower the regularity requirement for the finite elements (Brugnoli, 2020, Chap. 8). The weak formulation then looks for $(e_u, e_w, e_\kappa, w) \in H^1(\Omega), e_\varepsilon \in L^2(\Omega)$ such that the following system

$$\langle \psi_{u}, \rho A \dot{e}_{u} \rangle_{\Omega} = -\langle \partial_{x} \psi_{u}, e_{\varepsilon} \rangle_{\Omega} + \langle \psi_{u}, e_{\varepsilon} \rangle_{\partial \Omega}.$$

$$\langle \psi_{\varepsilon}, C_{a} \dot{e}_{\varepsilon} \rangle_{\Omega} = \langle \psi_{\varepsilon}, \partial_{x} e_{u} \rangle_{\Omega} + \langle \psi_{\varepsilon}, \partial_{x} w \partial_{x} e_{w} \rangle_{\Omega},$$

$$\langle \psi_{w}, \rho A \dot{e}_{w} \rangle_{\Omega} = -\langle \partial_{x} \psi_{w} \partial_{x} w, e_{\varepsilon} \rangle_{\Omega} + \langle \partial_{x} \psi_{w}, \partial_{x} e_{\kappa} \rangle_{\Omega}$$

$$+ \langle \psi_{w}, e_{\varepsilon} \partial_{x} w - \partial_{x} e_{\kappa} \rangle_{\partial \Omega},$$

$$\langle \psi_{\kappa}, C_{b} \dot{e}_{\kappa} \rangle_{\Omega} = -\langle \partial_{x} \psi_{\kappa}, \partial_{x} e_{w} \rangle_{\Omega} + \langle \psi_{\kappa}, \partial_{x} e_{w} \rangle_{\partial \Omega},$$

$$\langle \psi, \dot{w} \rangle_{\Omega} = \langle \psi, e_{w} \rangle_{\Omega}.$$
(19)

holds $\forall (\psi_u, \psi_w, \psi_\kappa, \psi) \in H^1(\Omega), \forall \psi_\varepsilon \in L^2(\Omega)$. In this formulation, the boundary axial forces $e_\varepsilon|_0^L$, vertical forces $e_{\varepsilon}\partial_x w - \partial_x e_{\kappa}|_0^L$ and rotations $\partial_x e_w|_0^L$ are enforced weakly. To obtain the associated finite-dimensional system, the following Galerkin approximation is considered

$$e_{u}^{h} = \sum_{i=1}^{n_{u}} \xi_{u}^{i}(x)e_{u}^{i}(t), \qquad \psi_{u}^{h} = \sum_{i=1}^{n_{u}} \xi_{u}^{i}(x)\psi_{u}^{i},$$

$$e_{\varepsilon}^{h} = \sum_{i=1}^{n_{\varepsilon}} \xi_{\varepsilon}^{i}(x)e_{\varepsilon}^{i}(t), \qquad \psi_{\varepsilon}^{h} = \sum_{i=1}^{n_{\varepsilon}} \xi_{\varepsilon}^{i}(x)\psi_{\varepsilon}^{i},$$

$$e_{w}^{h} = \sum_{i=1}^{n_{w}} \xi_{w}^{i}(x)e_{w}^{i}(t), \qquad \psi_{w}^{h} = \sum_{i=1}^{n_{w}} \xi_{w}^{i}(x)\psi_{w}^{i}, \qquad (20)$$

$$e_{\kappa}^{h} = \sum_{i=1}^{n_{\kappa}} \xi_{\kappa}^{i}(x)e_{\kappa}^{i}(t), \qquad \psi_{\kappa}^{h} = \sum_{i=1}^{n_{\kappa}} \xi_{\kappa}^{i}(x)\psi_{\kappa}^{i},$$

$$w^{h} = \sum_{i=1}^{n_{w}} \xi_{w}^{i}(x)w^{i}(t), \qquad \psi^{h} = \sum_{i=1}^{n_{w}} \xi_{w}^{i}(x)\psi^{i}.$$

Notice that w, e_w are discretized using the same test functions $\xi_w^i(x)$. Plugging (20) into (19), the followingfinite dimensional system is obtained

$$\begin{pmatrix} \rho A \dot{e}_{u} \\ C_{u} \dot{e}_{\varepsilon} \\ \rho A \dot{e}_{w} \\ C_{b} \dot{e}_{\kappa} \\ \dot{w} \end{pmatrix} = \begin{bmatrix} 0 & \partial_{x} & 0 & 0 & 0 & 0 \\ \partial_{x} & 0 & \partial_{x} w \, \partial_{x} & 0 & 0 & 0 \\ 0 & \partial_{x} (\cdot \partial_{x} w) & 0 & -\partial_{xx}^{2} & -1 \\ 0 & 0 & \partial_{xx}^{2} & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \end{bmatrix} \begin{pmatrix} e_{u} \\ e_{\varepsilon} \\ e_{w} \\ \delta_{w} H \end{pmatrix}.$$

$$(17)$$
To derive the discrete system, first (17) is put into weak orm. To this aim the test functions $(\psi_{u}, \psi_{\varepsilon}, \psi_{w}, \psi_{\kappa}, \psi)$ re introduced. For sake of simplicity, no dependency etween the displacements and the energy is considered, e. $\delta_{w} H = 0$:
$$(\psi_{u}, \rho A \dot{e}_{u})_{\Omega} = \langle \psi_{u}, \partial_{x} e_{\varepsilon} \rangle_{\Omega}.$$

$$(\psi_{u}, \rho A \dot{e}_{u})_{\Omega} = \langle \psi_{u}, \partial_{x} e_{\varepsilon} \rangle_{\Omega}.$$

$$(21)$$

The fifth entry in the coenergy variables column disappear because $\delta_w H = 0$. The mass matrices are defined as follows

$$M_{u}^{ij} = \langle \xi_{u}^{i}, \rho A \xi_{u}^{j} \rangle_{\Omega}, \qquad M_{w}^{ij} = \langle \xi_{w}^{i}, \rho A \xi_{w}^{j} \rangle_{\Omega}, M_{\varepsilon}^{ij} = \langle \xi_{\varepsilon}^{i}, C_{a} \xi_{\varepsilon}^{j} \rangle_{\Omega}, \qquad M_{\kappa}^{ij} = \langle \xi_{\kappa}^{i}, C_{b} \xi_{\kappa}^{j} \rangle_{\Omega}.$$
 (22)

The interconnection matrices are given by

$$D_{\varepsilon u}^{ij} = \left\langle \xi_{\varepsilon}^{i}, \, \partial_{x} \xi_{u}^{j} \right\rangle_{\Omega},$$

$$D_{\varepsilon w}^{ij}(\mathbf{w}) = \left\langle \xi_{\varepsilon}^{i}, \, \sum_{k=1}^{n_{w}} \partial_{x} \xi_{w}^{k}(x) w^{k}(t) \partial_{x} \xi_{w}^{j} \right\rangle_{\Omega}, \qquad (23)$$

$$D_{w\kappa}^{ij} = \left\langle \partial_{x} \xi_{w}^{i}, \, \partial_{x} \xi_{\kappa}^{j} \right\rangle_{\Omega}.$$

The boundary matrices and control inputs are given by

$$B_{u}^{i} = [\psi_{u}^{i}(L) \quad \psi_{u}^{i}(0)],$$

$$B_{w}^{i} = [\psi_{w}^{i}(L) \quad \psi_{w}^{i}(0)],$$

$$B_{\kappa}^{i} = [\psi_{\kappa}^{i}(L) \quad \psi_{\kappa}^{i}(0)],$$

$$\mathbf{u}_{1} = [e_{\varepsilon}(L) \quad -e_{\varepsilon}(0)]^{\top},$$

$$\mathbf{u}_{2} = [e_{\varepsilon}\partial_{x}w - \partial_{x}e_{\kappa}(L) \quad -e_{\varepsilon}\partial_{x}w - \partial_{x}e_{\kappa}(0)]^{\top},$$

$$\mathbf{u}_{3} = [\partial_{x}e_{w}(L) \quad -\partial_{x}e_{w}(L)]^{\top}.$$

$$(24)$$

The outputs \mathbf{y}_i are easily identified thanks to power balance (15), as they corresponds to the power collocated variables w.r.t. the inputs. The discrete energy definition is given by the mass matrices weighted inner product of the coenergy variable

$$H_d = \frac{1}{2} (\mathbf{e}_u \mathbf{M}_u \mathbf{e}_u + \mathbf{e}_{\varepsilon} \mathbf{M}_{\varepsilon} \mathbf{e}_{\varepsilon} + \mathbf{e}_w \mathbf{M}_w \mathbf{e}_w + \mathbf{e}_{\kappa} \mathbf{M}_{\kappa} \mathbf{e}_{\kappa}). \tag{25}$$

The energy rate is computed taking into account the lossless dynamics of System (21) and the discrete energy definition (25):

$$\dot{H}_d = \sum_{i=1}^3 \mathbf{y}_i^{\mathsf{T}} \mathbf{u}_i. \tag{26}$$

For the numerical study in Sec. 5, homogeneous boundary conditions will be considered for simplicity.

For what concerns the choice of the underlying finite elements, a simple selection conforming to the weak form (19) is given by continuous Galerkin to discretize the space $H^1(\Omega)$ and discontinuous Galerkin for the space $L^2(\Omega)$. Consider an interval mesh \mathcal{I}_h with elements E. The space of polynomials of order k on a mesh cell is denoted by P_k . The following conforming families of finite elements are selected for this problem:

$$\mathcal{E}_{u}^{h} = \left\{ e_{u}^{h} \in H^{1}(\Omega) | \forall E \in \mathcal{I}_{h}, \ e_{u}^{h}|_{E} \in P_{2k-1} \right\},
\mathcal{E}_{\varepsilon}^{h} = \left\{ e_{\varepsilon}^{h} \in L^{2}(\Omega) | \forall E \in \mathcal{I}_{h}, \ e_{\varepsilon}^{h}|_{E} \in P_{2k-2} \right\},
\mathcal{E}_{w}^{h} = \left\{ e_{w}^{h} \in H^{1}(\Omega) | \forall E \in \mathcal{I}_{h}, \ e_{w}^{h}|_{E} \in P_{k} \right\},
\mathcal{E}_{\kappa}^{h} = \left\{ e_{\kappa}^{h} \in H^{1}(\Omega) | \forall E \in \mathcal{I}_{h}, \ e_{\kappa}^{h}|_{E} \in P_{k} \right\},
\mathcal{W}^{h} = \left\{ w^{h} \in H^{1}(\Omega) | \forall E \in \mathcal{I}_{h}, \ w^{h}|_{E} \in P_{k} \right\},$$
(27)

where $k \geq 1$. The space \mathcal{E}_u^h corresponds to the space of continous Galerkin elements of order 2k-1 whereas \mathcal{E}_w^h , \mathcal{E}_κ^h , \mathcal{W}^h corresponds to the space of continous Galerkin elements of order k. The space $\mathcal{E}_\varepsilon^h$ corresponds to the space of discontinuous Galerkin elements of order 2k-2. This choice guarantees that

$$\partial_x \mathcal{E}_u^h \subset \mathcal{E}_\varepsilon^h, \partial_x \mathcal{W}^h \cdot \partial_x \mathcal{E}_w^h \subset \mathcal{E}_\varepsilon^h.$$
(28)

So the discrete weak formulation becomes: find

$$(e_u^h, e_{\varepsilon}^h, e_w^h, e_{\kappa}^h, w^h) \in \mathcal{E}_u^h \times \mathcal{E}_{\varepsilon}^h \times \mathcal{E}_w^h \times \mathcal{E}_{\kappa}^h \times \mathcal{W}^h$$
 such that system (19) holds

$$\forall (\psi_u^h, \ \psi_\varepsilon^h, \ \psi_w^h, \ \psi_\kappa^h, \ \psi^h) \in \mathcal{E}_u^h \times \mathcal{E}_\varepsilon^h \times \mathcal{E}_w^h \times \mathcal{E}_\kappa^h \times \mathcal{W}^h.$$

5. NUMERICAL TEST

In this section the convergence rate of the underlying finite elements is assessed by means of a numerical test. A manufactured solution is considered in order to compare the numerical with an analytical manufactured one.

Consider the following manufactured solution for the axial and vertical displacement

$$u^{\text{ex}} = x^3 [1 - (x/L)^3] \sin(2\pi t), w^{\text{ex}} = \sin(\pi x/L) \sin(2\pi t),$$
(29)

together with the boundary conditions

$$u|_{0}^{L} = 0, \quad w|_{0}^{L} = 0, \quad m_{xx}|_{0}^{L} = 0.$$
 (30)

For u and w given in (29) to be the solution of (1), appropriate forcing terms have to be introduced in the right hand side of (1). These are given by

$$f_u = \rho A \partial_{tt}^2 u^{\text{ex}} - \partial_x n_{xx}^{\text{ex}},$$

$$f_w = \rho A \partial_{tt}^2 w^{\text{ex}} + \partial_{xx}^2 m_{xx}^{\text{ex}} - \partial_x (n_{xx}^{\text{ex}} \partial_x w^{\text{ex}}),$$
(31)

The forcing term are computed considering the projection of the analytical forcing terms on the corresponding test functions ψ_u , ψ_w . Given the finite element selection (27), this test function corresponds to the Lagrange elements of order 2k-1 and k for ψ_u and ψ_w respectively.

The numerical values of the parameters for the simulation are reported in Table 2. The Firedrake library (Rathgeber et al. (2017)) is used to generate the matrices. To integrate the equations in time a Crank-Nicholson scheme is employed. The time step is set to $\Delta t = h/(2\pi)$ to balance the time discretization error with respect to the spatial error. The final time is set to $t_f = 1[\mathbf{s}]$. The non-linear system is solved using Newton-Krylov iterations with a line search and direct factorization of the linear system.

To measure the convergence rate, suitable norms have to be introduced. Let \mathcal{X} be a Hilbert space, and t_f a positive real number. We denote by $L^{\infty}([0,t_f];\mathcal{X})$ or $L^{\infty}(\mathcal{X})$ the space of functions $f:[0,t_f]\to\mathcal{X}$ for which the time-space norm $||\cdot||_{L^{\infty}([0,t_f];\mathcal{X})}$ satisfies

$$||f||_{L^{\infty}([0,t_f];\mathcal{X})} = \underset{t \in [0,t_f]}{\text{ess sup}} ||f||_{\mathcal{X}} < \infty.$$

To compute the $L^{\infty}(\mathcal{X})$ space-time dependent norm from the numerical results, the discrete norm $L^{\infty}_{\Delta t}(\mathcal{X})$ is used

$$||\cdot||_{L^{\infty}(\mathcal{X})} \approx ||\cdot||_{L^{\infty}_{\Delta t}(\mathcal{X})} = \max_{t \in t_i} ||\cdot||_{\mathcal{X}},$$

where t_i are the discrete simulation instants. The Hilbert space \mathcal{X} depends on the regularity of the considered variable. For variables (e_u, e_w, e_κ, w) the error is measured in the $L^\infty_{\Delta t}(H^1)$ norm, whereas for e_ε is measured using the $L^\infty_{\Delta t}(L^2)$ norm. The simulation results are reported in Fig. 2. It can be noticed that all variables converge at rate given by the polynomial degree k. For variable e_κ a superconvergent trend can be observed for case k=1 (cf. Fig 2d). The numerical solutions for the different variables is plotted in Fig. 3.

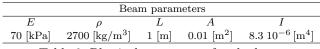


Table 2. Physical parameters for the beam.

6. CONCLUSION

In this contribution, a pH model for von Kármán beams is detailed. The resulting system of equations is then discretized using mixed finite elements. The validity of discretized model is assessed by comparison with an analytical solution. This demonstrates how pH formulations can be used in relevant non linear models arising from engineering. A natural outlook is the extension of this formulation to the two-dimensional case Brugnoli and Matignon (2021).

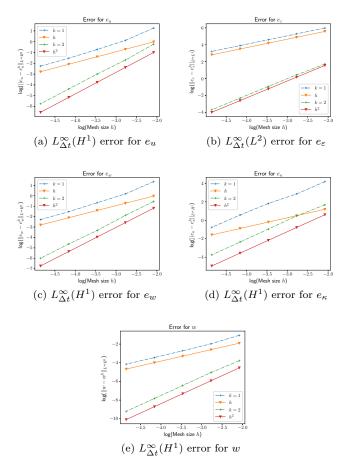


Fig. 2. Error trend for the different variables

The analysis of this model under a control theoretic perspective represents another interesting development. The numerical discrete model can be used to construct model based controllers.

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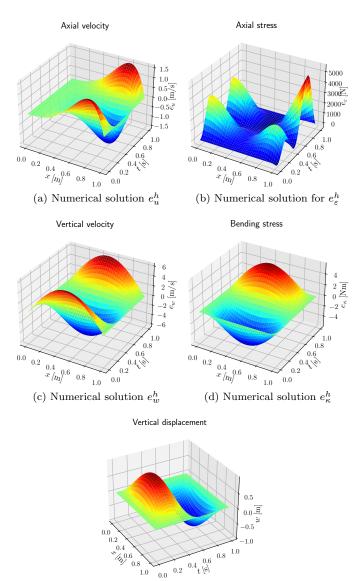


Fig. 3. Computed numerical solution for $h=2^{-5}$ and k=2 for the different variables

(e) Numerical solution w^h

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