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Structure-preserving coupling and decoupling of port-Hamiltonian systems

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ABSTRACT

The port-Hamiltonian framework is a structure-preserving modeling approach that preserves key physical properties such as energy conservation and dissipation. When subsystems are modeled as port-Hamiltonian systems (pHS) with linearly related inputs and outputs, their interconnection remains port-Hamiltonian. This paper introduces a systematic method for transforming coupled port Hamiltonian ordinary differential equations systems (pHODE) into a single monolithic formulation, and for decomposing a monolithic system into weakly coupled subsystems. The monolithic representation ensures stability and structural integrity, whereas the decoupled form enables efficient distributed simulation via operator splitting or dynamic iteration.

1. Introduction

In today's rapidly evolving technological landscape, the ability to model and analyze complex systems is more crucial than ever. Port-Hamiltonian systems (pHS) combine energy-based modeling with network interconnection structure to provide unparalleled insight into multi-component physical systems. First, pHS exhibit a Hamiltonian structure, i.e., pHS provide an *energy-based modeling approach*, essential for understanding the dynamics of modern systems (see related papers [1–4] and book [5]). Next, they allow for *dissipation* in the system. This accurately models energy losses, ensuring realistic simulations and designs. They have *ports* that allow for input/output connections to the outside world, i.e. this approach facilitates energy exchange with the external environment. Finally, pHS have an *interconnection property*: multiple subsystems can be coupled to a monolithic joint pH system, preserving the overall energy structure and increasing reliability of the system.

In this work, we investigate (a) how coupled port-Hamiltonian ordinary differential equation (pHODE) systems can be formulated as a single joint pHODE and (b) vice-versa, how a single joint pHODE system can be decoupled into subsystems of pHODE, and generalize some ideas mentioned in [6]. While (a) allows for monolithic simulation with excellent stability properties, (b) is the starting point for an efficient distributed simulation via operator splitting and/or dynamic iteration.

The paper is organized as follows: after an introduction, in Section 2 we define the setting of port-Hamiltonian ODE systems. Section 3 deals with the coupling of pHODE subsystems into a joint pHODE, and Section 4 discusses the decoupling of a single pHODE system into pHODE subsystems. Section 5 discusses briefly the advantages and disadvantages of both approaches. We conclude with final remarks in Section 6.

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2. Port-Hamiltonian ODE systems

Following the lines of [7, Definition 1], we define the general setting of port-Hamiltonian ordinary differential equations (pHODEs). Consider a time interval $I \subset \mathbb{R}$, a state space $\mathcal{X} \subset \mathbb{R}^n$. A pHODE is a (possibly implicit) system of ordinary differential equations (ODEs) of the form:

$$E(x)\dot{x} = (J(x) - R(x))z(x) + (B(x) - P(x))u, \quad (1a)$$

$$y = (B(x) + P(x))^\top z(x) + (S(x) - N(x))u, \quad (1b)$$

associated with a Hamiltonian function $H \in C^1(\mathcal{X}, \mathbb{R})$, where $x(t) \in \mathcal{X}$ is the state, $u(t), y(t) \in \mathbb{R}^m$ are the input and output functions, $r, z \in \mathcal{X}(S, \mathbb{R})$ are the *time-flow and effort functions*, $E \in C(\mathcal{X}, \mathbb{R}^{n \times n})$ is the *regular flow matrix*, $J, R \in \mathcal{X}(S, \mathbb{R}^{n \times n})$ are the *structure and dissipation matrices*, $B, P \in C(\mathcal{X}, \mathbb{R}^{n \times p})$ are the *port matrices* and $S, N \in C(\mathcal{X}, \mathbb{R}^{p \times p})$ are the *feed-through matrices*. Furthermore, the following property must hold:

$$\Gamma := \begin{bmatrix} J & B \\ -B^\top & N \end{bmatrix} = -\Gamma^\top, \quad W := \begin{bmatrix} R & P \\ P^\top & S \end{bmatrix} \geq 0, \quad (2)$$

i.e. Γ is a skew-symmetric matrix and W is a positive semidefinite matrix. The gradient of the Hamiltonian function satisfies $\nabla H = E(x)^\top z$ and $\partial_t H = z^\top r$ hold pointwise. The key property of such pHODE systems (1) is the following *dissipation equality*

$$\frac{d}{dt}H(x) = - \begin{pmatrix} z \\ u \end{pmatrix}^\top W \begin{pmatrix} z \\ u \end{pmatrix} + u^\top y, \quad (3)$$

for the Hamiltonian (we refer to [7, Section 2.2.1] for a proof).

Remark 1. In many engineering applications, the feed-through matrices S and N may vanish. Under the passivity requirement (2) this automatically implies for port matrix $P = 0$, because otherwise

$$W := \begin{bmatrix} R & P \\ P^\top & 0 \end{bmatrix}$$

will become indefinite for $P \neq 0$. Therefore we often work with the simplified formulation

$$E(x)\dot{x} = (J(x) - R(x))z(x) + B(x)u, \quad (4a)$$

$$y = B(x)^\top z(x). \quad (4b)$$

3. Coupling port-Hamiltonian ODE systems

To define *coupled port-Hamiltonian ODE systems*, we consider s subsystems of pHODEs (4)

$$E_i(x_i)\dot{x}_i(t) + r_i(x_i) = (J_i(x_i) - R_i(x_i))z_i(x_i) + \hat{B}_i(x_i)\hat{u}_i(t) + \bar{B}_i(x_i)\bar{u}_i(t),$$

$$\hat{y}_i(t) = \hat{B}_i(x_i)^\top z_i(x_i),$$

$$\bar{y}_i(t) = \bar{B}_i(x_i)^\top z_i(x_i),$$

$i = 1, \dots, s$, with $J_i = -J_i^\top$ and $R_i \geq 0$ and a Hamiltonian $H_i(x_i)$, which fulfills the compatibility conditions $E_i^\top z_i = \nabla_{x_i} H_i$, $i = 1, \dots, s$. Here, we distinguish inputs \hat{u}_i and outputs \hat{y}_i , which arise in the coupling between the subsystems, and inputs \bar{u}_i and outputs \bar{y}_i , which couple the individual system with the environment.

The subsystems are coupled via *external* inputs and outputs by

$$\begin{pmatrix} \hat{u}_1 \\ \vdots \\ \hat{u}_k \end{pmatrix} + \hat{C} \begin{pmatrix} \hat{y}_1 \\ \vdots \\ \hat{y}_k \end{pmatrix} = 0, \quad \hat{C} = -\hat{C}^\top. \quad (5)$$

These s systems can be condensed into a large monolithic port-Hamiltonian system (cf. [6] for the linear case)

$$E(x)\dot{x} = (J(x) - R(x))z(x) + \bar{B}(x)\bar{u}, \quad (6a)$$

$$\bar{y} = \bar{B}(x)^\top z(x), \quad (6b)$$

with $J(x) := \text{diag}(J_1(x_1), \dots, J_s(x_s)) - \hat{B}\hat{C}\hat{B}^\top$, $F = \text{diag}(F_1, \dots, F_s)$ for $F_i \in \{E, R, \bar{B}\}$ and $H(x) := \sum_{i=1}^s H_i(x_i)$. The matrix $J(x) := \text{diag}(J_1(x_1), \dots, J_s(x_s)) - \hat{B}\hat{C}\hat{B}^\top$ represents the Schur complement of the block matrix with diagonal blocks $\text{diag}(J_1(x_1), \dots, J_s(x_s))$, \hat{C}^{-1} , and off-diagonal blocks \hat{B}, \hat{B}^\top .

Remark 2. If the matrix \hat{C} in the input–output coupling (5) is not skew-symmetric, then one can split $\hat{C} = \hat{C}^{\text{symm}} + \hat{C}^{\text{skew-symm}}$ into its symmetric and skew-symmetric parts, and one gets an equivalence with the pHODE system (6) by replacing $J(x)$ and $R(x)$ by

$$J(x) := \text{diag}(J_1(x_1), \dots, J_s(x_s)) - \hat{B}\hat{C}^{\text{skew-symm}}\hat{B}^\top \quad \text{and} \quad R(x) := \text{diag}(R_1(x_1), \dots, R_s(x_s)) + \hat{B}\hat{C}^{\text{symm}}\hat{B}^\top,$$

provided that $R(x) \geq 0$ is still positive semidefinite. If not, one must use (7) to follow the approach discussed in the following Remark 3.

Remark 3. In many engineering applications, the inputs and outputs of the subsystems are coupled via a skew-symmetric coupling condition. In general, if an arbitrary linear relation $Mu + Ny = 0$ exists, then the coupled system of pHODEs is equivalent to a pHDAE system (cf. [7, Section 2.2.4]), but this comes at the cost of introducing additional dummy variables \hat{u} and \hat{y} that copy u and y :

$$\begin{bmatrix} E\dot{x} \\ 0 \\ 0 \\ 0 \end{bmatrix} = \begin{bmatrix} \text{diag}(J_1, \dots, J_s) - \text{diag}(R_1, \dots, R_s) & B & 0 & 0 \\ & -B^\top & 0 & I \\ & 0 & -I & 0 \\ & 0 & M & N \end{bmatrix} \begin{bmatrix} \nabla H(x) \\ \hat{u} \\ \hat{y} \\ 0 \end{bmatrix} + \begin{bmatrix} 0 \\ 0 \\ I \\ 0 \end{bmatrix} u, \tag{7a}$$

$$y = \hat{y}. \tag{7b}$$

In other words, one may end up with a ‘‘true’’ DAE even if all the subsystems are ODEs.

4. Decoupling pHODE systems

Decoupling is only possible if there is a diffeomorphic transformation of the state $w = F(x)$ such that the Hamiltonian becomes separable:

$$H(x) = \tilde{H}(w) = \sum_{i=1}^s \tilde{H}_i(w_i), \quad w = \begin{bmatrix} w_1 \\ w_2 \\ \vdots \\ w_s \end{bmatrix}, \quad n_i := \dim(w_i), \quad i = 1, \dots, s, \quad \sum_{i=1}^s n_i = n.$$

In this case, the pHODE system (4) can be transformed into the pHODE system

$$(F'(x)^{-1})^\top E(x)\dot{x} = \underbrace{(F'(x)^{-1})^\top E(x)F'(x)^{-1}}_{\tilde{E}(w)} \dot{w} \tag{8a}$$

$$= \underbrace{(F'(x)^{-1})^\top (J(x) - R(x))F'(x)^{-1}}_{\tilde{J}(w) - \tilde{R}(w)} \underbrace{F'(x)z(x)}_{\tilde{z}(w)} + \underbrace{(F'(x)^{-1})^\top B(x)u}_{\tilde{B}(w)},$$

$$y = \tilde{B}(w)^\top \tilde{z}(w), \tag{8b}$$

with $\tilde{E}^\top \tilde{z}(w) = \nabla_w \tilde{H}(w)$, as $\nabla_x H(x) = F'(x)^\top \nabla_w \tilde{H}(w)$ holds. Note that we have replaced x with $F^{-1}(w)$.

Next, we split \tilde{J} and \tilde{R} with respect to the splitting of w :

$$\tilde{J} = \text{diag}(\tilde{J}_{11}, \dots, \tilde{J}_{ss}) + \tilde{J}^{\text{offdiag}}, \quad \tilde{R} = \text{diag}(\tilde{R}_{11}, \dots, \tilde{R}_{ss}) + \tilde{R}^{\text{offdiag}}, \quad \tilde{B} = \begin{bmatrix} \tilde{B}_1 \\ \vdots \\ \tilde{B}_s \end{bmatrix},$$

where we have used the abbreviations for the $n_i \times n_j$ matrices \tilde{J}_{ij} and \tilde{R}_{ij}

$$\tilde{J}_{ij} := (J_{k,l})_{\substack{k=\sum_{l=0}^{i-1} n_l+1, \dots, \sum_{l=0}^i n_l \\ l=\sum_{l=0}^{j-1} n_l+1, \dots, \sum_{l=0}^j n_l}}, \quad \tilde{R}_{ij} := (R_{k,l})_{\substack{k=\sum_{l=0}^{i-1} n_l+1, \dots, \sum_{l=0}^i n_l \\ l=\sum_{l=0}^{j-1} n_l+1, \dots, \sum_{l=0}^j n_l}}, \quad n_0 := 0.$$

In addition, we assume that \tilde{J}_{ij} , \tilde{R}_{ij} and \tilde{B}_i depend only on w_i (for $i = 1, \dots, s$), and the transformed flow matrix $\tilde{E} = \text{diag}(\tilde{E}_{11}(w_1), \dots, \tilde{E}_{ss}(w_s))$, is block-diagonal, the system can be decomposed into s coupled pHODE subsystems as follows. Note that the block-diagonality of \tilde{E} is an essential structural condition, since off-diagonal blocks would introduce coupling through the time derivatives and prevent a decomposition into independent port-Hamiltonian subsystems.

Case 1: $\tilde{R}^{\text{offdiag}} = 0$

If we define $\hat{B} = I$ and $\hat{C} = -\tilde{J}^{\text{offdiag}}$, then the pHODE system (8) is *analytically equivalent* (i.e. there exists a diffeomorphic change of variables that transforms one system into the other) to the s pHODE subsystems

$$\tilde{E}_{ii}(w_i) \dot{w}_i = (\tilde{J}_{ii}(w_i) - \tilde{R}_{ii}(w_i)) \tilde{z}_i(w_i) + [I, \tilde{B}_i(w_i)] \begin{bmatrix} \hat{u}_i \\ u \end{bmatrix}, \quad \begin{bmatrix} \hat{y}_i \\ \hat{y}_i \end{bmatrix} = [I, \tilde{B}_i(w_i)]^\top \tilde{z}_i(w_i), \tag{9a}$$

coupled by the skew-symmetric coupling condition

$$\begin{bmatrix} \hat{u}_1 \\ \vdots \\ \hat{u}_k \end{bmatrix} + \hat{C} \begin{bmatrix} \hat{y}_1 \\ \vdots \\ \hat{y}_k \end{bmatrix} = 0, \quad \hat{C} = -\hat{C}^\top, \tag{9b}$$

and fulfilling the compatibility conditions $\tilde{E}_{ii}^\top \tilde{z}_i = \nabla_{w_i} \tilde{H}_{w_i}$. This is only one convenient choice. The choice $\hat{B} = I$ is not unique. Any factorization of the bilinear power exchange $-\tilde{J}^{\text{offdiag}} = \hat{B}\hat{C}\hat{B}^\top$ leads to an analytically equivalent interconnection, and alternative choices may reduce the port dimension or better reflect the physical coupling.

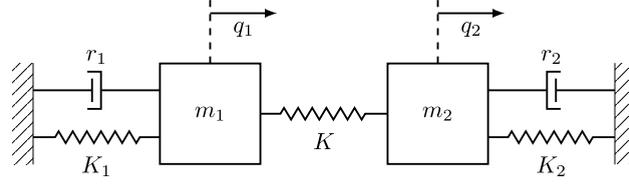


Fig. 1. ODE two masses oscillator with damping [6]. The coordinates q_1, q_2 describe the position of the masses.

Case 2: $\tilde{R}^{\text{offdiag}} \neq 0$

Here we define $\hat{B} = I$ and $\hat{C} = -\tilde{J}^{\text{offdiag}} + \tilde{R}^{\text{offdiag}}$ to obtain the decoupled system (9). However, the coupling condition is no longer skew-symmetric, but a general linear one with $M = I$ and $N = \hat{C}$ (see Remark 2).

Case 3: General coupling

We can use s different port matrices \hat{B}_i to define the equivalent pHODE system

$$\begin{aligned} \tilde{E}_{ii}(w_i)\dot{w}_i &= (\tilde{J}_{ii}(w_i) - \tilde{R}_{ii}(w_i)) \tilde{z}_i(w_i) + [\hat{B}_i, \tilde{B}_i(w_i)] \begin{bmatrix} \hat{u}_i \\ u \end{bmatrix}, \\ \begin{bmatrix} \hat{y}_i \\ \tilde{y}_i \end{bmatrix} &= [\hat{B}_i, \tilde{B}_i(w_i)]^\top \tilde{z}_i(w_i), \end{aligned} \tag{10a}$$

coupled by the input–output coupling condition

$$\begin{pmatrix} \hat{u}_1 \\ \vdots \\ \hat{u}_s \end{pmatrix} + \hat{C} \begin{pmatrix} \hat{y}_1 \\ \vdots \\ \hat{y}_s \end{pmatrix} = 0, \quad \hat{C} := \begin{pmatrix} 0 & \hat{C}_{12} & \dots & \hat{C}_{1s} \\ \hat{C}_{12}^\top & 0 & \dots & \hat{C}_{2s} \\ \vdots & \vdots & \ddots & \vdots \\ \hat{C}_{1s}^\top & \hat{C}_{2s}^\top & \dots & 0 \end{pmatrix} \tag{10b}$$

provided that (a)

$$\tilde{J}_{ij} - \tilde{R}_{ij} = -\hat{B}_i \hat{C}_{ij} \hat{B}_j^\top \Rightarrow \tilde{J}_{ij} = -\hat{B}_i \hat{C}_{ij}^{\text{symm}} \hat{B}_j^\top \text{ and } \tilde{R}_{ij} = \hat{B}_i \hat{C}_{ij}^{\text{skew-symm}} \hat{B}_j^\top,$$

holds with \hat{C}^{symm} and $\hat{C}^{\text{skew-symm}}$ corresponding to the symmetric and skew-symmetric part of \hat{C} , respectively, and (b) the port matrices \hat{B}_i depend only on w_i . The latter can always be achieved by defining $\hat{B}_i = I$ and

$$\hat{C}_{ij} := -(\tilde{J}_{ij} - R_{ij}).$$

Remark 4. In the case of a block-diagonal matrix \tilde{R} , i.e., $\tilde{R}_{ij} = 0$ for $i \neq j$, we obtain a skew-symmetric coupling with

$$\begin{pmatrix} 0 & \hat{C}_{12} & \dots & \hat{C}_{1p} \\ -\hat{C}_{12}^\top & 0 & \dots & \hat{C}_{2p} \\ \vdots & \vdots & \ddots & \vdots \\ -\hat{C}_{1p}^\top & -\hat{C}_{2p}^\top & \dots & 0 \end{pmatrix}.$$

For this case, the obvious choice in the last section is $\hat{B}_i = \hat{B}_j = I, \hat{C}_{ij} = -\tilde{J}_{ij}$.

Remark 5. Since R is positive semidefinite, all R_{ii} are also positive semidefinite. Additionally, if J is skew-symmetric, then the block-diagonal matrices J_{ii} are also skew-symmetric, which defines s pHODE systems.

Example 1. Consider the dissipative Hamiltonian system (pHODE system without external input, see Fig. 1)

$$\dot{x} = (J - R) \nabla H(x),$$

with

$$J = \begin{bmatrix} 0 & -1 & -1 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & -1 & 0 \\ 0 & 0 & 1 & 0 & -1 \\ 0 & 0 & 0 & 1 & 0 \end{bmatrix}, \quad R = \begin{bmatrix} r_1 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & r_2 & 0 \\ 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \quad H(x) = x^\top \begin{bmatrix} \frac{1}{m_1} & 0 & 0 & 0 & 0 \\ 0 & K_1 & 0 & 0 & 0 \\ 0 & 0 & K & 0 & 0 \\ 0 & 0 & 0 & \frac{1}{m_2} & 0 \\ 0 & 0 & 0 & 0 & K_2 \end{bmatrix} x.$$

Next, we want to decouple this pHODE into two pHODE subsystems

$$\begin{aligned} \dot{x}_1 &= (J_1 - R_1) \nabla_1 H_1(x_1) + B_1 u_1, & \dot{x}_2 &= (J_2 - R_2) \nabla_2 H_2(x_2) + B_2 u_2, \\ y_1 &= B_1^\top \nabla_1 H_1(x_1), & y_2 &= B_2^\top \nabla_2 H_2(x_2), \\ & & 0 &= u + C y. \end{aligned}$$

We have a lot of flexibility here: for example, if we split with respect to the first three and last two components, we have

$$J_1 = \begin{bmatrix} 0 & -1 & -1 \\ 1 & 0 & 0 \\ 1 & 0 & 0 \end{bmatrix}, \quad J_2 = \begin{bmatrix} 0 & -1 \\ 1 & 0 \end{bmatrix}, \quad R_1 = \begin{bmatrix} r_1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}, \quad R_2 = \begin{bmatrix} r_2 & 0 \\ 0 & 0 \end{bmatrix}$$

$$H_1(x_1) = x_1^\top \begin{bmatrix} \frac{1}{m_1} & 0 & 0 \\ 0 & K_1 & 0 \\ 0 & 0 & K \end{bmatrix} x_1, \quad H_2(x_2) = x_2^\top \begin{bmatrix} \frac{1}{m_1} & 0 \\ 0 & K_2 \end{bmatrix} x_2.$$

The flexibility lies in the definition of the port matrices B_1 and B_2 :

(a) We set $B_1 = I_3$ and $B_2 = I_2$ and

$$C = - \begin{bmatrix} 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 \end{bmatrix} = -J^{\text{offdiag}}.$$

(b) Here, we set

$$B_1 = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix}, \quad B_2 = \begin{bmatrix} -1 \\ 0 \end{bmatrix}, \quad C = \begin{bmatrix} 0 & C_{12} \\ -C_{12}^\top & 0 \end{bmatrix}, \quad C_{12} = -1,$$

and check

$$B_1 C_{12} B_2^\top = \begin{bmatrix} 0 \\ 0 \\ 1 \end{bmatrix} \cdot (-1) \cdot [-1, \quad 0] = \begin{bmatrix} 0 & 0 \\ 0 & 0 \\ 1 & 0 \end{bmatrix} = -J_{12} \quad \text{with } J = \begin{bmatrix} J_1 & J_{12} \\ -J_{12}^\top & J_2 \end{bmatrix}.$$

This splitting is in accordance with the modeling of a two-mass and three-spring system with damping, as described in [6]. This is achieved by setting $x_1 = [p_1, q_1, q_1 - q]^\top$ and $x_2 = [p_2, q_2]^\top$. Here $p_i = m_i \dot{q}_i$ denotes the momentum of mass i , and q represents the displacement of the internal node connecting the middle spring.

Remark 6. Properly choosing the port matrices minimizes the dimension of the coupling condition, i.e., the coupling matrix \hat{C} . With the former, we need only one scalar input for each subsystem. However, choosing B_1 and B_2 as identities requires as many inputs as there are state variables.

Another splitting splits with respect to components 1, 4 and components 2, 3, 5. In this case we have

$$J_1 = \begin{bmatrix} 0 & 0 \\ 0 & 0 \end{bmatrix}, \quad J_2 = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}, \quad R_1 = \text{diag}(r_1, r_2), \quad R_2 = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix},$$

$$H_1(x_1) = x_1^\top \text{diag}\left(\frac{1}{m_1}, \frac{1}{m_2}\right) x_1, \quad H_2 = x_2^\top \text{diag}(K_1, K, K_2) x_2.$$

A nontrivial choice of the port matrices is

$$B_1 = \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix}, \quad B_2 = \begin{bmatrix} -1 & 1 \\ 0 & 1 \\ 1 & 0 \end{bmatrix}, \quad C_{12} = \begin{bmatrix} 0 & 1 \\ -1 & 0 \end{bmatrix},$$

and we check $B_1 C_{12} B_2^\top = -J_{12}$.

One may also split the system into 3, 4 and 5 subsystems as well.

Example 2 (Poroelastic Network Model). A discretized version of the poroelastic network model [8] reads

$$\begin{aligned} E \dot{x} &= (J - R)x + Bu, \\ y &= B^\top x, \end{aligned}$$

with

$$E = \begin{bmatrix} \rho M_u & 0 & 0 \\ 0 & K_u(\mu, \lambda) & 0 \\ 0 & 0 & \frac{1}{M} M_p \end{bmatrix}, \quad J = \begin{bmatrix} 0 & -K_u(\mu, \lambda) & \alpha D^\top \\ K_u(\mu, \lambda)^\top & 0 & 0 \\ -\alpha D & 0 & 0 \end{bmatrix}, \quad R = \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & \frac{\kappa}{v} K_p \end{bmatrix},$$

$$B = \begin{bmatrix} B_f & 0 \\ 0 & 0 \\ 0 & B_g \end{bmatrix}, \quad x = \begin{bmatrix} w(t) \\ v(t) \\ p(t) \end{bmatrix}, \quad u = \begin{bmatrix} f(t) \\ g(t) \end{bmatrix},$$

where w, v, p are the discretized velocity, displacement, and pressure fields, f, g are the discretized volume-distributed forces and the external injection, μ and λ are the Lamé coefficients, α is the Biot–Willes fluid–solid coupling coefficient, κ is the permeability and ρ is the density. M_u and M_p are mass matrices with respect to u and p , K_u is a stiffness matrix with respect to u and D is a damping matrix, see [9] for more details. We can easily split this implicit pHODE system with respect to the first two and the third block and obtain

$$\begin{bmatrix} \rho M_u & 0 \\ 0 & K_u(\mu, \lambda) \end{bmatrix} \cdot \begin{bmatrix} \dot{w}(t) \\ \dot{v}(t) \end{bmatrix} = \begin{bmatrix} 0 & -K_u(\mu, \lambda) \\ K_u(\mu, \lambda)^\top & 0 \end{bmatrix} \cdot \begin{bmatrix} w(t) \\ v(t) \end{bmatrix} + \begin{bmatrix} \alpha D^\top \\ 0 \end{bmatrix} u_1, \tag{11a}$$

$$y_1 = \begin{bmatrix} \alpha D^\top \\ 0 \end{bmatrix}^\top \begin{bmatrix} w(t) \\ v(t) \end{bmatrix} = \alpha D w(t), \tag{11b}$$

$$\frac{1}{M} M_p \dot{p} = -\frac{\kappa}{v} K_p p + u_2, \tag{11c}$$

$$0 = \begin{bmatrix} y_1 \\ u_2 \end{bmatrix} = \begin{bmatrix} 0 & I \\ -I & 0 \end{bmatrix} \begin{bmatrix} y_1 \\ y_2 \end{bmatrix}. \tag{11e}$$

Because the first subsystem lacks a dissipative part and the second lacks a skew-symmetric part, symplectic time integration can be used for the first part and dissipative time integration can be used for the second part in a dynamic iteration procedure.

Example 3 (Discretized Full Set of Maxwell’s Equations). The full set of Maxwell’s equations formulated with electrodynamic potentials after discretization forms a pHDAE system with

$$\underbrace{\begin{bmatrix} M_\epsilon & M_\epsilon G & 0 \\ G^\top M_\epsilon & G^\top M_\epsilon G & 0 \\ 0 & 0 & M_\mu \end{bmatrix}}_{:=E_{\text{Maxwell}}} d_t \underbrace{\begin{bmatrix} d_t a \\ \varphi \\ h \end{bmatrix}}_{:=x} = \underbrace{\begin{bmatrix} 0 & 0 & -C^\top \\ 0 & 0 & 0 \\ C & 0 & 0 \end{bmatrix}}_{:=J(=-J^\top)} - \underbrace{\begin{bmatrix} M_\kappa & M_\kappa G & 0 \\ G^\top M_\kappa & G^\top M_\kappa G & 0 \\ 0 & 0 & 0 \end{bmatrix}}_{:=R \geq 0} \underbrace{\begin{bmatrix} d_t a \\ \varphi \\ h \end{bmatrix}}_{:=z} + \begin{bmatrix} I \\ G^\top \\ 0 \end{bmatrix} j_s, \tag{12}$$

where C, G are discrete curl and gradient matrices, $M_\kappa, M_\epsilon, M_\mu$ are discrete material matrices for electric conductivity κ , permittivity ϵ and permeability μ , respectively, a, φ, h are the degree of freedom vectors and $u = j_s$ is the input vector of source currents and the output vector $y = d_t a + G\varphi (= -e)$ is the negated vector of electric grid voltages. The Hamiltonian associated with (12) is the discrete electromagnetic grid energy given by

$$H = \frac{1}{2} (\varphi^\top G^\top M_\epsilon G \varphi + h^\top M_\mu h) + \varphi^\top G^\top M_\epsilon d_t a + d_t a^\top M_\epsilon d_t a, \tag{13}$$

and the compatibility condition holds with

$$\nabla H = (M_\epsilon(d_t a + G\varphi), G^\top M_\epsilon(d_t a + G\varphi), M_\mu h)^\top = E_{\text{Maxwell}}^\top x. \tag{14}$$

Again, we can easily split with respect to the first two and the last block with Hamiltonians

$$H_1 = \frac{1}{2} \varphi^\top G^\top M_\epsilon G \varphi + \varphi^\top G^\top M_\epsilon d_t a + d_t a^\top M_\epsilon d_t a \quad \text{and} \quad H_2 = \frac{1}{2} h^\top M_\mu h,$$

$$\begin{bmatrix} M_\epsilon & M_\epsilon G \\ G^\top M_\epsilon & G^\top M_\epsilon G \end{bmatrix} d_t \begin{bmatrix} d_t a \\ \varphi \end{bmatrix} = \begin{bmatrix} M_\kappa & M_\kappa G \\ G^\top M_\kappa & G^\top M_\kappa G \end{bmatrix} \begin{bmatrix} d_t a \\ \varphi \end{bmatrix} + \begin{bmatrix} I \\ G^\top \end{bmatrix} j_s + \begin{bmatrix} C^\top \\ 0 \end{bmatrix} u_1,$$

$$\bar{y}_1 = \begin{bmatrix} I \\ G^\top \end{bmatrix}^\top \begin{bmatrix} d_t a \\ \varphi \end{bmatrix}, \quad y_1 = \begin{bmatrix} C^\top \\ 0 \end{bmatrix}^\top \begin{bmatrix} d_t a \\ \varphi \end{bmatrix},$$

$$M_\mu d_t h = u_2, \quad y_2 = h, \quad 0 = \begin{bmatrix} u_1 \\ u_2 \end{bmatrix} + \begin{bmatrix} 0 & I \\ -I & 0 \end{bmatrix} \begin{bmatrix} y_1 \\ y_2 \end{bmatrix}.$$

5. Numerical aspects of monolithic and decoupled formulations

In numerical simulations, the monolithic and decoupled formulations of port-Hamiltonian ODE systems offer distinct advantages. The monolithic formulation preserves the full system structure and typically exhibits superior stability properties, particularly when integrated with structure-preserving or symplectic time-stepping schemes. This makes it ideal for stiff or highly coupled

systems where global energy conservation is crucial. Nevertheless, one may reduce computational time by applying operator splitting techniques to the general systems which preserve the port-Hamiltonian structure, for example, $J - R$ coupling [10] or diagonal/off-diagonal coupling [11], jointly with multirate approaches.

In contrast, the decoupled formulation offers larger flexibility and computational efficiency in multiphysics or large-scale contexts. It allows for the use of specialized numerical solvers, such as mimetic or energy-consistent methods, which are tailored to individual subsystems and respect local conservation laws. This partitioned approach also enables parallel or distributed computation via dynamic iteration [12], although with potentially weaker global stability. The monolithic formulation favors global stability, whereas the decoupled approach provides modularity and solver flexibility, at the cost of potentially reduced overall stability.

6. Conclusion

We have presented a unified framework for coupling and decoupling nonlinear port-Hamiltonian ordinary differential equation systems. The analysis demonstrates how multiple pH subsystems can be reformulated as a single monolithic system, and conversely, how a monolithic pHODE can be decomposed into weakly coupled subsystems without losing the underlying Hamiltonian structure. The results provide a theoretical basis for structure-preserving simulation strategies that balance stability and flexibility. Future research will address adaptive coupling strategies, structure-preserving time discretizations, and extensions to port-Hamiltonian differential–algebraic [13] and stochastic systems [14].

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Data availability

No data was used for the research described in the article.

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