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Stability of the Rao-Nakra sandwich beam with a dissipation of fractional derivative type: theoretical and numerical study

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STABILITY OF THE RAO–NAKRA SANDWICH BEAM WITH A DISSIPATION OF FRACTIONAL DERIVATIVE TYPE: THEORETICAL AND NUMERICAL STUDY

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ABSTRACT. This paper is devoted to the solution and stability of a one-dimensional model depicting Rao–Nakra sandwich beams, incorporating damping terms characterized by fractional derivative types within the domain, specifically a generalized Caputo derivative with exponential weight. To address existence, uniqueness, stability, and numerical results, fractional derivatives are substituted by diffusion equations relative to a new independent variable, ξ , resulting in an augmented model with a dissipative semigroup operator. Polynomial decay of energy is achieved, with a decay rate depending on the fractional derivative parameters. Both the polynomial decay and its dependency on the parameters of the generalized Caputo derivative are numerically validated. To this end, an energy-conserving finite difference numerical scheme is employed.

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Discussion and Conclusions References

1. INTRODUCTION

In 1999, Liu, Trogdon and Yong [28] developed the following general three-layer laminated beam–plate model:

(1.1)
$$\begin{cases} \varrho_1 h_1 u_{\mathbf{tt}} - E_1 h_1 u_{xx} - \tau = 0, \\ \varrho_3 h_3 v_{\mathbf{tt}} - E_3 h_3 u_{xx} + \tau = 0, \\ \varrho h w_{\mathbf{tt}} + E I w_{xxxx} - G_1 h_1 (w_x + \phi_1)_x - G_3 h_3 (w_x + \phi_3)_x - h_2 \tau_x = 0, \\ \varrho_1 I_1 \phi_{1\mathbf{tt}} - E_1 I_1 \phi_{1xx} - \frac{h_1}{2} \tau + G_1 h_1 (w_x + \phi_1) = 0, \\ \varrho_3 I_3 \phi_{3\mathbf{tt}} - E_3 I_3 \phi_{3xx} - \frac{h_3}{2} \tau + G_3 h_3 (w_x + \phi_3) = 0. \end{cases}$$

The physical parameters h_i , ϱ_i , E_i , G_i , $I_i > 0$ are the thickness, density, Young's modulus, shear modulus and moments of inertia of the *i*-th layer for i = 1, 2, 3, from bottom to top, respectively. In addition, $\varrho h = \varrho_1 h_1 + \varrho_3 h_3$ and $EI = E_1 I_1 + E_3 I_3$.

The Rao–Nakra sandwich beam model [45] is the following:

(1.2)
$$\begin{cases} \rho_1 u_{\mathbf{tt}}(x, \mathbf{t}) - \vartheta u_{xx}(x, \mathbf{t}) - k(-u(x, \mathbf{t}) + v(x, \mathbf{t}) + \gamma w_x(x, \mathbf{t})) = 0, \\ \rho_2 v_{\mathbf{tt}}(x, \mathbf{t}) - \chi v_{xx}(x, \mathbf{t}) + k(-u(x, \mathbf{t}) + v(x, \mathbf{t}) + \gamma w_x(x, \mathbf{t})) = 0, \\ \rho_3 w_{\mathbf{tt}}(x, \mathbf{t}) + \zeta w_{xxxx}(x, \mathbf{t}) - k(-u(x, \mathbf{t}) + v(x, \mathbf{t}) + \gamma w_x(x, \mathbf{t}))_x = 0, \\ u(x, 0) = u_0(x), \ u_{\mathbf{t}}(x, 0) = u_1(x), \\ v(x, 0) = v_0(x), \ v_{\mathbf{t}}(x, 0) = v_1(x), \\ w(x, 0) = w_0(x), \ w_{\mathbf{t}}(x, 0) = w_1(x), \end{cases}$$

with $0 < x < \ell$ and $\mathbf{t} > 0$. Here $u = u(x, \mathbf{t})$ and $v = v(x, \mathbf{t})$ are the longitudinal displacement. $w = w(x, \mathbf{t})$ is the transverse displacement of the beam. This system is derived from the system (1.1) when we consider the core material to be linearly elastic, that is, $\tau = 2G_2\gamma$ with the shear strain

$$\gamma = \frac{1}{2h_2}(-u + v + \alpha w_x)$$
 and $\alpha = h_2 = \frac{1}{2}(h_1 + h_3)$

where $k = \frac{G_2}{h_2}$, the shear modulus $G_2 = \frac{E_2}{2(1+\nu)}$, and $-1 < \nu < \frac{1}{2}$ is the Poisson ratio.

Many authors have been working on the system (1.2) from different points of view: see [27, 39, 40, 46] and the references therein. In the past three decades there has been a growing interest by many reserchers for the study of fractional calculus in different fields of sciences [32, 48, 49]. Several aspects of engineering, applied sciences, and mathematical physics benefitted from this ascending wave of applications. Space sciences, fluid mechanics, porous media flows, viscoelastic and biological processes are but a few areas in which fractional order differential equations have become a favored tool to tread new path. To give some examples, fractional derivatives have

been used to model frequency dependent damping behavior of many viscoelastic materials as well modeling many chemical processes. Many problems in several scientific applied areas, including the analysis of viscoelastic materials, heat conduction in materials with memory, electrodynamics with memory, signal processing, among others, can be modeled by fractional differential calculus. Indeed, many investigations have shown that models involving fractional derivatives are more realistic to represent some natural phenomena than models involving classical derivatives. For more information we refer to [24, 34, 47] and the references therein. For example, the wave equation with boundary fractional damping has been treated in [34, 35] where the strong stability and the lack of uniform stabilization were proved. Applications of this kind are also found in [33, 36, 42, 44, 50] and the references therein.

In [50] the authors studied the Rao–Nakra system with a boundary dissipation of fractional derivative type. They established the polynomial stability of the system. In this article we are interested in studying the polynomial stabilization to the system (1.3) from a theoretical and numerical point of view. In order to achieve this, we eliminate the dissipation given on the border and place the dissipation in the domain. The main idea to reach this result is to use an idea given in [4]. It is important to note that the imposed dissipation is of a fractional derivative type. Our method is based on the fact that the input–output relationship in a certain diffusion equation introduced by Mbodje in [34] is realized by a fractional derivative operator.

Numerical approximations of derivatives and fractional integrals have been extensively studied, developed, and refined. Comprehensive reviews on this topic are available in [11, 26]. Many of these approaches involve direct approximations of the Riemann–Liouville, Caputo [18, 31, 13, 14, 15], or Grünwald–Letnikov integral [26], using quadrature methods, finite differences, and truncated summations [19]. While some progress has been made in enhancing the convergence rate, such as through trapezoidal approximation [26] or spline interpolation [17], these methods do not inherently preserve the energy conservation property of the Rao–Nakra model under study, nor do they maintain stability in the presence of dissipative terms. In this work, we address this challenge by employing a suitable combination of the β -Newmark method [38], coupled with a Crank–Nicolson method for the Mbodje diffusion equation, to achieve our objective. Without lost of generality, we will consider the following system for $\rho_1 = \rho_2 = \rho_3 = 1$:

$$(1.3) \begin{cases} u_{\mathbf{tt}} - \vartheta u_{xx} - k(-u+v+\gamma w_x) + \partial_{\mathbf{t}}^{\alpha,\,\eta} u = 0, \, (0,\ell) \times (0,+\infty), \\ v_{\mathbf{tt}} - \chi v_{xx} + k(-u+v+\gamma w_x) + \partial_{\mathbf{t}}^{\alpha,\,\eta} v = 0, \, (0,\ell) \times (0,+\infty), \\ w_{\mathbf{tt}} + \zeta w_{xxxx} - k(-u+v+\gamma w_x)_x + \partial_{\mathbf{t}}^{\alpha,\,\eta} w = 0, \, (0,\ell) \times (0,+\infty), \\ u(x,\,0) = u_0(x), \, u_{\mathbf{t}}(x,\,0) = u_1(x), \, (0,\ell), \\ v(x,\,0) = v_0(x), \, v_{\mathbf{t}}(x,\,0) = v_1(x), \, (0,\ell), \\ w(x,\,0) = w_0(x), \, w_{\mathbf{t}}(x,\,0) = w_1(x), \, (0,\ell), \\ u(0,\,\mathbf{t}) = u(\ell,\,\mathbf{t}) = 0, \, v(0,\,\mathbf{t}) = v(\ell,\,\mathbf{t}) = 0, \, (0,+\infty), \\ w(0,\,\mathbf{t}) = w(\ell,\,\mathbf{t}) = w_x(0,\,\mathbf{t}) = w_x(\ell,\,\mathbf{t}) = 0, \, (0,+\infty), \end{cases}$$

with $0 < x < \ell$, $\mathbf{t} > 0$, and real-valued functions $u = u(x, \mathbf{t})$ and $v = v(x, \mathbf{t})$, $w = w(x, \mathbf{t})$.

The rest of the paper is divided into five sections. In Section 1 we show that the system (1.3) may be replaced by an augmented system (2.10) obtained by coupling an equation with a suitable diffusion, and we study the energy functional associated to system. In Section 3, we establish the existence and uniqueness of solutions of the system (2.10); for this we use [5, 4]. In section 4 we prove the strong stability, the lack of uniform stabilization and the polynomial stability of the system (2.10). In Section 5, we give a numerical study of the polynomial stability.

Throughout this paper, C is a generic constant, not necessarily the same at each occasion (it may change from line to line) and depending on the indicated quantities.

2. Augmented Model

This section is devoted to the reformulation of the model (1.3) with boundary conditions $(1.3)_{7,8}$ as an augmented system. First we provide a brief review of fractional calculus. There are many definitions of fractional derivatives [18], among them those of Riemann–Liouville and Caputo are the most widely used [31]. The latter has the same Laplace transform as the integer order derivative, so it is widely used in control theory. In this paper the fractional derivative damping force is regarded as a control force to study the properties of free damped vibrations of the system, hence the Caputo definition [13, 14, 15] is used here. Let $0 < \alpha < 1$. The Caputo fractional integral of order α is defined by the formula

(2.1)
$$I^{\alpha}f(\mathbf{t}) = \frac{1}{\Gamma(\alpha)} \int_0^{\mathbf{t}} (\mathbf{t} - s)^{\alpha - 1} f(s) ds,$$

where Γ is the Gamma function, $f \in L^1(0, +\infty)$ and $\mathbf{t} > 0$.

The Caputo fractional derivative operator of order α is defined by the expression

(2.2)
$$D^{\alpha}f(\mathbf{t}) = I^{1-\alpha}Df(\mathbf{t}) := \frac{1}{\Gamma(1-\alpha)} \int_0^{\mathbf{t}} (\mathbf{t}-s)^{-\alpha}f'(s)ds.$$

The Caputo definition of fractional derivative possesses a very simple but interesting interpretation: if the function $f(\mathbf{t})$ represents the strain history within a viscoelastic material whose relaxation function is $[\Gamma(1-\alpha)\mathbf{t}^{\alpha}]^{-1}$, then the material will experience at any time \mathbf{t} a total stress given by the expression $D^{\alpha}f(\mathbf{t})$. Also, it easy to show that D^{α} is a left inverse of I^{α} , but in general it is not a right inverse. More precisely, we have

$$D^{\alpha}I^{\alpha}f = f, \qquad I^{\alpha}D^{\alpha}f(\mathbf{t}) = f(\mathbf{t}) - f(0).$$

See [47] for the proof of above equalities and for more properties of fractional calculus.

In this work, we consider a slightly different version of (2.1) and (2.4). In [16], Choi and MacCamy introduced the following definition of fractional integro-differential operators with an

exponential weight. Given $0 < \alpha < 1$ and $\eta \ge 0$, the exponential fractional integral of order α is defined by

(2.3)
$$I^{\alpha, \eta} f(\mathbf{t}) = \frac{1}{\Gamma(\alpha)} \int_0^{\mathbf{t}} e^{-\eta(\mathbf{t}-s)} (\mathbf{t}-s)^{\alpha-1} f(s) ds,$$

and the exponential fractional derivative operator of order α is defined by

(2.4)
$$\partial_{\mathbf{t}}^{\alpha, \eta} f(\mathbf{t}) = \frac{1}{\Gamma(1-\alpha)} \int_0^{\mathbf{t}} e^{-\eta (\mathbf{t}-s)} (\mathbf{t}-s)^{-\alpha} f'(s) ds.$$

Note that $\partial_{\mathbf{t}}^{\alpha, \eta} f(\mathbf{t}) = I^{1-\alpha, \eta} f'(\mathbf{t}).$

We will need later the following results:

Theorem 2.1. [34] Let μ be the function

(2.5)
$$\mu(\xi) = |\xi|^{(2\alpha - 1)/2}, \quad \xi \in \mathbb{R}, \quad 0 < \alpha < 1.$$

Then the relation between the Input U and the Output O is given by the following system:

(2.6)
$$\varphi_{\mathbf{t}}(\mathbf{t},\xi) + |\xi|^2 \varphi(\mathbf{t},\xi) = \mu(\xi) U(\mathbf{t}), \qquad \xi \in \mathbb{R}, \qquad \mathbf{t} > 0,$$

(2.7)
$$\varphi(0,\,\xi) = 0,$$

(2.8)
$$O = \pi^{-1} \sin(\alpha \pi) \int_{\mathbb{R}} \mu(\xi) \varphi(\mathbf{t}, \, \xi) d\xi$$

This implies that

$$(2.9) O = I^{1-\alpha}U,$$

where $U \in C([0, +\infty))$.

Lemma 2.2. [1, Lemma 2.1] If $\lambda \in D = \{\lambda \in \mathbb{C} : Re \ \lambda + \eta > 0 \text{ or } Im \ \lambda \neq 0\}$, then

$$\int_{\mathbb{R}} \frac{\mu^2(\xi)}{\xi^2 + \eta + \lambda} \, d\xi = \frac{\pi}{\sin(\alpha \pi)} (\eta + \lambda)^{\alpha - 1}.$$

Our strategy requires the elimination of the fractional derivatives in time from the domain condition in the system (1.3). To this, setting $\mu(\xi) = |\xi|^{(2\alpha-1)/2}$, $\xi \in \mathbb{R}$, $\mathfrak{C} = \pi^{-1} \sin(\alpha \pi)$, and exploiting the technique from [21], we reformulate the system (1.3) by using Theorem 2.1. This leads to the following augmented model, where $(x, t, \xi) \in (0, \ell) \times (0, +\infty) \times \mathbb{R}$ and we introduce

the constant $\mathfrak{C} := \pi^{-1} \sin(\alpha \pi)$:

$$(2.10) \begin{cases} u_{\mathbf{tt}}(x,\mathbf{t}) - \vartheta u_{xx}(x,\mathbf{t}) - k(-u(x,\mathbf{t}) + v(x,\mathbf{t}) + \gamma w_{x}(x,\mathbf{t})) \\ + \mathfrak{C} \int_{\mathbb{R}} \mu(\xi)\varphi(x,\mathbf{t},\xi)d\xi = 0, \\ \varphi_{\mathbf{t}}(x,\mathbf{t},\xi) + (\xi^{2} + \eta)\varphi(x,\mathbf{t},\xi) = \mu(\xi)u_{\mathbf{t}}(x,\mathbf{t}), \\ v_{\mathbf{tt}}(x,\mathbf{t}) - \chi v_{xx}(x,\mathbf{t}) + k(-u(x,\mathbf{t}) + v(x,\mathbf{t}) + \gamma w_{x}(x,\mathbf{t})) \\ + \mathfrak{C} \int_{\mathbb{R}} \mu(\xi)\phi(x,\mathbf{t},\xi)d\xi = 0, \\ \phi_{\mathbf{t}}(x,\mathbf{t},\xi) + (\xi^{2} + \eta)\phi(x,\mathbf{t},\xi) = \mu(\xi)v_{\mathbf{t}}(x,\mathbf{t}), \\ w_{\mathbf{tt}}(x,\mathbf{t}) + \zeta w_{xxxx}(x,\mathbf{t}) - k(-u(x,\mathbf{t}) + v(x,\mathbf{t}) + \gamma w_{x}(x,\mathbf{t})))_{x} \\ + \mathfrak{C} \int_{\mathbb{R}} \mu(\xi)\psi(x,\mathbf{t},\xi)d\xi = 0, \\ \psi_{\mathbf{t}}(x,\mathbf{t},\xi) + (\xi^{2} + \eta)\psi(x,\mathbf{t},\xi) = \mu(\xi)w_{\mathbf{t}}(x,\mathbf{t}), \\ u(x,0) = u_{0}(x), u_{\mathbf{t}}(x,0) = u_{1}(x), \\ v(x,0) = v_{0}(x), v_{\mathbf{t}}(x,0) = u_{1}(x), \\ w(x,0) = w_{0}(x), w_{\mathbf{t}}(x,0) = w_{1}(x), \\ u(0,t) = u(\ell,\mathbf{t}) = 0, v(0,\mathbf{t}) = v(\ell,\mathbf{t}) = 0, \\ w(0,t) = w(\ell,\mathbf{t}) = w_{x}(0,\mathbf{t}) = w_{x}(\ell,\mathbf{t}) = 0, \\ \varphi(x,0,\xi) = 0, \phi(x,0,\xi) = 0, \psi(x,0,\xi) = 0. \end{cases}$$

3. Setting of the Semigroup

In this section, we discuss the existence and uniqueness of solutions for the coupled system (2.10) by the semigroup theory [41, 43].

We will use the standard $L^2(0, \ell)$ space; the scalar product and the norm are denoted by

$$\langle \varphi, \psi \rangle_{L^2(0,\ell)} = \int_0^\ell \varphi \,\overline{\psi} \, dx \text{ and } \|\psi\|_{L^2(0,\ell)}^2 = \int_0^\ell |\psi|^2 dx.$$

We introduce the Hilbert space

equipped with the following inner product:

$$\begin{aligned} \langle \mathcal{U}, \tilde{\mathcal{U}} \rangle_{\mathcal{H}} &= \int_{0}^{\ell} U \overline{\tilde{U}} dx + \int_{0}^{\ell} V \overline{\tilde{V}} dx + \int_{0}^{\ell} W \overline{\tilde{W}} dx \\ &+ \vartheta \int_{0}^{\ell} u_{x} \overline{\tilde{u}}_{x} dx + \chi \int_{0}^{\ell} v_{x} \overline{\tilde{v}}_{x} dx + \zeta \int_{0}^{\ell} w_{xx} \overline{\tilde{w}}_{xx} dx \\ &+ k \int_{0}^{\ell} (-u + v + \gamma w_{x}) (-\overline{\tilde{u}} + \overline{\tilde{v}} + \gamma \overline{\tilde{w}}_{x}) dx \\ &+ \mathfrak{E} \int_{\mathbb{R}} \langle \varphi, \, \tilde{\varphi} \rangle_{L^{2}(0,\,\ell)} d\xi + \mathfrak{E} \int_{\mathbb{R}} \langle \phi, \, \tilde{\phi} \rangle_{L^{2}(0,\,\ell)} d\xi + \mathfrak{E} \int_{\mathbb{R}} \langle \psi, \, \tilde{\psi} \rangle_{L^{2}(0,\,\ell)} d\xi, \end{aligned}$$

$$(3.2)$$

where $\mathcal{U} = (u, v, w, U, V, W, \varphi, \phi, \psi)^T$ and $\widetilde{\mathcal{U}} = (\tilde{u}, \tilde{v}, \tilde{w}, \tilde{U}, \tilde{V}, \tilde{W}, \tilde{\varphi}, \tilde{\phi}, \tilde{\psi})^T$. We wish to transform the initial boundary value problem (2.10) into an abstract Cauchy problem in the Hilbert space \mathcal{H} . For this we introduce the functions $u_{\mathbf{t}} = U, v_{\mathbf{t}} = V, w_{\mathbf{t}} = W$, and we rewrite the system (2.10) as the following initial value problem:

(3.3)
$$\frac{d\mathcal{U}}{d\mathbf{t}}(\mathbf{t}) = \mathcal{A}\mathcal{U}(\mathbf{t}), \quad \mathcal{U}(0) = \mathcal{U}_0, \quad \forall \mathbf{t} > 0.$$

Here \mathcal{U} is as above, $\mathcal{U}_0 = (u_0, v_0, w_0, u_1, v_1, w_0, 0, 0, 0)^T$, and the operator $\mathcal{A} : \mathcal{D}(\mathcal{A}) \subset \mathcal{H} \to \mathcal{H}$ is given by the formula

.

with the domain

$$\begin{split} \mathcal{D}(\mathcal{A}) &= \big\{ \mathcal{U} = (u, v, w, U, V, W, \varphi, \phi, \psi) \in \mathcal{H} : \ U, V \in H_0^1(0, \ell), W \in H_0^2(0, \ell), \\ &\quad \vartheta u_{xx} + k(-u + v + \gamma w_x) - \mathfrak{C} \int_{\mathbb{R}} \mu(\xi) \varphi(x, \xi) d\xi \in L^2(0, \ell), \\ &\quad \chi v_{xx} - k(-u + v + \gamma w_x) - \mathfrak{C} \int_{\mathbb{R}} \mu(\xi) \phi(x, \xi) d\xi \in L^2(0, \ell), \\ &\quad - \zeta w_{xxxx} + k\gamma(-u + v + \gamma w_x)_x - \mathfrak{C} \int_{\mathbb{R}} \mu(\xi) \psi(x, \xi) d\xi \in L^2(0, \ell), \\ &\quad - (\xi^2 + \eta) \varphi(x, \xi) + \mu(\xi) U(x) \in L^2(\mathbb{R}, L^2(0, \ell)) \\ &\quad - (\xi^2 + \eta) \psi(x, \xi) + \mu(\xi) V(x) \in L^2(\mathbb{R}, L^2(0, \ell)) \\ &\quad - (\xi^2 + \eta) \psi(x, \xi) + \mu(\xi) W(x) \in L^2(\mathbb{R}, L^2(0, \ell)) \\ &\quad - (\xi^2 + \eta) \psi(x, \xi) + \mu(\xi) W(x) \in L^2(\mathbb{R}, L^2(0, \ell)) \big\}. \end{split}$$

Let the operator $A: \mathcal{D}(A) \subset H \to H$ which is given by

(3.5)
$$A\begin{pmatrix} u\\ v\\ w \end{pmatrix} = \begin{pmatrix} -\vartheta u_{xx} - k(-u+v+\gamma w_x)\\ -\chi v_{xx} + k(-u+v+\gamma w_x)\\ \zeta w_{xxxx} - k(-u+v+\gamma w_x)_x \end{pmatrix},$$

with the domain $\mathcal{D}(A) = \left[H^2(0,\ell) \cap H^1_0(0,\ell)\right] \times \left[H^2(0,\ell) \cap H^1_0(0,\ell)\right] \times \left[H^4(0,\ell) \cap H^2_0(0,\ell)\right],$ $H = (L^2(0,\ell))^3.$

We note that the operator A is self-adjoint and strictly positive.

So, the problem (1.3) can be rewritten as following, as in [4, 5]:

(3.6)
$$\begin{cases} Z_{\mathbf{tt}}(\mathbf{t}) + AZ(\mathbf{t}) + BB^* \partial_{\mathbf{t}}^{\alpha,\eta} Z(\mathbf{t}) = 0, \ \mathbf{t} > 0\\ Z(0) = Z_0, Z_{\mathbf{t}}(0) = Z_1, \end{cases}$$

where $B = B^* = I_H, Z(\mathbf{t}) = (u(\mathbf{t}), v(\mathbf{t}), w(\mathbf{t}))^T$ and $Z_0 = (u_0, v_0, w_0)^T, \quad Z_1 = (u_1, v_1, w_1)^T.$ We define

(3.7)
$$\mathcal{H}_0 = H_0^1(0,\,\ell) \times H_0^1(0,\,\ell) \times H_0^2(0,\,\ell) \times L^2(0,\,\ell) \times L^2(0,\,\ell) \times L^2(0,\,\ell)$$

equipped with the inner product given by

$$\langle \mathcal{U}, \tilde{\mathcal{U}} \rangle_{\mathcal{H}_{0}} = \int_{0}^{\ell} U \overline{\tilde{\mathcal{U}}} dx + \int_{0}^{\ell} V \overline{\tilde{\mathcal{V}}} dx + \int_{0}^{\ell} W \overline{\tilde{\mathcal{W}}} dx + \vartheta \int_{0}^{\ell} u_{x} \overline{\tilde{u}}_{x} dx + \chi \int_{0}^{\ell} v_{x} \overline{\tilde{v}}_{x} dx + \zeta \int_{0}^{\ell} w_{xx} \overline{\tilde{w}}_{xx} dx + k \int_{0}^{\ell} (-u + v + \gamma w_{x}) (-\overline{\tilde{u}} + \overline{\tilde{v}} + \gamma \overline{\tilde{w}}_{x}) dx,$$

$$(3.8)$$

where $\mathcal{U} = (u, v, w, U, V, W)^T$ and $\tilde{\mathcal{U}} = (\tilde{u}, \tilde{v}, \tilde{w}, \tilde{U}, \tilde{V}, \tilde{W})^T$.

Then, the operator $\mathcal{A}_0 : \mathcal{D}(\mathcal{A}_0) \subset \mathcal{H}_0 \to \mathcal{H}_0$ given by

(3.9)
$$\mathcal{A}_{0} \begin{pmatrix} u \\ v \\ w \\ U \\ U \\ V \\ W \end{pmatrix} = \begin{pmatrix} U \\ V \\ \vartheta u_{xx} + k(-u+v+\gamma w_{x}) - U \\ \chi v_{xx} - k(-u+v+\gamma w_{x}) - V \\ -\zeta w_{xxxx} + k(-u+v+\gamma w_{x})_{x} - W \end{pmatrix}.$$

with the domain

$$\mathcal{D}(\mathcal{A}_0) = [H^2(0,\,\ell) \times H^1_0(0,\,\ell)]^2 \times [H^4(0,\,\ell) \cap H^2_0(0,\,\ell)] \times H^2_0(0,\,\ell).$$

We recall two propositions from [4, 5]:

Proposition 3.1. [5, Proposition 2.6.2] The operator \mathcal{A}_0 generates a C_0 -semigroup of contractions in the Hilbert space \mathcal{H}_0 . Moreover, the following auxiliary problem:

$$(3.10) \begin{cases} u_{\mathbf{tt}}(\mathbf{t}) - \vartheta u_{xx}(\mathbf{t}) - k(-u(\mathbf{t}) + v(\mathbf{t}) + \gamma w_{x}(\mathbf{t})) + u_{\mathbf{t}}(\mathbf{t}) = 0, \\ v_{\mathbf{tt}}(\mathbf{t}) - \chi v_{xx}(\mathbf{t}) + k(-u(\mathbf{t}) + v(\mathbf{t}) + \gamma w_{x}(\mathbf{t})) + v_{\mathbf{t}}(\mathbf{t}) = 0, \\ w_{\mathbf{tt}}(\mathbf{t}) + \zeta w_{xxxx}(\mathbf{t}) - k\gamma(-u(\mathbf{t}) + v(\mathbf{t}) + \gamma w_{x}(\mathbf{t}))_{x} + w_{\mathbf{t}}(\mathbf{t}) = 0, \\ u(0, \mathbf{t}) = u(\ell, \mathbf{t}) = 0, \ v(0, \mathbf{t}) = v(\ell, \mathbf{t}) = 0, \\ w(0, \mathbf{t}) = w(\ell, \mathbf{t}) = w_{x}(0, \mathbf{t}) = w_{x}(\ell, \mathbf{t}) = 0, \\ u(x, 0) = u_{0}(x), \quad u_{\mathbf{t}}(x, 0) = u_{1}(x), \\ v(x, 0) = v_{0}(x), \quad w_{\mathbf{t}}(x, 0) = v_{1}(x), \\ w(x, 0) = w_{0}(x), \quad w_{\mathbf{t}}(x, 0) = w_{1}(x), \end{cases}$$

admits a unique solution $(u(x, \mathbf{t}), v(x, \mathbf{t}), w(x, \mathbf{t}))$ such that if $(u_0, v_0, w_0, u_1, v_1, w_1) \in \mathcal{D}(\mathcal{A}_0)$, then the solution $(u(x, \mathbf{t}), v(x, \mathbf{t}), w(x, \mathbf{t}))$ of (3.10) verifies the following regularity property:

$$\mathcal{U} = (u, v, w, u_{\mathbf{t}}, v_{\mathbf{t}}, w_{\mathbf{t}}) \in C([0, +\infty) : \mathcal{D}(\mathcal{A}_0)) \cap C^1([0, +\infty), \mathcal{H}_0),$$

and when $(u_0, v_0, w_0, u_1, v_1, w_1) \in \mathcal{H}_0$, then

$$\mathcal{U} = (u, v, w, u_{\mathbf{t}}, v_{\mathbf{t}}, w_{\mathbf{t}}) \in C([0, +\infty), \mathcal{H}_0).$$

The energy of the solution of the system (3.10) is defined as follows:

$$E_{0}(\mathbf{t}) = \frac{1}{2} \left[\|u_{\mathbf{t}}(\mathbf{t})\|_{L^{2}(0,\ell)}^{2} + \|v_{\mathbf{t}}(\mathbf{t})\|_{L^{2}(0,\ell)}^{2} + \|w_{\mathbf{t}}(\mathbf{t})\|_{L^{2}(0,\ell)}^{2} \\ + \vartheta \|u(\mathbf{t})\|_{H_{0}^{1}(0,\ell)}^{2} + \chi \|v(\mathbf{t})\|_{H_{0}^{1}(0,\ell)}^{2} + \zeta \|w(\mathbf{t})\|_{H_{0}^{2}(0,\ell)}^{2} \\ + k\| - u(\mathbf{t}) + v(\mathbf{t}) + \gamma w_{x}(\mathbf{t})\|_{L^{2}(0,\ell)}^{2} \right],$$

$$(3.11)$$

and it is decreasing function of the time \mathbf{t} . In particular, we have

(3.12)
$$\frac{dE_0(\mathbf{t})}{d\mathbf{t}} = -\|u_{\mathbf{t}}(\mathbf{t})\|_{L^2(0,\ell)}^2 - \|v_{\mathbf{t}}(\mathbf{t})\|_{L^2(0,\ell)}^2 - \|w_{\mathbf{t}}(\mathbf{t})\|_{L^2(0,\ell)}^2.$$

Proposition 3.2. [5, Theorem 2.3.1] The operator \mathcal{A} generates a C_0 -semigroup of contractions in the Hilbert space \mathcal{H} . Hence the problem (3.10) admits a unique solution

$$(u(x, \mathbf{t}), v(x, \mathbf{t}), w(x, \mathbf{t}), \varphi(x, \mathbf{t}, \xi), \phi(x, \mathbf{t}, \xi), \psi(x, \mathbf{t}, \xi))$$

such that if $(u_0, v_0, w_0, u_1, v_1, w_1, 0, 0, 0) \in \mathcal{D}(\mathcal{A})$, then the solution $(u, v, w, \varphi, \phi, \psi)$ of (3.10) satisfies the following regularity property:

$$\mathcal{U} = (u, v, w, u_{\mathbf{t}}, v_{\mathbf{t}}, w_{\mathbf{t}}, \varphi, \phi, \psi) \in C([0, +\infty), \mathcal{D}(\mathcal{A})) \cap C^{1}([0, +\infty), \mathcal{H}),$$

and when $(u_0, v_0, w_0, u_1, v_1, w_1, 0, 0, 0) \in \mathcal{H}$, then

$$\mathcal{U} = (u, v, w, u_{\mathbf{t}}, v_{\mathbf{t}}, w_{\mathbf{t}}, \varphi, \phi, \psi) \in C([0, +\infty), \mathcal{H}).$$

The energy of the system (3.10), defined by

(3.13)
$$E(\mathbf{t}) := \frac{1}{2} \| (u, v, w, u_{\mathbf{t}}, v_{\mathbf{t}}, w_{\mathbf{t}}, \varphi, \phi, \psi) \|_{\mathcal{H}}^{2},$$

is decreasing in time t. In particular, we have

(3.14)
$$\begin{aligned} \frac{dE(\mathbf{t})}{d\mathbf{t}} &= -\mathfrak{C} \int_{\mathbb{R}} (|\xi|^2 + \eta) \|\varphi(\mathbf{t},\,\xi)\|_{L^2(0,\,\ell)}^2 d\xi - \mathfrak{C} \int_{\mathbb{R}} (|\xi|^2 + \eta) \|\phi(\mathbf{t},\,\xi)\|_{L^2(0,\,\ell)}^2 d\xi \\ &- \mathfrak{C} \int_{\mathbb{R}} (|\xi|^2 + \eta) \|\psi(\mathbf{t},\,\xi)\|_{L^2(0,\,\ell)}^2 d\xi. \end{aligned}$$

4. Non-uniform stabilization

4.1. Polynomial stability for $\eta = 0$. Assume that $\eta = 0$, then the operator \mathcal{A} is not onto and consequently $0 \notin \rho(\mathcal{A})$ the resolvent set of \mathcal{A} .

Since the embedding $H_0^1(0, \ell) \times H_0^1(0, \ell) \times H_0^2(0, \ell) \subset L^2(0, \ell) \times L^2(0, \ell) \times L^2(0, \ell)$ is compact and the only solution of the following problem:

$$\begin{cases} u_{\mathbf{t}\mathbf{t}} - \vartheta u_{xx} - k(-u+v+\gamma w_x) = 0, \quad (0,\,\ell) \times (0,\,+\infty), \\ v_{\mathbf{t}\mathbf{t}} - \chi v_{xx} + k(-u+v+\gamma w_x) = 0, \quad (0,\,\ell) \times (0,\,+\infty), \\ w_{\mathbf{t}\mathbf{t}} + \zeta w_{xxxx} - k(-u+v+\gamma w_x)_x = 0, \quad (0,\,\ell) \times (0,\,+\infty), \\ u_{\mathbf{t}}(x,\mathbf{t}) = 0, v_{\mathbf{t}}(x,\mathbf{t}) = 0, w_{\mathbf{t}}(x,\mathbf{t}) = 0, \quad (0,\,\ell) \times (0,\,+\infty), \\ u(0,\,\mathbf{t}) = u(\ell,\,\mathbf{t}) = 0, \, v(0,\,\mathbf{t}) = v(\ell,\,\mathbf{t}) = 0, \quad (0,\,+\infty), \\ w(0,\,\mathbf{t}) = w(\ell,\,\mathbf{t}) = w_x(0,\,\mathbf{t}) = w_x(\ell,\,\mathbf{t}) = 0, \quad (0,\,+\infty), \\ u(x,\,0) = u_0(x), \quad u_{\mathbf{t}}(x,\,0) = u_1(x), \quad (0,\,\ell), \\ v(x,\,0) = v_0(x), \quad w_{\mathbf{t}}(x,\,0) = v_1(x), \quad (0,\,\ell), \\ w(x,\,0) = w_0(x), \quad w_{\mathbf{t}}(x,\,0) = w_1(x), \quad (0,\,\ell), \end{cases}$$

is the trivial one.

(4.1)

Then according to [5, Sect. 2.4] the semigroup $(e^{\mathbf{t}\mathcal{A}})_{t\geq 0}$ is strongly stable, i.e.,

$$\|e^{\mathbf{t}\mathcal{A}}\mathcal{U}_0\|_{\mathcal{H}} \longrightarrow 0 \quad \text{as} \quad \mathbf{t} \to +\infty$$

for all initial data $\mathcal{U}_0 \in \mathcal{H}$.

To determine the rate of stability in this case, we need the following result [7, Theorem 8.4]:

Theorem 4.1 ([7]). Let $\mathcal{T}(t)$ be a bounded C_0 -semigroup on a Hilbert space X with generator H. Assume that $\sigma(H) \cap i\mathbb{R} = \{0\}$ and that there exist $\beta \ge 1, \gamma > 0$ such that

$$\left\| (i\lambda I - H)^{-1} \right\|_{\mathcal{L}(X)} \le \begin{cases} O(|\lambda|^{-\beta}), \, \lambda \to 0, \\ O(|\lambda|^{\gamma}), \, \lambda \to \infty. \end{cases}$$

Then, there exists a constant C > 0 such that

$$\|\mathcal{T}(t)z\|_{X} \leq \frac{C}{\mathbf{t}^{\frac{1}{\max(\beta,\gamma)}}} \|z\|_{\mathcal{D}(H)}, \ \forall t > 0, z \in \mathcal{D}(H) \cap R(H),$$

where $\mathcal{D}(H)$ is the domain of H and R(H) is the range of H.

Based on Theorem 4.1, a simple adaptation of the proofs of [8, Theorems 5.8 and 5.9] (see also [4, 5]) lead to the following stability result:

Proposition 4.2. If $\eta = 0$, then there exists a C > 0 such that

(4.2)
$$\left\| e^{\mathbf{t}\mathcal{A}}U_0 \right\|_{\mathcal{H}} \leq \frac{C}{\sqrt{\mathbf{t}}} \left\| U_0 \right\|_{\mathcal{D}(\mathcal{A})}, \, \forall \mathbf{t} > 0, \, U_0 \in \mathcal{D}(\mathcal{A}) \cap R(\mathcal{A}),$$

where $R(\mathcal{A})$ is the range of \mathcal{A} .

4.2. Polynomial stability for $\eta > 0$. In this case there exists δ , C > 0, such that the auxiliary dissipative operator satisfies the following relation:

$$\left\| e^{\mathbf{t}\mathcal{A}_0} \right\|_{\mathcal{L}(\mathcal{H}_0)} \leq C \, e^{-\delta \mathbf{t}}, \quad \forall \ \mathbf{t} \geq 0.$$

Applying [5, Corollary 2.6.1] hence we obtain the following polynomial decay result for the systems (1.3) and (2.10):

Proposition 4.3. If $\eta > 0$, then the semigroup $(e^{tA})_{t\geq 0}$ is polynomially stable, namely there exists a constant C > 0 such that

$$\begin{split} \|e^{\mathbf{t}\mathcal{A}}(u_0, v_0, w_0, u_1, v_1, w_1, \varphi_0, \phi_0, \psi_0)\|_{\mathcal{H}} \\ &\leq \frac{C}{(1+\mathbf{t})^{\frac{1}{1-\alpha}}} \|(u_0, v_0, w_0, u_1, v_1, w_1, \varphi_0, \phi_0, \psi_0)\|_{\mathcal{D}(\mathcal{A})}, \quad \forall \ \mathbf{t} \geq 0, \end{split}$$

for all initial data $(u_0, v_0, w_0, u_1, v_1, w_1, \varphi_0, \phi_0, \psi_0) \in \mathcal{D}(\mathcal{A})$. In particular, the energy of the strong solution of (1.3) and (2.10) satisfies the following estimate:

$$E(\mathbf{t}) \le \frac{C}{(1+\mathbf{t})^{\frac{2}{1-\alpha}}} \| (u_0, v_0, w_0, u_1, v_1, w_1, 0, 0, 0) \|_{\mathcal{D}(\mathcal{A})}^2, \quad \forall \, \mathbf{t} > 0.$$

5. Numerical study

In this section we will verify numerically the polynomial rate of decay obtained in the previous section.

5.1. Linear equations of Motion. First, we approximate the longitudinal and transverse displacement vector $[u, v, w]^{\top}$ in space using an energy-conservative finite difference method. For $J \in \mathbb{N}$ and $\delta x = \ell/J$, we define x_j , with $j = 1, \ldots, J$, as a uniform discretization of the interval $(0, \ell)$, obtaining a vector $\mathbf{U} = [\mathbf{u}(t), \mathbf{v}(t), \mathbf{w}(t)]^{\top}$ approximation of $[u, v, w]^{\top}$ in \mathbb{R}^{3J} . We obtain the linear equation of motion

(5.1)
$$\mathbf{M}\ddot{\mathbf{U}}(t) + \mathbf{K}\mathbf{U}(t) + \mathbf{C}\overset{\alpha,\eta}{\mathbf{U}}(t) = 0$$

where $\overset{\alpha,\eta}{\mathbf{U}}(t) = D^{\alpha,\eta}\mathbf{U}(t)$ is the generalized Caputo fractional derivative defined in (2.2). In addition, and according to a discretization consistent with (1.3), we choose $\mathbf{M} = \mathbf{C} = \mathbf{I}_{3J}$ (identity matrix). The stiffness matrix is given by $\mathbf{K} = \mathbf{K}_{\text{stress}} + \mathbf{K}_{\text{coupling}}$, where

$$\mathbf{K}_{\text{stress}} = \begin{pmatrix} -\vartheta \mathbf{D}^2 & & \\ & -\chi \mathbf{D}^2 & \\ & & \zeta \mathbf{D}^4 \end{pmatrix}, \qquad \mathbf{K}_{\text{coupling}} = k \begin{pmatrix} \mathbf{I}_J & -\mathbf{I}_J & \gamma \mathbf{R} \\ -\mathbf{I}_J & \mathbf{I}_J & -\gamma \mathbf{R} \\ \gamma \mathbf{R}^T & -\gamma \mathbf{R}^T & -\gamma^2 \mathbf{D}^2. \end{pmatrix}$$

Here, \mathbf{D}^2 and \mathbf{D}^4 are the finite difference matrix approximation of ∂_{xx} and ∂_{xxxx} respectively with the boundary conditions $(1.3)_{7,8}$, and \mathbf{R} is the upper triangular matrix, obtained from the Cholesky decomposition $\mathbf{D}^2 = -\mathbf{R}^T \mathbf{R}$. Using uniform finite differences we have

According to the boundary conditions $(1.3)_{7,8}$, we seek the solution of Finite Difference in

$$X_J := \left\{ \mathbf{U} = [\mathbf{u}(t), \mathbf{v}(t), \mathbf{w}(t)]^\top \in \mathbb{R}^{3J} \right\},\$$

considering $\mathbf{u}_0 = \mathbf{u}_{J+1} = \mathbf{v}_0 = \mathbf{v}_{J+1} = \mathbf{w}_0 = \mathbf{w}_{J+1} = \mathbf{v}_{-1} = \mathbf{v}_{J+2} = 0.$

5.2. Time discretization. In order to preserve the energy with a second order scheme in time, we choose a β -Newmark scheme for w. The method consists of updating the displacement, velocity and acceleration vectors at the current time $t^n = n\delta t$ to the time $t^{n+1} = (n+1)\delta t$, a small time interval δt later. The Newmark algorithm [38] is based on a set of two relations expressing the forward displacement \mathbf{U}^{n+1} and velocity $\dot{\mathbf{U}}^{n+1}$ in terms of their current values and the forward and current values of the acceleration:

(5.2)
$$\dot{\mathbf{U}}^{n+1} = \dot{\mathbf{U}}^n + (1 - \widetilde{\gamma})\delta t \, \ddot{\mathbf{U}}^n + \widetilde{\gamma} \delta t \, \ddot{\mathbf{U}}^{n+1},$$

(5.3)
$$\mathbf{U}^{n+1} = \mathbf{U}^n + \delta t \dot{\mathbf{U}}^n + \left(\frac{1}{2} - \widetilde{\beta}\right) \delta t^2 \, \ddot{\mathbf{U}}^n + \widetilde{\beta} \delta t^2 \, \ddot{\mathbf{U}}^{n+1},$$

where $\tilde{\beta}$ and $\tilde{\gamma}$ are parameters of the methods that will be fixed later. Replacing (5.2)–(5.3) in the equation of motion (5.1), we obtain

(5.4)
$$\left(\mathbf{M} + \widetilde{\beta}\delta t^{2} \mathbf{K}\right) \ddot{\mathbf{U}}^{n+1} + \mathbf{C} \overset{\alpha,\eta}{\mathbf{U}^{n+1}} = -\mathbf{K} \left(\mathbf{U}^{n} + \delta t \dot{\mathbf{U}}^{n} + \left(\frac{1}{2} - \widetilde{\beta}\right) \delta t^{2} \ddot{\mathbf{U}}^{n}\right).$$

5.3. Approximation of fractional derivatives. At this state we have two possibilities: (1) approximate the fractional derivative using a classical finite difference method such as the truncation of the Grünwald—Letnikov derivative [19], which in turn is an equivalent form of the Riemann–Liouville derivative; or, (2) propose a conservative numerical scheme for the Mbodje augmented model (2.10).

5.3.1. Approximation using a truncation of the Grünwald--Letnikov derivative. In the case of the classical Grünwald-Letnikov derivative we consider a fractional trapezoidal formula [26], in order to approximate

(5.5)
$$\left[{}_{C}D_{0,t}^{\alpha,\eta}f \right]_{t=t_{n}} \approx \sum_{k=0}^{n} a_{k,n}e^{-\eta\delta t(n-k)}f'(t_{k}),$$

where

(5.6)
$$a_{k,n} = \frac{\delta t^{1-\alpha}}{\Gamma(3-\alpha)} \begin{cases} (n-1)^{2-\alpha} - (n-2+\alpha)n^{1-\alpha}, & k = 0, \\ (n-k+1)^{2-\alpha} + (n-1-k)^{2-\alpha} - 2(n-k)^{2-\alpha}, & 1 \le k \le n-1 \\ 1, & k = n. \end{cases}$$

Replacing the term $\overset{\alpha,\eta}{\mathbf{U}}_{n+1}^{n+1}$ by the numerical approximation $\left[{}_{C}D_{0,t}^{\alpha,\eta}f\right]_{t=t_{n}}$ given by (5.5)–(5.6), then the equation (5.4) becomes

(5.7)
$$\left(\mathbf{M} + \widetilde{\beta}\delta t^{2}\mathbf{K} + \frac{\widetilde{\gamma}\delta t^{2-\alpha}}{\Gamma(3-\alpha)}\mathbf{C}\right)\ddot{\mathbf{U}}^{n+1} = -\mathbf{K}\left(\mathbf{U}^{n} + \delta t\dot{\mathbf{U}}^{n} + \left(\frac{1}{2} - \widetilde{\beta}\right)\delta t^{2}\ddot{\mathbf{U}}^{n}\right) - \frac{\delta t^{1-\alpha}}{\Gamma(3-\alpha)}\mathbf{C}\left(\dot{\mathbf{U}}^{n} + (1-\widetilde{\gamma})\,\delta t\ddot{\mathbf{U}}^{n} + \sum_{k=0}^{n}a_{k,n+1}\dot{\mathbf{U}}^{k}\right).$$

5.3.2. Approximation using the Mbodje augmented model. On the other hand, in case of the Mbodje augmented model (2.10) we propose the following conservative numerical scheme:

(5.8)
$$\begin{cases} \left(\mathbf{M} + \widetilde{\beta}\delta t^{2}\,\mathbf{K}\right)\ddot{\mathbf{U}}^{n+1} + \mathfrak{C}\sum_{\ell=1}^{M}\mu_{\ell}\boldsymbol{\Phi}_{\ell}^{n+1} = -\mathbf{K}\left(\mathbf{U}^{n} + \delta t\dot{\mathbf{U}}^{n} + \left(\frac{1}{2} - \widetilde{\beta}\right)\delta t^{2}\ddot{\mathbf{U}}^{n}\right),\\ \mathbf{\Phi}_{\ell}^{n+1} = \mathbf{\Phi}_{\ell}^{n} - \delta t\left(\xi_{\ell}^{2} + \eta\right)\mathbf{\Phi}_{\ell}^{n+\frac{1}{2}} + \delta t\mu_{\ell}\dot{\mathbf{U}}^{n+\frac{1}{2}},\end{cases}$$

where $\mathbf{\Phi}_{\ell}^{n+\frac{1}{2}} = \frac{\mathbf{\Phi}_{\ell}^{n} + \mathbf{\Phi}_{\ell}^{n+1}}{2}$, with $\mathbf{\Phi}_{\ell}^{n}$ denoting the approximation of $[\phi, \varphi, \psi]^{\top} \in \mathbb{R}^{3J}$ evaluated in $\xi_{\ell} := \ell \delta \xi$ and $t = n \delta t$, for $\ell = 1, \ldots, M$, $n = 1, \ldots, N$ and a fixed $\delta \xi > 0$. Using (5.2) again and replacing in (5.8)₂, we can rewrite (5.8) in the following more explicit and computable way:

(5.9)
$$\begin{cases} \left(\mathbf{M} + \widetilde{\gamma}\delta t \, \mathbf{C}_{augm} + \widetilde{\beta}\delta t^2 \, \mathbf{K}\right) \ddot{\mathbf{U}}^{n+1} = -\mathfrak{C} \sum_{\ell=1}^{M} \widetilde{\mu}_{\ell} \boldsymbol{\Phi}_{\ell}^{n} - \mathbf{C}_{augm} \left(\dot{\mathbf{U}}^{n} + (1-\widetilde{\gamma})\,\delta t \ddot{\mathbf{U}}^{n}\right) \\ -\mathbf{K} \left(\mathbf{U}^{n} + \delta t \dot{\mathbf{U}}^{n} + \left(\frac{1}{2} - \widetilde{\beta}\right)\delta t^2 \ddot{\mathbf{U}}^{n}\right), \\ \mathbf{\Phi}_{\ell}^{n+1} = \frac{2 - \delta t \left(\xi_{\ell}^{2} + \eta\right)}{2 + \delta t \left(\xi_{\ell}^{2} + \eta\right)} \boldsymbol{\Phi}_{\ell}^{n} + \frac{2\delta t \mu_{\ell}}{2 + \delta t \left(\xi_{\ell}^{2} + \eta\right)} \dot{\mathbf{U}}^{n+\frac{1}{2}}, \end{cases}$$

where $\mu_{\ell} = \xi_{\ell}^{(2\alpha-1)/2}$, $\tilde{\mu}_{\ell} = \frac{2 - \delta t \left(\xi_{\ell}^2 + \eta\right)}{2 + \delta t \left(\xi_{\ell}^2 + \eta\right)} \mu_{\ell}$ and $\mathbf{C}_{augm} = \delta t \mathfrak{C} \left(\sum_{\ell=1}^{M} \frac{2\tilde{\mu}_{\ell}^2}{2 + \delta t \left(\xi_{\ell}^2 + \eta\right)}\right) \mathbf{I}_{3J}$. Although the two approximations are consistently valid from the point of view of finite differences,

we will opt for the second approximation because it conserves the energy in the non-dissipative case, and its decrease is consistent with the continuous case. It also involves the calculation of Φ which is inherent to the definition of energy itself.

5.4. Decay of the discrete energy using the Mbodje augmented model. Evaluating $(5.8)_1$ in $t = t_{n+\frac{1}{2}}$, multiplying by $\ddot{\mathbf{U}}^{n+\frac{1}{2}}$, and summing $(5.8)_2$ multiplied by $\mathfrak{C}\Phi_{\ell}^{n+\frac{1}{2}}$, we obtain that

$$\begin{bmatrix} E_{\Delta}^{\text{Mbodje}} \end{bmatrix}_{n}^{n+1} := \begin{bmatrix} \frac{1}{2} \dot{\mathbf{U}}^{T} \mathbf{M} \dot{\mathbf{U}} + \frac{1}{2} \mathbf{U}^{T} \mathbf{K} \mathbf{U} + \frac{\mathfrak{C}}{2} \sum_{\ell=1}^{M} \mu_{\ell} |\boldsymbol{\Phi}_{\ell}|^{2} \end{bmatrix}_{n}^{n+1}$$

$$= \begin{bmatrix} -\frac{\vartheta}{2} \dot{\mathbf{u}}^{T} \mathbf{D}^{2} \dot{\mathbf{u}} - \frac{\chi}{2} \dot{\mathbf{v}}^{T} \mathbf{D}^{2} \dot{\mathbf{v}} + \frac{\zeta}{2} \dot{\mathbf{w}}^{T} \mathbf{D}^{4} \dot{\mathbf{w}} + \frac{k}{2} \| - \mathbf{u} + \mathbf{v} + \gamma \mathbf{R} \mathbf{w} \|^{2}$$

$$+ \frac{\mathfrak{C}}{2} \sum_{\ell=1}^{M} \mu_{\ell} \left(|\phi_{\ell}|^{2} + |\varphi_{\ell}|^{2} + |\psi_{\ell}|^{2} \right) \end{bmatrix}_{n}^{n+1}$$

$$(5.10) \qquad = -\mathfrak{C} \sum_{\ell=1}^{M} \left(\xi_{\ell}^{2} + \eta \right) \left(\left| \phi_{\ell}^{n+\frac{1}{2}} \right|^{2} + \left| \varphi_{\ell}^{n+\frac{1}{2}} \right|^{2} + \left| \psi_{\ell}^{n+\frac{1}{2}} \right|^{2} \right)$$

which is consistent with the estimates (3.13)–(3.14), and constitutes a correct approximation of the energy and its decreasing behavior.

Remark 5.1. In the case of an approximation using a truncation of the Grünwald—Letnikov derivative, we have the following estimate of the energy. Evaluating (5.1) in $t = t_{n+\frac{1}{2}}$ and



FIGURE 1. Comparison between the Grünwald–Letnikov derivative approximation and the discretized Mbodje augmented model. At the top left: $E_{\Delta}^{GL}(t) = \frac{1}{2} \dot{\mathbf{U}}^T \mathbf{M} \dot{\mathbf{U}} + \frac{1}{2} \mathbf{U}^T \mathbf{K} \mathbf{U}$; At the top right: $E_{\Delta}^{GL,\Phi}(t) = E_{\Delta}^{GL}(t) + \frac{\mathfrak{C}}{2} \sum_{\ell=1}^{M} \mu_{\ell} |\Phi_{\ell}|^2$; At the bottom: $E_{\Delta}^{Mbodje}(t)$ defined in (5.10).

multiplying by $\ddot{\mathbf{U}}^{n+\frac{1}{2}}$ we obtain

$$\begin{split} \left[E_{\Delta}^{GL}\right]_{n}^{n+1} &:= \left[\frac{1}{2}\dot{\mathbf{U}}^{T}\mathbf{M}\dot{\mathbf{U}} + \frac{1}{2}\mathbf{U}^{T}\mathbf{K}\mathbf{U}\right]_{n}^{n+1} \\ &= \left[-\frac{\vartheta}{2}\dot{\mathbf{u}}^{T}\mathbf{D}^{2}\dot{\mathbf{u}} - \frac{\chi}{2}\dot{\mathbf{v}}^{T}\mathbf{D}^{2}\dot{\mathbf{v}} + \frac{\zeta}{2}\dot{\mathbf{w}}^{T}\mathbf{D}^{4}\dot{\mathbf{w}} + \frac{k}{2}\|-\mathbf{u}+\mathbf{v}+\gamma\mathbf{R}\mathbf{w}\|^{2}\right]_{n}^{n+1} \\ &= -\frac{\delta t^{2-\alpha}}{\Gamma(3-\alpha)}\left(\dot{\mathbf{U}}^{n+\frac{1}{2}}\right)^{T}\mathbf{C}\dot{\mathbf{U}}^{n+\frac{1}{2}} - (2^{1-\alpha}-1)\frac{\delta t^{2-\alpha}}{\Gamma(3-\alpha)}\left(\dot{\mathbf{U}}^{n+\frac{1}{2}}\right)^{T}\mathbf{C}\dot{\mathbf{U}}^{n} \\ (5.11) \qquad -\sum_{k=0}^{n-1}a_{k,n+\frac{1}{2}}\left(\dot{\mathbf{U}}^{n+\frac{1}{2}}\right)^{T}\mathbf{C}\dot{\mathbf{U}}^{k} \end{split}$$

where $a_{k,n+\frac{1}{2}} = \frac{a_{k,n+1} + a_{k,n}}{2}$. The first term on the right hand side of (5.11) is obviously strictly negative. However, the sign of the second and third terms in (5.11) cannot be guaranteed. In fact, it is verified numerically that the energy discretized in this way is not necessarily strictly decreasing, unless the solution is sufficiently smooth and delta t is small enough. Figure 1 shows a comparison of both models. It is proved that the calculation of energy using the classical

Grünwald—Letnikov derivatives fails, either by using the derivative defined in (5.11), (in top left graph), as well as by artificially adding the calculation of $\mathbf{\Phi}$ as seen in the top right graph. Therefore, the only acceptable method for all our calculations should be the discretization of the augmented Mbodje model (bottom graph).

5.5. Numerical examples. We consider here the value of the parameters $\vartheta = \chi = \zeta = k = \gamma = 1$, a beam of length $\ell = 1$, and the initial conditions

(5.12)
$$\begin{cases} u(x,0) = x \left(\frac{\ell^3}{8} - \left|x - \frac{\ell}{2}\right|^3\right), & u_t(x,0) = 0, \\ v(x,0) = 0, & v_t(x,0) = 0, \\ w(x,0) = x \left(x - \frac{\ell}{2}\right) \left(\frac{\ell^3}{8} - \left|x - \frac{\ell}{2}\right|^3\right), & w_t(x,0) = 0. \end{cases}$$

We observe that $u(0,t) = u(\ell,t) = 0$ and $u(\cdot,0)$ is of class $\mathcal{C}^1(0,\ell)$ but its second derivative has a discontinuity in $x = \ell/2$, and in turn, $w(0,t) = w(\ell,t) = 0$ with $w(\cdot,0)$ of class $\mathcal{C}^3(0,\ell)$ and its fourth derivative has a discontinuity at $x = \ell/2$. Therefore, we have just $\mathcal{U}_0 \in \mathcal{D}(A)$, but no greater regularity than that.

In Figure 2, we consider a fractional derivative with $\alpha = 0.5$, and we do a simulation of the numerical scheme (5.2), (5.3), (5.7) with J = 500, and $\delta = T/N_t$ with T = 100 and $N_t = 1000$. The graphs in (A) of Figure 2 shows the state of the instant of the deformations for t = 10. The beam is represented by the thick yellow graph and its deformation corresponds to the transversal displacements. The longitudinal displacements of u(x, 10) and v(x, 10), are represented by the blue and red lines respectively. The graph (B) of the figure 2 shows the evolution of the transverse displacements w in term of space and time. Graphs (C) and (D) of the same figure show the longitudinal displacements u and v (respectively), also as a function of space and time. In this set of figures, the decay of the quantities, and therefore of the energy, is visually observed due to the dispative terms of the fractional derivative type.

On the other hand, we show in figure 3 the graphs of the numerical variables that approximate the auxiliary functions $\phi(x, t; \xi)$, $\varphi(x, t; \xi)$, and $\psi(x, t; \xi)$ of the augmented model (2.10), for the instant t = 2 of the simulations with the same parameters and data as the example in figure 2. We numerically observe a decay of these functions with respect to the variable ξ , although not so evident in the case of the variable ψ associated with the term of the fractional derivative of the transverse displacement (when truncating at $\xi = 10$). In this sense, we must be careful with the numerical truncation of the variable $\xi \in (0, +\infty)$. For this reason we will consider a truncation at $\xi = 10,000$ for the examples that will follow, and that will show the polynomial decay of energy for different values of α and η .

Finally, in order to numerically appreciate the polynomial decay of the energy, we do the simulation with the same parameters $\vartheta = \chi = \zeta = k = \gamma = 1$, a beam of length $\ell = 1$, and the initial

0.02

0.01 w(x,t)

-0.01

0

0

0.5

х

1 0

(B) Transversal deformations w



(A) State of the deformations at instant t = 10





(D) Longitudinal deformations v

FIGURE 2. Simulation of the displacements. On the left: the displacements at t = 2. On the right: the evolution in time of the transverse displacements.

conditions (5.12), but for a longer time T = 10000, and a time step $\delta t = 0.1$. Likewise, in order to obtain better precision in the results, we truncate the variable ξ of the auxiliary functions at $\xi = 10000.$

In Figure 4, various plots depict energy on a log-log scale, where curves of the form $C(t+1)^p$ exhibit nearly linear behavior over sufficiently long times.

The graph in Figure 4(A) illustrates energy curves for 9 different values of η and 5 values of α , totaling 45 combinations. Similar curves with the same α are grouped into 5 distinct colors. While there are variations in slopes p and constants C, graphical observation reveals a close relationship between these parameters, particularly in their bounded ranges. The numerical estimation of C and p using a least squares method is presented in Table 1.

10

8

6

2



FIGURE 3. Behavior of the auxiliary variables ϕ , φ and ψ of the augmented model (2.10)

Additionally, from the graph, it is conjectured that energy decays more rapidly to zero for $\alpha = 0.5$ (or similar values) and certain $\eta \neq 0$. Conversely, $\alpha = 0.25$ exhibits intriguing behavior, accentuating differences across various η values. Thus, the curves corresponding to $\alpha = 0.25$ are separately plotted in Figure 4(B). Here, a pattern emerges with curves lying between two quasi-linear lines with distinct slopes. The steepest negative slope corresponds to $\eta = 0.3$, while the smallest absolute slope corresponds to $\eta = 0$, with intermediate values approaching the former asymptotically.

Figures 4(C) and 4(D) compare energy curves with their respective asymptotic forms resembling straight lines in a log-log graph. These asymptotic curves, of the form $C(1 + t)^p$, are derived through least squares, effectively capturing the data's behavior. In Figure 4(C), distinct η values for the same α exhibit notably different slopes (i.e., decay rates p). Conversely, in Figure 4(D), variations in the multiplicative constant C outweigh changes in the decay rate p across different α values.



(A) Energies for different parameters α and η regrouped in terms of the values of α .



(C) Comparison between energy curves with a same $\alpha = 0.25$ as a function of time, and their asymptotes $C(t+1)^p$



(B) Zoom: Energies for $\alpha = 0.25$,

and different values of η



(D) Comparison between energy curves with a same $\eta = 0.3$ as a function of time, and their asymptotes $C(t+1)^p$

FIGURE 4. Energy with different values of η and α in loglog scale.

DISCUSSION AND CONCLUSIONS

We have demonstrated the existence, uniqueness, and stability of a Rao–Nakra model with dissipative terms of the generalized Caputo fractional derivative type within the domain, building upon ideas from a proven case for a broader class of evolution systems [4]. While focusing on the Rao–Nakra beam model, our treatment of the fractional term through an input–output relationship with a diffusion equation [34] has illuminated a numerical method of finite-conservative differences type that ensures stable energy decay, unlike standard methods for approximating fractional derivatives. This enabled us to numerically simulate energy decay, confirming its polynomial nature.

				α		
$C(t+1)^p$		0.01	0.25	0.50	0.75	0.99
η	0	C = 0.0300	0.0250	0.0600	0.0186	0.1521
		p = -0.6465	-0.4136	-0.7497	-0.9147	-0.7968
	1.10^{-4}	C = 0.0869	0.0408	0.0250	0.0198	0.1521
		p = -0.8136	-0.8630	-0.9773	-0.9255	-0.7968
	3.10^{-4}	C = 0.1089	0.4175	0.0338	0.0198	0.1521
		p = -0.8600	-1.2750	-1.0469	0.9266	-0.7968
	1.10^{-3}	C = 0.0894	0.0832	0.0338	0.0198	0.1521
		p = -0.8413	-1.1394	-0.9985	-0.9236	-0.7968
	3.10^{-3}	C = 0.0822	0.0258	0.0159	0.0190	0.1522
		p = -0.8317	-1.0075	-0.9688	-0.9217	-0.7968
	1.10^{-2}	C = 0.0803	0.0170	0.0144	0.0188	0.1525
		p = -0.8290	-0.9586	-0.9569	-0.9207	-0.7967
	3.10^{-2}	C = 0.0801	0.0155	0.0141	0.0188	0.1533
		p = -0.8283	-0.9469	-0.9534	-0.9201	-0.7964
	$1 10^{-1}$	C = 0.0806	0.0152	0.0140	0.0190	0.1562
	1.10	p = -0.8279	-0.9431	-0.9514	-0.9190	-0.7955
	$3 10^{-1}$	C = 0.0823	0.0154	$0.01\overline{43}$	0.0196	0.1653
	0.10	p = -0.8269	-0.9404	-0.9485	-0.9161	-0.7925

TABLE 1. Constant C and rate p of polynomial decay $C(t+1)^p$ for the energy by a least square method, for different values of the fractional derivative parameters α and η .

One key observation from the data and interpretation of Figure 4 and Table 1 is the clear asymptotic polynomial decay behavior of energy. We note that the decay rate is notably greater in absolute value when $\eta \neq 0$, and lower when $\eta = 0$, validating the theoretical propositions (Proposition 4.3 and Theorem 4.2). However, we acknowledge discrepancies between the theoretically derived rates and the numerically estimated ones. We attribute this to factors such as the discretization step sizes δt and δx , the order of convergence of the method introducing additional numerical dissipation, and the limited simulation times for observing true theoretical decay rates.

Future work should delve into numerical methods enhancing convergence order to obtain numerical values closer to theoretical ones. Additionally, exploring inverse or control problems, such as identifying parameter pairs (α, η) for quicker energy decay, merits attention. Notably, from Table 1, we observe a minimum in the rate p near $(\alpha, \eta) = (0.5; 0.0003)$. An inverse problem could involve minimizing this slope with respect to (α, η) . Moreover, considering the importance of the constant C, as evidenced by the curve nearest to zero at t = 10,000 corresponding to $(\alpha, \eta) = (0.5; 0.3)$ in Figure 4(A), remains intriguing. However, beyond t = 10,000[s], maintaining a constant p to ensure optimal decay is unclear. While minimizing energy decay speed is of interest, defining a suitable cost function characterizing this decay speed requires further exploration. Thus, identifying the appropriate cost function to minimize for solving the inverse problem remains an open task.

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