A port-Hamiltonian formulation for the full von-Kármán plate model

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<u>Summary</u>. In this contribution, a port-Hamiltonian reformulation of the full von-Kármán dynamical model for geometrically non-linear plates is detailed, including the collocated boundary control and observation. Starting from the canonical equations, a set of variables is chosen so as to make the total energy quadratic. The model, reformulated in these variables, highlights a port-Hamiltonian structure ruled by a state-modulated interconnection operator.

Classical model

The classical full von-Kármán dynamical model is presented in Bilbao et al. [2015]. The problem, defined on an open connected set $\Omega \subset \mathbb{R}^2$, takes the dimensionless form

$$\ddot{\boldsymbol{u}} = \operatorname{Div} \boldsymbol{N},$$
 $\boldsymbol{N} = \boldsymbol{\Phi}(\boldsymbol{\varepsilon}),$ $\boldsymbol{\varepsilon} = \operatorname{Grad} \boldsymbol{u} + 1/2 \operatorname{grad} \boldsymbol{w} \otimes \operatorname{grad} \boldsymbol{w},$ $\ddot{\boldsymbol{w}} = -\operatorname{div} \operatorname{Div} \boldsymbol{M} + \operatorname{div} (\boldsymbol{N} \operatorname{grad} \boldsymbol{w}),$ $\boldsymbol{M} = \boldsymbol{\Phi}(\boldsymbol{\kappa}),$ $\boldsymbol{\kappa} = \operatorname{Grad} \operatorname{grad} \boldsymbol{w},$ (1)

where $\boldsymbol{u} \in \mathbb{R}^2$ is the in-plane displacement, \boldsymbol{w} is the vertical displacement, $\boldsymbol{\varepsilon}$ is the in-plane strain tensor, $\boldsymbol{\kappa}$ is the curvature tensor, \boldsymbol{N} is the in-plane stress resultant and \boldsymbol{M} is the bending stress resultant. The notation $\boldsymbol{a} \otimes \boldsymbol{b} = \boldsymbol{a} \boldsymbol{b}^{\top}$ denotes the dyadic product of two vectors. The div operator is the divergence of a vector field, and grad the gradient of a scalar field. The operator $\operatorname{Grad} = \frac{1}{2} \left(\nabla + \nabla^{\top} \right)$ designates the symmetric part of the gradient (i. e. the deformation gradient in continuum mechanics). For a tensor field $\boldsymbol{U}: \Omega \to \mathbb{R}^{2\times 2}$, with components U_{ij} , the divergence $\operatorname{Div}(\boldsymbol{U})$ is a vector, defined column-wise as

$$Div(U) := \sum_{i=1}^{2} \partial_{x_i} U_{ij}, \quad \forall j = \{1, 2\}.$$

The linear tensor mapping Φ is positive and preserves symmetry:

$$\Phi(A) = \nu \operatorname{Tr}(A)\mathbf{1} + (1 - \nu)A, \qquad A = A^{\top} \implies \Phi(A) = \Phi(A)^{\top}, \quad \text{where} \quad \mathbf{1} = \operatorname{Diag}(1, 1).$$

The total energy of the model (Hamiltonian functional)

$$H = \frac{1}{2} \int_{\Omega} \left\{ \|\dot{\boldsymbol{u}}\|^2 + \dot{w}^2 + \boldsymbol{N} : \boldsymbol{\varepsilon} + \boldsymbol{M} : \boldsymbol{\kappa} \right\} d\Omega, \quad \text{where} \quad \boldsymbol{A} : \boldsymbol{B} = \text{Tr}(\boldsymbol{A}^{\top} \boldsymbol{B})$$
 (2)

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).

The equivalent port-Hamiltonian system (pHs)

To find a suitable port-Hamiltonian system, we first select a set of new energy variables to make the Hamiltonian functional quadratic. The selection is the same as for both the linear plate problems in Brugnoli et al. [2019a,b]:

$$\alpha_u = \dot{u}, \qquad \alpha_w = \dot{w}, \qquad A_\varepsilon = \varepsilon, \qquad A_\kappa = \kappa.$$
 (3)

The energy is quadratic in these variables

$$H = \frac{1}{2} \int_{\Omega} \left\{ \|\boldsymbol{\alpha}_{u}\|^{2} + \alpha_{w}^{2} + \boldsymbol{\Phi}(\boldsymbol{A}_{\varepsilon}) : \boldsymbol{A}_{\varepsilon} + \boldsymbol{\Phi}(\boldsymbol{A}_{\kappa}) : \boldsymbol{A}_{\kappa} \right\}. \tag{4}$$

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

$$e_u := \delta_{\alpha_w} H = \dot{u}, \qquad e_w := \delta_{\alpha_w} H = \dot{w}, \qquad E_{\varepsilon} := \delta_{A_{\varepsilon}} H = \Phi(A_{\varepsilon}), \qquad E_{\kappa} := \delta_{A_{\kappa}} H = \Phi(A_{\kappa}).$$
 (5)

Before stating the final formulation, consider the operator $\mathcal{C}(w)(\cdot): L^2(\Omega, \mathbb{R}^{2\times 2}_{\mathrm{sym}}) \to L^2(\Omega)$ acting on symmetric tensors

$$C(w)(T) = \operatorname{div}(T\operatorname{grad} w). \tag{6}$$

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

$$C(w)^*(\cdot) = -\frac{1}{2} \left[\operatorname{grad}(\cdot) \otimes \operatorname{grad}(w) + \operatorname{grad}(w) \otimes \operatorname{grad}(\cdot) \right]. \tag{7}$$

Proof 1 Consider a smooth scalar field $v \in C_0^\infty(\Omega)$ and a smooth symmetric tensor field $U \in C_0^\infty(\Omega, \mathbb{R}^{2 \times 2}_{sym})$ with compact support. The formal adjoint of $\mathcal{C}(w)(\cdot)$ satisfies the relation

$$\langle v, \mathcal{C}(w)(U) \rangle_{L^2(\Omega)} = \langle \mathcal{C}(w)(v)^*, U \rangle_{L^2(\Omega, \mathbb{R}_{\text{cym}}^{2\times 2})}.$$
 (8)

The proof follows from the computation

$$\langle v, \mathcal{C}(w)(\boldsymbol{U}) \rangle_{L^{2}(\Omega)} = \langle v, \operatorname{div}(\boldsymbol{U} \operatorname{grad} w) \rangle_{L^{2}(\Omega)}, \quad \text{Integration by parts,}$$

$$= \langle -\operatorname{grad} v, \boldsymbol{U} \operatorname{grad} w \rangle_{L^{2}(\Omega, \mathbb{R}^{2})}, \quad \text{Dyadic product properties,}$$

$$= \langle -\operatorname{grad} v \otimes \operatorname{grad} w, \boldsymbol{U} \rangle_{L^{2}(\Omega, \mathbb{R}^{2 \times 2}_{sym})}, \quad \text{Symmetry of } \boldsymbol{U},$$

$$= \langle -1/2(\operatorname{grad} v \otimes \operatorname{grad} w + \operatorname{grad} w \otimes \operatorname{grad} v), \boldsymbol{U} \rangle_{L^{2}(\Omega, \mathbb{R}^{2 \times 2}_{sym})}.$$

$$(9)$$

This means

$$C(w)^*(\cdot) = -\frac{1}{2} \left[\operatorname{grad}(\cdot) \otimes \operatorname{grad}(w) + \operatorname{grad}(w) \otimes \operatorname{grad}(\cdot) \right], \tag{10}$$

leading to the final result.

The pH realization is then given by the following system

$$\frac{\partial}{\partial t} \begin{pmatrix} \boldsymbol{\alpha}_{u} \\ \boldsymbol{A}_{\varepsilon} \\ \boldsymbol{\alpha}_{w} \\ \boldsymbol{A}_{\kappa} \end{pmatrix} = \begin{bmatrix} \mathbf{0} & \text{Div} & \mathbf{0} & \mathbf{0} \\ \text{Grad} & \mathbf{0} & -\mathcal{C}(w)^{*} & \mathbf{0} \\ 0 & \mathcal{C}(w) & 0 & -\text{div} \, \text{Div} \\ \mathbf{0} & \mathbf{0} & \text{Grad} \, \text{grad} & \mathbf{0} \end{bmatrix} \begin{pmatrix} \delta_{\boldsymbol{\alpha}_{u}} H \\ \delta_{\boldsymbol{A}_{\varepsilon}} H \\ \delta_{\boldsymbol{\alpha}_{w}} H \\ \delta_{\boldsymbol{A}_{\kappa}} H \end{pmatrix}, \tag{11}$$

The second line of system (11) represents the time derivative of the membrane strain tensor. To close the system, variable w has to be accessible. For this reason, its dynamics has to be included. The augmented system reads

$$\frac{\partial}{\partial t} \begin{pmatrix} \boldsymbol{\alpha}_{u} \\ \boldsymbol{A}_{\varepsilon} \\ w \\ \boldsymbol{\alpha}_{w} \\ \boldsymbol{A}_{\kappa} \end{pmatrix} = \underbrace{\begin{bmatrix} \mathbf{0} & \text{Div} & \mathbf{0} & \mathbf{0} & \mathbf{0} \\ \text{Grad} & \mathbf{0} & \mathbf{0} & -\mathcal{C}(w)^{*} & \mathbf{0} \\ 0 & 0 & 0 & 1 & 0 \\ 0 & \mathcal{C}(w) & -1 & 0 & -\text{div} \, \text{Div} \\ \mathbf{0} & \mathbf{0} & \mathbf{0} & \text{Grad} \, \text{grad} & \mathbf{0} \end{bmatrix}}_{\mathcal{J}} \begin{pmatrix} \delta_{\boldsymbol{\alpha}_{u}} H \\ \delta_{\boldsymbol{A}_{\varepsilon}} H \\ \delta_{w} H \\ \delta_{\boldsymbol{\alpha}_{w}} H \end{pmatrix} . \tag{12}$$

Given the results in Brugnoli et al. [2019a,b] and Proposition 1, 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 terms allows accommodating other potentials, for example the gravitational one. Suitable boundary variables are then obtained considering the power balance

$$\dot{H} = \langle \gamma_0 \boldsymbol{e}_u, \, \gamma_\perp \boldsymbol{E}_\varepsilon \rangle_{\partial\Omega} + \langle \gamma_0 \boldsymbol{e}_w, \, \gamma_{\perp\perp,1} \boldsymbol{E}_\kappa + \gamma_0 (\boldsymbol{E}_\varepsilon \boldsymbol{n} \cdot \operatorname{grad} w) \rangle_{\partial\Omega} + \langle \gamma_1 \boldsymbol{e}_w, \, \gamma_{\perp\perp} \boldsymbol{E}_\kappa \rangle_{\partial\Omega},$$
(13)

where $\gamma_0 e_u = e_u|_{\partial\Omega}$ is the Dirichlet trace, $\gamma_\perp E_\varepsilon = E_\varepsilon n|_{\partial\Omega}$ is the normal trace (n is the outward normal vector), $\gamma_{\perp\perp,1} E_\kappa = -n \cdot \text{Div } E_\kappa - \partial_s (n^\top E_\kappa s)|_{\partial\Omega}$ is the effective shear force at the boundary (s is the tangent versor at the boundary), $\gamma_1 e_w = \partial_n e_w|_{\partial\Omega}$ is the normal derivative trace and $\gamma_{\perp\perp} E_\kappa = n^\top E_\kappa n$ is the normal to normal trace. The boundary conditions are consistent with the ones assumed in Puel and Tucsnak [1996] for deriving a global existence result for this model.

Conclusions

We have presented a pHs formulation of the full von-Kármán model. The dynamics of the system exhibits a state modulated interconnection operator, while the energy remains quadratic in the chosen variables. Of particular interest is the discretization of such a model for simulation and control purposes. The Partitioned Finite Element Method (PFEM), an extension of mixed finite elements to pHs, seems to be particularly suitable to achieve a structure-preserving discretization of this model, as in Cardoso-Ribeiro et al. [2020] for the 2D Shallow Water Equation, which exhibits the same kind of polynomial nonlinearity.

References

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