

A port-Hamiltonian formulation for the full Von-Karman plate model

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Summary. In this contribution, a port-Hamiltonian reformulation of the full von-Karman dynamical model for geometrically non-linear plates is detailed. Starting from the canonical equations, a set of variables is chosen so that that 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-Karman dynamical model is detailed in Bilbao et al. [2015]. The problem, defined on an open connected set $\Omega \subset \mathbb{R}^2$, takes the dimensionless form

$$\begin{aligned} \ddot{\mathbf{u}} &= \text{Div } \mathbf{N}, & \mathbf{N} &= \Phi(\boldsymbol{\varepsilon}), & \boldsymbol{\varepsilon} &= \text{Grad } \mathbf{u} + 1/2 \text{grad } w \otimes \text{grad } w, \\ \ddot{w} &= -\text{div Div } \mathbf{M} + \text{div}(\mathbf{N} \text{grad } w), & \mathbf{M} &= \Phi(\boldsymbol{\kappa}), & \boldsymbol{\kappa} &= \text{Grad grad } w, \end{aligned} \quad (1)$$

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

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

The tensor mapping Φ is positive and preserves the symmetry

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

Its inverse is given by

$$\Phi^{-1}(\mathbf{A}) = \frac{1}{1 - \nu}\mathbf{A} - \frac{\nu}{1 - \nu^2} \text{Tr}(\mathbf{A})\mathbf{1}. \quad (2)$$

The total energy of the model (Hamiltonian functional)

$$H = \frac{1}{2} \int_{\Omega} \left\{ \|\dot{\mathbf{u}}\|^2 + \dot{w}^2 + \mathbf{N} : \boldsymbol{\varepsilon} + \mathbf{M} : \boldsymbol{\kappa} \right\} d\Omega, \quad \text{where} \quad \mathbf{A} : \mathbf{B} = \text{Tr}(\mathbf{A}^\top \mathbf{B}) \quad (3)$$

consists of the kinetic energy and membrane and bending deformation energies. This model is conservative 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

To find a suitable port-Hamiltonian (pH) system, we first select a set of new variables to make the Hamiltonian functional quadratic. The selection is the same as for a linear plate problem Brugnoli [2020]

$$\boldsymbol{\alpha}_u = \dot{\mathbf{u}}, \quad \alpha_w = \dot{w}, \quad \mathbf{A}_\varepsilon = \boldsymbol{\varepsilon}, \quad \mathbf{A}_\kappa = \boldsymbol{\kappa}. \quad (4)$$

The energy is quadratic in these variables

$$H = \frac{1}{2} \int_{\Omega} \left\{ \|\boldsymbol{\alpha}_u\|^2 + \alpha_w^2 + \Phi(\mathbf{A}_\varepsilon) : \mathbf{A}_\varepsilon + \Phi(\mathbf{A}_\kappa) : \mathbf{A}_\kappa \right\}. \quad (5)$$

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

$$\mathbf{e}_u := \delta_{\boldsymbol{\alpha}_u} H = \dot{\mathbf{u}}, \quad e_w := \delta_{\alpha_w} H = \dot{w}, \quad \mathbf{E}_\varepsilon := \delta_{\mathbf{A}_\varepsilon} H = \Phi(\mathbf{A}_\varepsilon), \quad \mathbf{E}_\kappa := \delta_{\mathbf{A}_\kappa} H = \Phi(\mathbf{A}_\kappa) \quad (6)$$

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Conclusions

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References

Stefan Bilbao, Olivier Thomas, Cyril Touzé, and Michele Ducceschi. Conservative numerical methods for the full von kármán plate equations. *Numerical Methods for Partial Differential Equations*, 31(6):1948–1970, 2015. doi: 10.1002/num.21974. URL <https://onlinelibrary.wiley.com/doi/abs/10.1002/num.21974>.

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