# PYPHS DOCUMENTATION Version 0.1.9b2

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# 1 Introduction

The python package pyphs is dedicated to the treatment of passive multiphysical systems in the Port-Hamiltonian Systems (PHS) formalism. This formalism structures physical systems into

- energy conserving parts,
- power dissipating parts and
- source parts.

This guarantees a *power balance* is fulfilled, including for numerical *simulations* based on an adapted *numerical method*.

- 1. Systems are described by directed multi-graphs (networkx.MultiDiGraph).
- 2. The time-continuous port-Hamiltonian structure is build from an  $auto-mated\ graph\ analysis.$
- 3. The discrete-time port-Hamiltonian structure is derived from a *structure* preserving numerical method.
- 4. LATEX description code and C++ simulation code are automatically generated.

## 1.1 Installation

## Notice only python 2.7 is supported.

It is recommanded to install pyphs using PyPI (the Python Package Index). In terminal :

pip install pyphs

Mac OSX only: An installation for *Anaconda* users is also available. In terminal:

conda install -c afalaize pyphs

## 1.2 The PHS formalism

Below is a recall of the Port-Hamiltonian Systems (PHS) formalism. For details, the reader is referred to the e.g. the acaemic reference [?]. We consider

systems that can be described by the following time-continuous non-linear state-space representation :  $\frac{1}{2}$ 

$$\underbrace{\begin{pmatrix} \frac{\mathrm{d}\mathbf{x}}{\mathrm{d}t} \\ \mathbf{w} \\ \mathbf{y} \end{pmatrix}}_{\mathbf{b}} = \underbrace{\begin{pmatrix} \mathbf{M}_{xx} & \mathbf{M}_{xw} & \mathbf{M}_{xy} \\ \mathbf{M}_{wx} & \mathbf{M}_{ww} & \mathbf{M}_{wy} \\ \mathbf{M}_{yx} & \mathbf{M}_{yw} & \mathbf{M}_{yy} \end{pmatrix}}_{\mathbf{M}} \cdot \underbrace{\begin{pmatrix} \nabla \mathbf{H}(\mathbf{x}) \\ \mathbf{z}(\mathbf{w}) \\ \mathbf{u} \end{pmatrix}}_{\mathbf{a}} \tag{1}$$

where

$$\mathbf{M} = \underbrace{\begin{pmatrix} \mathbf{J}_{xx} & \mathbf{J}_{xw} & \mathbf{J}_{xy} \\ \mathbf{J}_{yx} & \mathbf{J}_{yw} & \mathbf{J}_{yy} \\ \mathbf{J}_{yx} & \mathbf{J}_{yw} & \mathbf{J}_{yy} \end{pmatrix}}_{\mathbf{I}} - \underbrace{\begin{pmatrix} \mathbf{R}_{xx} & \mathbf{R}_{xw} & \mathbf{R}_{xy} \\ \mathbf{R}_{wx} & \mathbf{R}_{ww} & \mathbf{R}_{wy} \\ \mathbf{R}_{yx} & \mathbf{R}_{yw} & \mathbf{R}_{yy} \end{pmatrix}}_{\mathbf{R}}$$
(2)

and

—  $\mathbf{J}: \mathbf{x} \mapsto \mathbf{J}(\mathbf{x})$  is a skew-symmetric matrix :

$$\mathbf{J}_{\alpha\beta} = -\mathbf{J}_{\beta\alpha}^{\mathsf{T}} \text{ for } (\alpha,\beta) \in \{\mathsf{x},\mathsf{w},\mathsf{y}\}^2,$$

- $\mathbf{R}: \mathbf{x} \mapsto \mathbf{R}(\mathbf{x}) \succeq 0$  is a positive definite matrix,
- $\mathbf{x}: t \mapsto \mathbf{x}(t) \in \mathbb{R}^{n_{\mathbf{x}}}$  is the state vector,
- $H: \mathbf{x} \mapsto H(\mathbf{x}) \in \mathbb{R}_+$  is a *storage function* (convex and positive-definite scalar function with H(0) = 0),
- $\nabla H : \mathbf{x} \mapsto \nabla H(\mathbf{x}) \in \mathbb{R}^{n_{\mathbf{x}}}$  denote the gradient of the storage function with the *storage power*

$$P_{\mathbf{x}} = \frac{\mathrm{d}\mathbf{x}}{\mathrm{d}t} \cdot \nabla \mathbf{H}(\mathbf{x}),$$

- $\mathbf{w}: t \mapsto \mathbf{w}(t) \in \mathbb{R}^{n_{\mathbf{w}}}$  is the dissipation vector variable,
- $-\mathbf{z}: \mathbf{w} \mapsto \mathbf{z}(\mathbf{w}) \in \mathbb{R}^{n_{\mathbf{v}}}$  is a dissipation function (with positive definite jacobian matrix and  $\mathbf{z}(0) = 0$ ) for the dissipated power

$$P_{\mathbf{w}} = \mathbf{w} \cdot \mathbf{z}(\mathbf{w}) + \mathbf{a} \cdot \mathbf{R} \cdot \mathbf{a},$$

- $\mathbf{u}: t \mapsto \mathbf{u}(t) \in \mathbb{R}^{n_y}$  is the input vector,
- $\mathbf{y}: t \mapsto \mathbf{y}(t) \in \mathbb{R}^{n_y}$  is the output vector,
- the power received by the sources from the system is

$$P = \mathbf{u} \cdot \mathbf{v}$$
.

The state is split according to  $\mathbf{x} = (\mathbf{x}_1^\intercal, \, \mathbf{x}_{n1}^\intercal)^\intercal$  with

 $\mathbf{x}_1 = (x_1, \cdots, x_{n_{\mathbf{x}1}})^\intercal$  the states associated with the quadratic components of the storage function  $\mathbf{H}_1(\mathbf{x}_1) = \frac{\mathbf{x}_1 \cdot \mathbf{Q} \cdot \mathbf{x}_1}{2}$ 

 $\mathbf{x}_{\mathtt{nl}} = (x_{n_{\mathtt{xl}}+1}, \cdots, x_{n_{\mathtt{x}}})^{\intercal}$  the states associated with the non-quadratic components of the storage function with  $n_{\mathtt{x}} = n_{\mathtt{xl}} + n_{\mathtt{xnl}}$  and

$$H(\mathbf{x}) = H_1(\mathbf{x}_1) + H_{n1}(\mathbf{x}_{n1})$$

.

The set of dissipative variables is split according to  $\mathbf{w} = (\mathbf{x}_1^\intercal, \, \mathbf{w}_{\mathtt{n}1}^\intercal)^\intercal$  with

 $\mathbf{w}_1 = (w_1, \dots, w_{n_{\mathbf{v}_1}})^{\mathsf{T}}$  the variables associated with the linear components of the dissipative relation  $\mathbf{z}_1(\mathbf{w}_1) = \mathbf{Z}_1 \mathbf{w}_1$ 

 $\mathbf{w}_{\mathtt{nl}} = (w_{n_{\mathtt{vl}}+1}, \cdots, w_{n_{\mathtt{v}}})^{\intercal}$  the variables associated with the nonlinear components of the dissipative relation  $\mathbf{z}_{\mathtt{nl}} : \mathbf{w}_{\mathtt{nl}} \mapsto \mathbf{z}_{\mathtt{nl}}(\mathbf{w}_{\mathtt{nl}}) \in \mathbb{R}^{n_{\mathtt{wnl}}}$  with  $n_{\mathtt{w}} = n_{\mathtt{wl}} + n_{\mathtt{wnl}}$  and

$$\mathbf{z}(\mathbf{w}) = \left( \begin{array}{c} \mathbf{Z}_{\text{l}} \, \mathbf{w}_{\text{l}} \\ \mathbf{z}_{\text{nl}}(\mathbf{w}_{\text{nl}}) \end{array} \right).$$

Accordingly, the structure matrices are split as

$$\underbrace{\begin{pmatrix} \frac{\mathrm{d} x_1}{\mathrm{d} t} \\ \frac{\mathrm{d} x_{\mathrm{nl}}}{\mathrm{d} t} \\ \frac{\mathrm{d} t}{\mathrm{w}_1} \\ \underline{y} \end{pmatrix}}_{b} = \underbrace{\begin{pmatrix} M_{\text{xlx1}} & M_{\text{xlxnl}} & M_{\text{xlw1}} & M_{\text{xlw1}} & M_{\text{xlwnl}} & M_{\text{xly}} \\ M_{\text{xnlx1}} & M_{\text{xnlxnl}} & M_{\text{xnlw1}} & M_{\text{xnlwnl}} & M_{\text{xnly}} \\ M_{\text{wlx1}} & M_{\text{wlxnl}} & M_{\text{wlxnl}} & M_{\text{wlwnl}} & M_{\text{wly}} \\ M_{\text{wnlw1}} & M_{\text{wnlw1}} & M_{\text{wnlw1}} & M_{\text{wnlwnl}} & M_{\text{wnly}} \\ M_{\text{yx1}} & M_{\text{yxnl}} & M_{\text{yxnl}} & M_{\text{ywl}} & M_{\text{ywnl}} & M_{\text{yy}} \end{pmatrix}}_{M} \cdot \underbrace{\begin{pmatrix} \mathbf{Q} \cdot \mathbf{x} \\ \nabla H_{n1}(\mathbf{x}_{n1}) \\ \mathbf{Z}_{1} \cdot \mathbf{w}_{1} \\ \mathbf{Z}_{n1}(\mathbf{w}_{n1}) \\ \mathbf{u} \end{pmatrix}}_{\mathbf{a}}_{a}$$

$$(3)$$

# Table des matières

1	Introduction				
	1.1	Installation	1		
	1.2	The PHS formalism	1		
<b>2</b>	Structure of the pyphs.PortHamiltonianObject				
	2.1	The symbs module	4		
	2.2	The exprs module	4		
	2.3	The dims module	5		
3	Alg	orithms	6		
	3.1	Graph analysis	6		
	3.2	- *	6		
		3.2.1 Split of the linear part from the nonlinear part	6		
		3.2.2 Presolve numerical nonlinear subsystem	7		
		3.2.3 Algorithm	7		
	3.3	Realizability solver	8		
4	Connectors : Gyrators and Transformers				
	4.1	Gyrator	9		
	4.2	Transformer	9		
5	Fractional calculus				
	5.1	Fractional integrator	9		
	5.2	~	10		
	5.3		10		

# 2 Structure of the pyphs.PortHamiltonianObject

Below is a list of each module of practical use in the object pyphs.PortHamiltonianObject, along with a short description. We consider the following instantiation:

```
# import of (pre-installed) pyphs package:
import pyphs

# instantiate the PortHamiltonianObject:
phs = pyphs.PortHamiltonianObject(label='mylabel')
```

# 2.1 The symbs module

Container for all the SYMPY symbolic variables (sympy.Symbol).

**Attributes** are ordered *list of symbols* associated with the system's vectors components.

```
\begin{array}{l} {\tt phs.symbs.x: state \ vector \ symbols \ } {\bf x} \in \mathbb{R}^{n_{\rm x}}, \\ {\tt phs.symbs.w: dissipative \ vector \ variable \ symbols \ } {\bf w} \in \mathbb{R}^{n_{\rm w}}, \\ {\tt phs.symbs.u: input \ vector \ symbols \ } {\bf u} \in \mathbb{R}^{n_{\rm y}}, \\ {\tt phs.symbs.y: output \ vector \ symbols \ } {\bf y} \in \mathbb{R}^{n_{\rm y}}, \\ {\tt phs.symbs.cu: input \ vector \ symbols \ for \ connectors \ } {\bf c_u} \in \mathbb{R}^{n_{\rm y}}, \\ {\tt phs.symbs.cy: output \ vector \ symbols \ for \ connectors \ } {\bf c_y} \in \mathbb{R}^{n_{\rm y}}, \\ {\tt phs.symbs.p: Time-varying \ parameters \ symbols \ } {\bf p} \in \mathbb{R}^{n_{\rm y}}. \end{array}
```

## Methods:

phs.symbs.dx(): Returns the symbols associated with the state differential dx formed by appending the prefix d to each symbol in x.

phs.symbs.args() : Return the list of symbols associated with the vector
 of all arguments of the symbolic expressions (expr module).

## 2.2 The exprs module

Container for all the  $\operatorname{SYMPY}$  symbolic expressions  $\operatorname{sympy}$ .  $\operatorname{Exprassociated}$  with the system's functions.

Attributes: For scalar function (e.g. the storage function H), arguments of phs.exprs are SYMPY expressions (sympy.Expr); for vector functions (e.g. the disipative function z), arguments are ordered lists of SYMPY expressions; for matrix functions (e.g. the Jacobian matrix of disipative function z), arguments are sympy.Matrix objects. Notice the expressions arguments 1 must belong either to (i) the elements of phs.symbs.args(), or (ii) the keys of the dictionary phs.symbs.subs.

```
phs.exprs.H : storage function H \in \mathbb{R},
phs.exprs.z : dissipative function \mathbf{z} \in \mathbb{R}^{n_{\mathbf{v}}},
phs.exprs.g : input/output gains vector function \mathbf{g} \in \mathbb{R}^{n_{\mathbf{g}}},
```

<sup>1.</sup> Accessed through the sympy.Expr.free\_symbols (e.g. phs.exprs.H.free\_symbols to recover the arguments of the Storage function H).

The following expression are computed from the exprs.build() method (see below):

phs.exprs.dxH: the continuous gradient vector of storage scalar function  $\nabla \mathbf{H}(\mathbf{x}) \in \mathbb{R}^{n_{\mathbf{x}}}$ ,

phs.exprs.dxHd : the discrete gradient vector of storage scalar function  $\overline{\nabla} H(\mathbf{x}, \delta \mathbf{x}) \in \mathbb{R}^{n_{\mathbf{x}}}$ ,

phs.exprs.hessH : the continuous hessian matrix of storage scalar function (computed as  $\nabla \nabla H(\mathbf{x}) \in \mathbb{R}^{n_{\mathbf{x}} \times n_{\mathbf{x}}}$ ),

phs.exprs.jacz : the continuous jacobian matrix of dissipative vector function  $\nabla \mathbf{z}(\mathbf{w}) \in \mathbb{R}^{n_{\mathbf{w}} \times n_{\mathbf{w}}}$ .

phs.exprs.y : the expression of the continuous output vector function  $\mathbf{y}(\nabla \mathbf{H}, \mathbf{z}, \mathbf{u}) \in \mathbb{R}^{n_y}$ ,

phs.exprs.yd : the expression of the discrete output vector function  $\overline{\mathbf{y}}(\overline{\nabla}\mathbf{H},\mathbf{z},\mathbf{u}) \in \mathbb{R}^{n_y}$ ,

### Methods:

phs.exprs.build() : Build the following system functions as SYMPY expressions and append them as attributes to the phs.exprs module :
 phs.exprs.dxH, phs.exprs.dxHd, phs.exprs.hessH, phs.exprs.jacz, phs.exprs.y,
 and phs.exprs.yd.

phs.exprs.setexpr(name, expr) : Add the SYMPY expression expr to the phs.exprs module, with argument name, and add name to the set of phs.exprs.\_names.

phs.exprs.freesymbols() : Return a python set of all the free symbols
 (sympy.Symbol) that appear at least once in all expressions with names
 in phs.exprs.\_names.

## 2.3 The dims module

Container for accessors to the system's dimensions. No attributes should be changed manually. To split the system into its linear and nonlinear part, use phs.split\_linear() which organize the system vectors as

$$\mathbf{x} = \begin{pmatrix} \mathbf{x}_1 \\ \mathbf{x}_{n1} \end{pmatrix}, \quad \dim(\mathbf{x}_1) =$$
 (4)

Attributes: phs.dims.xl: Number of state vector components associated with a quadratic storage function:  $H_1(\mathbf{x}_1) = \mathbf{x}_1^\intercal \cdot \frac{\mathbf{Q}}{2} \cdot \mathbf{x}_1$ , and phs.dims.x() is equal to phs.dims.xl + phs.dims.xnl().

phs.dims.wl: Number of dissipative vector variable components associated with a linear dissipative function:  $\mathbf{z}_1(\mathbf{w}_1) = \mathbf{Z}_1 \cdot \mathbf{w}_1$ , and phs.dims.w() is equal to phs.dims.wl + phs.dims.wnl().

## Methods:

phs.dims.x() : Return the dimension of state vector len(phs.symbs.x).

phs.dims.xnl() : Return the number of state vector components associated with a nonlinear storage function

len(phs.symbs.x).

setexpr(name, expr): Add the SYMPY expression expr to the exprs module, with argument name, and add name to the set of exprs.\_names.

freesymbols() : Retrun a python set of all the free symbols (sympy.symbols)
that appear at least once in all expressions with names in exprs.\_names.

# 3 Algorithms

This section details the algorithms actually implemented for

- 1. the graph analysis and
- 2. the different simulation methods

## 3.1 Graph analysis

The graph analysis method that derives the port-Hamiltonian system's differential-algebraic equations from with a given netlist is detailed in the reference [?]. The algorithm implemented in PYPHS is exactly that in [?, algorithm 1].

## 3.2 Simulation methods

The discrete gradient method is used in conjunction with the port-Hamiltonian structure to produce a passive-guaranteed numerical scheme (see [?] for details). In the sequel, quantities are defined on the current time step  $\mathbf{x} \equiv \mathbf{x}(t_k)$ , with  $k \in \mathbb{N}_+^*$ .

## 3.2.1 Split of the linear part from the nonlinear part

The dicrete gradient for the quadratic part of the Hamiltonian is  $\nabla H_1 = \frac{1}{2} \mathbf{Q} (2\mathbf{x}_1 + \delta \mathbf{x}_1)$  and the discret linear subsystem is

$$\mathbf{D}_{1}^{-1} = \mathbf{i}\mathbf{D}_{1} = \begin{pmatrix} \frac{\mathbf{I}_{d}}{\delta t} & 0 \\ 0 & \mathbf{I}_{d} \end{pmatrix} - \begin{pmatrix} \mathbf{M}_{x1x1} & \mathbf{M}_{x1w1} \\ \mathbf{M}_{w1x1} & \mathbf{M}_{w1w1} \end{pmatrix} \begin{pmatrix} \frac{1}{2} \mathbf{Q} & 0 \\ 0 & \mathbf{Z}_{1} \end{pmatrix}, \\ \underbrace{\begin{pmatrix} \delta \mathbf{x}_{1} \\ \mathbf{w}_{1} \end{pmatrix}}_{\mathbf{v}_{1}} = \mathbf{D}_{1} \underbrace{\begin{pmatrix} \mathbf{M}_{x1x1} \\ \mathbf{M}_{w1x1} \end{pmatrix} \mathbf{Q}}_{\mathbf{N}_{1x1}} \mathbf{x}_{1} + \mathbf{D}_{1} \underbrace{\begin{pmatrix} \mathbf{M}_{x1xn1} & \mathbf{M}_{x1wn1} \\ \mathbf{M}_{w1xn1} & \mathbf{M}_{w1wn1} \end{pmatrix}}_{\mathbf{N}_{1n1}} \underbrace{\begin{pmatrix} \nabla \mathbf{H}_{n1} \\ \mathbf{z}_{n1} \end{pmatrix}}_{\mathbf{N}_{1y}} + \mathbf{D}_{1} \underbrace{\begin{pmatrix} \mathbf{M}_{x1y} \\ \mathbf{M}_{w1y} \end{pmatrix}}_{\mathbf{N}_{1y}} \mathbf{v}_{1y}$$

$$(5)$$

and the nonlinear subsystem is

$$\begin{pmatrix} \frac{\mathbf{I_{d}}}{\delta t} & 0 \\ 0 & \mathbf{I_{d}} \end{pmatrix} \underbrace{\begin{pmatrix} \delta \mathbf{x_{nl}} \\ \mathbf{w_{nl}} \end{pmatrix}}_{\mathbf{v_{nl}}} = \underbrace{\begin{pmatrix} \mathbf{M_{xnlxnl}} & \mathbf{M_{xnlwnl}} \\ \mathbf{M_{wnlxnl}} & \mathbf{M_{wnlwnl}} \end{pmatrix}}_{\mathbf{N_{mln}}} \mathbf{f_{nl}} + \underbrace{\begin{pmatrix} \mathbf{M_{xnly}} \\ \mathbf{M_{wnly}} \end{pmatrix}}_{\mathbf{N_{mly}}} \mathbf{u}$$

$$+ \underbrace{\begin{pmatrix} \mathbf{M_{xnlxnl}} & \mathbf{M_{xnlwnl}} \\ \mathbf{M_{wnlxnl}} & \mathbf{M_{wnlwl}} \end{pmatrix}}_{\mathbf{N_{mln}}} \begin{pmatrix} \frac{1}{2} \mathbf{Q} & 0 \\ 0 & \mathbf{Z_{l}} \end{pmatrix}}_{\mathbf{N_{nlx}}} \mathbf{v_{l}} + \underbrace{\begin{pmatrix} \mathbf{M_{xnlxnl}} \\ \mathbf{M_{wnlxnl}} \end{pmatrix}}_{\mathbf{N_{nlxnl}}} \mathbf{Q} \mathbf{x_{l}}$$

$$(6)$$

## 3.2.2 Presolve numerical nonlinear subsystem

$$\begin{pmatrix}
\frac{\mathbf{I_d}}{\delta t} & 0 \\
0 & \mathbf{I_d}
\end{pmatrix} \mathbf{v_{nl}} = \underbrace{(\overline{\mathbf{N}_{nlxl}} + \overline{\mathbf{N}_{nl1}} \, \mathbf{N}_{lxl})}_{\mathbf{N_{nlxl}}} \mathbf{N}_{1xl} \underbrace{\mathbf{x}_1 + \underbrace{(\overline{\mathbf{N}_{nlnl}} + \overline{\mathbf{N}_{nl1}} \, \mathbf{N}_{lnl})}_{\mathbf{N}_{nlnl}} \mathbf{f_{nl}}}_{\mathbf{N}_{nlnl}} \mathbf{f_{nl}}$$

$$\underbrace{(\overline{\mathbf{N}_{nly}} + \overline{\mathbf{N}_{nlxl}} \, \mathbf{N}_{ly})}_{\mathbf{N}_{nly}} \mathbf{u}$$
(7)

#### 3.2.3Algorithm

Inputs

$$\begin{array}{lll} \mathbf{i} \mathbf{D_{1}} & = & \left( \begin{array}{c} \frac{\mathbf{I_{d}}}{\delta t} & 0 \\ 0 & \mathbf{I_{d}} \end{array} \right) - \left( \begin{array}{c} \mathbf{M_{x1x1}} & \mathbf{M_{x1w1}} \\ \mathbf{M_{w1x1}} \end{array} \right) \left( \begin{array}{c} \frac{1}{2} \mathbf{Q} & 0 \\ 0 & \mathbf{Z_{1}} \end{array} \right) \\ \overline{\mathbf{N_{1x1}}} & = & \left( \begin{array}{c} \mathbf{M_{x1x1}} \\ \mathbf{M_{w1x1}} \end{array} \right) \mathbf{Q} \\ \overline{\mathbf{N_{1n1}}} & = & \left( \begin{array}{c} \mathbf{M_{x1x1}} & \mathbf{M_{x1wn1}} \\ \mathbf{M_{w1xn1}} & \mathbf{M_{x1wn1}} \\ \mathbf{M_{w1xn1}} & \mathbf{M_{w1wn1}} \end{array} \right) \\ \overline{\mathbf{N_{1y}}} & = & \left( \begin{array}{c} \mathbf{M_{x1x1}} & \mathbf{M_{xn1wn1}} \\ \mathbf{M_{w1xn1}} & \mathbf{M_{xn1wn1}} \\ \mathbf{M_{wn1xn1}} & \mathbf{M_{xn1wn1}} \\ \mathbf{M_{wn1x1}} & \mathbf{M_{xn1w1}} \end{array} \right) \\ \overline{\mathbf{N_{n11}}} & = & \left( \begin{array}{c} \mathbf{M_{xn1x1}} \\ \mathbf{M_{xn1x1}} \\ \mathbf{M_{wn1x1}} \\ \mathbf{M_{wn1x1}} \end{array} \right) \mathbf{Q} \\ \overline{\mathbf{N_{n1y}}} & = & \left( \begin{array}{c} \mathbf{M_{xn1x1}} \\ \mathbf{M_{wn1x1}} \\ \mathbf{M_{wn1y}} \\ \mathbf{M_{wn1y}} \end{array} \right) \\ \mathcal{J}_{\mathbf{fn1}}(\mathbf{v_{n1}}) & = & \left( \begin{array}{c} \mathbf{J}_{\nabla \mathbf{H_{n1}}} & 0 \\ 0 & \mathbf{J}_{\mathbf{z_{n1}}} \\ \end{array} \right) \\ \mathbf{I_{n1}} & = & \left( \begin{array}{c} \frac{\mathbf{I_{d}}}{\delta t} & 0 \\ 0 & \mathbf{I_{d}} \end{array} \right) \end{array} \right) \end{array}$$

Process

$$\begin{array}{lll} D_{1} & = & iD_{1}^{-1} \\ N_{1x1} & = & D_{1} \, \overline{N_{1x1}} \\ N_{1n1} & = & D_{1} \, \overline{N_{1n1}} \\ N_{1y} & = & D_{1} \, \overline{N_{1y}} \\ N_{n1x1} & = & \overline{N_{n1x1}} + \overline{N_{n11}} \, N_{1x1} \\ N_{n1x1} & = & \overline{N_{n1x1}} + \overline{N_{n11}} \, N_{1x1} \\ N_{n1y} & = & \overline{N_{n1y}} + \overline{N_{n11}} \, N_{1y} \\ c & = & N_{n1x1} \, x_{1} + N_{n1y} \, u \\ Iterate & : & F_{n1}(v_{n1}) = I_{n1} \, v_{n1} - N_{n1n1} \, f_{n1} - c \\ & & \mathcal{J}_{F_{n1}}(v_{n1}) = I_{n1} - N_{n1n1} \, \mathcal{J}_{fn1}(v_{n1}) \\ & & v_{n1} = v_{n1} - \mathcal{J}_{F_{n1}}^{-1}(v_{n1}) \, F_{n1}(v_{n1}) \\ v_{1} & = & N_{1x1} \, x_{1} + N_{1n1} \, f_{n1} + N_{1y} \, u \\ y & = & M_{yx1} \, \nabla H_{1} + M_{yxn1} \, \nabla H_{n1} M_{yy1} \, Z_{1} \, w_{1} + M_{yyn1} \, z_{n1} + M_{yy} \, u \\ x & = & x + \delta x \end{array} \label{eq:D1}$$

$$\mathbf{y} = \mathbf{M}_{yx1} \nabla \mathbf{H}_{1} + \mathbf{M}_{yxn1} \nabla \mathbf{H}_{n1} \mathbf{M}_{yw1} \mathbf{Z}_{1} \mathbf{w}_{1} + \mathbf{M}_{ywn1} \mathbf{z}_{n1} + \mathbf{M}_{yy} \mathbf{u}$$
(10)  
$$= \mathbf{M}_{yx1} \nabla \mathbf{H}_{1} + \mathbf{M}_{yxn1} \nabla \mathbf{H}_{n1} \mathbf{M}_{yw1} \mathbf{Z}_{1} \mathbf{w}_{1} + \mathbf{M}_{ywn1} \mathbf{z}_{n1} + \mathbf{M}_{yy} \mathbf{u}$$
(11)  
(12)

# 3.3 Realizability solver

### Connections

Serial 
$$\sum_{n=1}^{N} e_n = 0$$
,  $f_1 = \cdots = f_N = \phi$  (variable commune)

parallel  $\sum_{n=1}^{N} f_n = 0$ ,  $e_1 = \cdots = e_N = \phi$  (variable commune) Storage

## Realizable

$$\begin{cases} \phi = u = \frac{d}{dt}x = \frac{dx_1}{dt} = \dots = \frac{d}{dt}x_N \\ y = \nabla H_(x) = \sum_{i=1}^N \nabla H_i(x_i) \end{cases}$$

alors 
$$x = x_1 = x_2$$
 et  $H(x) = \left(\sum_{i=1}^{N} H_i\right)(x)$ 

## Non-Realizable

$$\begin{cases} \phi = u = \nabla \mathbf{H}_1(x) = \nabla \mathbf{H}_1(x_1) = \dots = \nabla \mathbf{H}_N(x_N), \\ y = \frac{\mathbf{d}}{\mathbf{d}t}x = \sum_{i=1}^N \frac{\mathbf{d}}{\mathbf{d}t}x_i \end{cases}$$

alors 
$$x = \sum_{i=1}^{N} x_i$$
 et  $\mathbf{H}(x) = \left(\sum_{i=1}^{N} H_i \nabla \mathbf{H}_i^{-1} G\right)(x)$  avec  $G^{-1}(x) = \sum_{i=1}^{N} \nabla \mathbf{H}_i(x_i)$  Dissipatives

### Realizable

$$\begin{cases} \phi = u = w &= w_1 = \dots = w_N, \\ y = z(w) &= \sum_{i=1}^{N} z_i(x_i) \end{cases}$$

# Non-Realizable

$$\begin{cases} \phi = u = z(w) = z_1(w_1) = \dots = z_N(w_N), \\ y = w = \sum_{i=1}^{N} w_i \end{cases}$$

alors 
$$w = \sum_{i=1}^{N} w_i = \left(\sum_{i=1}^{N} z_i^{-1}\right)(\phi)$$
 et  $z^{-1}(\phi) = w \Rightarrow z(w) = \left(\sum_{i=1}^{N} z_i\right)^{-1}(w)$ 

# 4 Connectors : Gyrators and Transformers

A connector is set of two edges  $(e_1, e_2)$ , the efforts and flux of which are coupled depending on the connector's type.

# 4.1 Gyrator

$$\begin{array}{rcl}
\mathfrak{e}_2 & = & \alpha \, \mathfrak{f}_1 \\
\mathfrak{e}_1 & = & -\alpha \, \mathfrak{f}_1
\end{array}$$

or equivalently:

$$\begin{array}{rcl}
\mathfrak{f}_2 & = & -\alpha^{-1} \, \mathfrak{e}_1 \\
\mathfrak{f}_1 & = & \alpha \, \mathfrak{e}_2
\end{array}$$

## 4.2 Transformer

$$\begin{array}{rcl}
\mathfrak{f}_2 & = & \alpha \, \mathfrak{f}_1 \\
\mathfrak{e}_1 & = & -\alpha \, \mathfrak{e}_2
\end{array}$$

or equivalently:

$$\begin{array}{rcl}
\mathfrak{f}_2 & = & -\alpha^{-1} \, \mathfrak{e}_1 \\
\mathfrak{f}_1 & = & \alpha \, \mathfrak{e}_2
\end{array}$$

# 5 Fractional calculus

The diffusive process in loudspeakers suspension (creep phenomenon ??) and loudspeakers ferromagnetic path (eddy current phenomenon ??) can be described by linear models that include fractional order dynamics (see [?, ?, ?, ?] for fractional modeling of viscoelasticity, and [?, ?, ?] for fractional modeling of eddy currents). A well established formalism for the realization of fractional transfer functions is the so called diffusive representations, recalled thereafter (see detailed developements in [?, ?], and [?] for a port-Hamiltonian formulation).

# 5.1 Fractional integrator

Defining  $s = \rho e^{i\theta}$ , with  $\rho \geq 0$  and  $\theta \in [-\pi, \pi[$ , the transfer function of the fractional integrator  $\mathcal{I}_{\beta}(s) = s^{-\beta}$  exhibits a cut  $\mathcal{C} = \mathbb{R}_{-}$ . The residue theorem gives the realization of  $\mathcal{I}_{\beta}$  as the continuous aggregation of linear damping along the cut  $\mathcal{C}$ . This leads to the following diffusive representation [?, §2]:

$$\mathcal{I}_{\beta}(s) : \mathbb{C} \setminus \mathbb{R}_{-} \to \mathbb{C}$$

$$s \mapsto \int_{0}^{\infty} \mu_{\beta}(\xi) \frac{1}{s+\xi} d\xi$$
(13)

where the weights  $\mu_{\beta}(\xi) = \frac{\mathcal{I}_{\beta}(-\xi-i0^+)-\mathcal{I}_{\beta}(-\xi+i0^+)}{2i\pi} = \frac{\sin(\beta\pi)}{\pi}\xi^{-\beta}$  correspond to the jump of  $\mathcal{I}_{\beta}$  across  $\mathcal{C} \equiv \{-\xi \in \mathbb{R}^-\}$ ). A state-space representation with output  $y_{\beta}(s) = \mathcal{I}_{\beta}(s)u_{\beta}(s)$  is:

$$\begin{cases} \frac{\mathrm{d}x_{\xi}}{\mathrm{d}t} &= -\xi x_{\xi} + u_{\beta}, \quad x_{\xi}(0) = 0, \\ y_{\beta} &= \int_{0}^{+\infty} \mu_{\beta}(\xi) x_{\xi} \mathrm{d}\xi. \end{cases}$$
(14)

The system (14) is recast as an infinite dimensional pH system (??), defining the hamiltonian density  $H_{\xi}(x_{\xi}) = \mu_{\beta}(\xi) \frac{x_{\xi}^2}{2}$  and the resistance density  $r_{\xi} = \frac{\xi}{\mu_{\beta}(\xi)}$ 

with  $z_{\xi}(w_{\xi}) = r_{\xi}w_{\xi}$ :

$$\begin{pmatrix} \frac{\mathrm{d}x_{\xi}}{\mathrm{d}t}w_{\xi} \\ y_{\beta} \end{pmatrix} = \begin{pmatrix} 0 & -1 & 1 \\ 1 & 0 & 0 \\ \mathbb{1}_{\infty} & 0 & 0 \end{pmatrix} \begin{pmatrix} \frac{\partial H_{\xi}}{\mathrm{d}x_{\xi}} \\ z_{\xi}(w_{\xi}) \\ u_{\beta} \end{pmatrix}$$
(15)

where  $\mathbb{1}_{\infty}$  denotes an infinite dimensional row vector of 1, that is  $y_{\beta} = \int_{0}^{\infty} \frac{\partial H_{\xi}}{\partial x_{\xi}} d\xi$ . Notice the total energy is  $H_{\beta}(\mathbf{x}_{\beta}) = \int_{\xi \in \mathcal{C}} H_{\xi}(x_{\xi}) d\xi$  with infinite dimensional state  $\mathbf{x}_{\beta} \in \mathbb{R}^{\mathbb{R}_{+}}$ . The realization of the dynamical element with parameter p and transfer function  $\mathcal{I}_{p,\beta}(s) = (ps^{\beta})^{-1}$  is given by (15), with  $\tilde{\mu}_{\beta,p}(\xi) = \frac{\mu_{\beta}(\xi)}{p}$ .

## 5.2 Fractional differentiator

Fractional damping can be modeled as combination of fractional integrators and differentiators (see [?, ?, ?, ?]). The realization of fractional differentiator of order  $\alpha$  with input  $u_{\alpha}$ , transfer function  $\mathcal{D}_{\alpha}(s) = s^{\alpha}$  and output  $y_{\alpha} = \mathcal{D}_{\alpha}u_{\alpha}$ , is built on the diffusive representation (14) as follows [?, ?]:

$$\begin{cases} \frac{\mathrm{d}x_{\xi}}{\mathrm{d}t} &= -\xi \cdot x_{\xi} + u_{\alpha}, \quad x_{\xi}(0) = 0, \\ y_{\alpha} &= \int_{0}^{+\infty} \mu_{1-\alpha}(\xi) (u_{\alpha} - \xi \cdot x_{\xi}) \mathrm{d}\xi. \end{cases}$$
(16)

Defining the hamiltonian density  $H_{\xi}(x_{\xi}) = \mu_{1-\alpha}(\xi)\xi\frac{x_{\xi}^2}{2}$ , the resistance density  $r_{\xi} = \mu_{1-\alpha}(\xi)$  and  $z_{\xi}(w_{\xi}) = r_{\xi}w_{\xi}$ , the pH formulation of the fractional differentiator (16) is

$$\begin{pmatrix} \frac{\mathrm{d}x_{\xi}}{\mathrm{d}t} \\ w_{\xi} \\ y_{\alpha} \end{pmatrix} = \begin{pmatrix} 0 & -\frac{1}{\mu_{1-\alpha}(\xi)} & 0 \\ \frac{1}{\mu_{1-\alpha}(\xi)} & 0 & -1 \\ 0 & -\mathbb{1}_{\infty} & 0 \end{pmatrix} \begin{pmatrix} \frac{\partial H_{\xi}}{\mathrm{d}x_{\xi}} \\ z_{\xi}(w_{\xi}) \\ u_{\alpha} \end{pmatrix}. \tag{17}$$

## 5.3 Finite order approximation

For implementation purpose, a finite approximation of diffusive representation (15) is built based on a finite set of  $n_{\xi}$  poles  $(\xi_1, \dots, \xi_{n_{\xi}})$  localized on the cut  $\mathcal{C}$ . The weights  $\boldsymbol{\mu} = (\mu_1 \cdots \mu_{n_{\xi}})^T$  are obtained from a least square optimization as detailed in [?, sec. 5.1.2], by minimizing an appropriate distance between  $\mathcal{I}_{\beta}$  and its discretisation  $\widehat{\mathcal{I}}_{\beta}$ :

$$\widehat{\mathcal{I}}_{\beta}(s) = \sum_{n=1}^{n_{\xi}} \frac{\mu_n}{s + \xi_n} = \mathbf{E}(s) \cdot \boldsymbol{\mu} \quad \text{with} \quad \mathbf{E}(s) = \left(\frac{1}{s + \xi_1} \cdots \frac{1}{s + \xi_{n_{\xi}}}\right)^{\mathsf{T}}.$$
 (18)

The poles  $\xi_n$ 's are chosen as  $\xi_n = 10^{\ell_n} \in \mathcal{C}$ , for  $0 \le n \le n_\xi + 1$ , where the  $\ell_n$ 's are equally spaced, with step  $\delta = \frac{\ell_{n_\xi + 1} - \ell_0}{n_\xi + 1}$ , from  $\ell_0$  to  $\ell_{N+1}$ . Since gain deviations are perceived relatively to the reference gains on the audio range, the weights  $\mu$  are optimized with respect to the objective function

$$\mathcal{O}(\boldsymbol{\mu}) = \int_{\omega_{-}}^{\omega_{+}} \left| 1 - \frac{\widehat{\mathcal{I}}_{\beta}(s = i\omega)}{\mathcal{I}_{\beta}(s = i\omega)} \right|^{2} d\ln \omega.$$
 (19)

where  $\omega_{-}=2\pi f_{-}$ ,  $\omega_{+}=2\pi f_{+}$  for  $[f_{-},f_{+}]=[20\mathrm{Hz},20\mathrm{kHz}]$ . In practice, the integral in (19) is approximated by a finite sum on a frequency grid, here,  $\ln \omega_{k} = \ln \omega_{-} + \frac{k}{n_{\omega}} \ln \frac{\omega_{+}}{\omega_{-}}$  for  $0 \le k \le n_{\omega}$ . This yields the following practical objective function

$$\widehat{\mathcal{O}}(\boldsymbol{\mu}) = \overline{(\mathbf{M}\boldsymbol{\mu} - \mathbf{T})}^{\mathsf{T}} \mathbf{W} (\mathbf{M}\boldsymbol{\mu} - \mathbf{T}), \tag{20}$$

where matrix  $\mathbf{M}$  is composed of the rows  $[\mathbf{M}]_{k,:} = \mathbf{E}(s = i\omega_{k-\frac{1}{2}})^T$  defined in (18), where  $\omega_{k-\frac{1}{2}} = \sqrt{\omega_{k-1}\omega_k}$  denotes the mean of  $\omega_{k-1}$  and  $\omega_k$  for  $1 \le k \le n_\omega$ . Vector  $\mathcal{I}$  is composed of  $[\mathcal{I}]_k = \mathcal{I}_\beta(s = i\omega_{k-\frac{1}{2}})$  and the diagonal matrix  $\mathbf{W}$  is defined by  $[\mathbf{W}]_{k,k} = (\ln \omega_k - \ln \omega_{k-1}) / |\mathcal{I}|_k|^2$ . The minimization of (20) is achieved by off-the-shelf optimization algorithm, imposing the weights to be positive:

$$\widehat{\boldsymbol{\mu}} = \{ \min_{\boldsymbol{\mu}} \widehat{\mathcal{O}}(\boldsymbol{\mu}) : \boldsymbol{\mu} > 0 \}$$
 (21)

The finite dimensional pH system realizing the weighted fractional integrator with transfer function  $\mathcal{I}_{p,\beta} = (ps^{\beta})^{-1}$  is given in table 1 with:

$$\begin{cases}
p_n &= \frac{\widehat{\mu}_n}{p}, \\
r_n &= \frac{\xi_n}{p_n},
\end{cases} \quad n \in (1, \dots n_{\xi}). \tag{22}$$

According to section 5.2, the finite dimensional approximation of the weighted

State: $\mathbf{x}_{\beta} = (x_1, \cdots, x_{n_{\xi}})^{T}$	Energy: $H_{\beta}(\mathbf{x}_{\beta}) = \frac{1}{2} \mathbf{x}_{\beta}^{T} \operatorname{diag}(p_1, \dots, p_{n_{\xi}}) \mathbf{x}_{\beta}$	
Dissipation variable: $\mathbf{w}_{\beta} = (w_1, \cdots, w_{n_{\xi}})^{T}$	Dissipation law: $\mathbf{z}_{\beta}(\mathbf{w}_{\beta}) = \operatorname{diag}(r_1, \cdots, r_{n_{\xi}}) \mathbf{w}_{\beta}$	
$egin{array}{c}  ext{Input}: \ u_eta \end{array}$	$egin{aligned}  ext{Output}: \ \widehat{y}_{eta} \end{aligned}$	
Structure: $\mathbf{J}_{xx} = \mathbb{O}_{n_{\xi} \times n_{\xi}}, \ \mathbf{J}_{xw} = -\mathbf{I}_{\mathbf{d}n_{\xi}}, \ \mathbf{J}_{xy} = \mathbb{1}_{n_{\xi} \times 1}, $ $\mathbf{J}_{ww} = \mathbb{O}_{n_{\xi} \times n_{\xi}}, \ \mathbf{J}_{wy} = \mathbb{O}_{n_{\xi} \times 1}, \ \mathbf{J}_{yy} = 0.$		

TABLE 1 – Port-Hamiltonian formulation (??) for the approximation of the fractional integrator  $y_{\beta}(s) = (ps^{\beta})^{-1}u_{\beta}(s)$  on a finite set of  $n_{\xi}$  poles. The parameters  $p_n, r_n$  for  $n \in (1, \dots, n_{\xi})$  are defined in (22) based on the minimization of (20). As an example, if  $u_{\beta} \equiv i$  and  $y_{\beta} \equiv v$ , this structure corresponds to the serial connection of  $n_{\xi}$  parallel RC cells; if  $y_{\beta} \equiv i$  and  $u_{\beta} \equiv v$ , this structure corresponds to the parallel connection of  $n_{\xi}$  serial LC cells.

fractional differentiator with transfer function  $\mathcal{D}_{\alpha,p} = ps^{\alpha}$  is obtained from the minimization of (20) for the transfer function  $\mathcal{I}_{1-\alpha}$ . The corresponding pH formulation is given in table 2 with:

$$\begin{cases}
p_n = p \widehat{\boldsymbol{\mu}}_n \xi_n, \\
r_n = p_n \widehat{\boldsymbol{\mu}}_n,
\end{cases} \qquad n \in (1, \dots n_{\xi}).$$
(23)

State: $\mathbf{x}_{\alpha} = (x_1, \cdots, x_{n_{\xi}})^{T}$	Energy: $H_{\alpha}(\mathbf{x}_{\alpha}) = \frac{1}{2} \mathbf{x}_{\alpha}^{T} \operatorname{diag}(p_1, \dots, p_{n_{\xi}}) \mathbf{x}_{\alpha}$	
Dissipation variable: $\mathbf{w}_{\alpha} = (w_1, \cdots, w_{n_{\xi}})^{T}$	Dissipation law: $\mathbf{z}_{\alpha}(\mathbf{w}_{\alpha}) = \operatorname{diag}(r_1, \cdots, r_{n_{\xi}}) \mathbf{w}_{\alpha}$	
$\begin{array}{c} \text{Input}: \\ u_{\alpha} \end{array}$	$egin{aligned}  ext{Output} : \ \widehat{y}_{lpha} \end{aligned}$	
Structure: $ \mathbf{J}_{\mathtt{xx}} = \mathbb{O}_{n_{\xi} \times n_{\xi}},  \mathbf{J}_{\mathtt{xw}} = -\operatorname{diag}(\widehat{\boldsymbol{\mu}})^{-1},  \mathbf{J}_{\mathtt{xy}} = \mathbb{O}_{n_{\xi} \times 1}, \\ \mathbf{J}_{\mathtt{ww}} = \mathbb{O}_{n_{\xi} \times n_{\xi}},  \mathbf{J}_{\mathtt{wy}} = -\mathbb{1}_{n_{\xi} \times 1},  \mathbf{J}_{\mathtt{yy}} = 0. $		

Table 2 – Port-Hamiltonian formulation (??) for the approximation of the fractional differentiator  $y_{\alpha}(s) = ps^{\alpha} u_{\alpha}(s)$  on a finite set of  $n_{\xi}$  poles. The parameters  $p_n, r_n$  for  $n \in (1, \dots, n_{\xi})$  are defined in (23) based on the minimization of (20) for the transfer function  $\mathcal{I}_{1-\alpha}$ . The interpretation is less intuitive than for the fractional integrator of table 1 due to the coefficients in  $\mathbf{J}_{xw}$  that involves transformers.