MODEL STRUCTURES AND FITTING CRITERIA FOR SYSTEM IDENTIFICATION WITH NEURAL NETWORKS

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1/15

Motivations

Recurrent Neural Networks are commonly used to model dynamical systems. However, they seldom exploit available a priori knowledge.

- We present tailor-made model structures for system identification
- We develop efficient algorithms to fit these model structures to data.

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Recurrent Neural Networks are commonly used to model dynamical systems. However, they seldom exploit available a priori knowledge.

In this work:

- We present tailor-made model structures for system identification with Neural Networks
- We develop efficient algorithms to fit these model structures to data.

Settings

The data-generating system S_o is assumed to have the discrete-time state-space representation:

$$x_{k+1} = f(x_k, u_k)$$
$$y_k^{\circ} = g(x_k)$$
$$y_k = y_k^{\circ} + e_k$$

Training dataset \mathcal{D} consisting of N input samples $U = \{u_0, u_1, \ldots, u_{N-1}\}$ and output samples $Y = \{y_0, y_1, \ldots, y_{N-1}\}$ available.

Objective: estimate a dynamical model of S_o .



A very generic neural model structure:

$$x_{k+1} = \mathcal{N}_f(x_k, u_k; \theta)$$
$$y_k = \mathcal{N}_g(x_k; \theta)$$

where \mathcal{N}_f , \mathcal{N}_g are feed-forward neural networks. Can be specialized:

Linear approximation available ⇒

$$x_{k+1} = Ax_k + Bu_k + \mathcal{N}_f(x_k, u_k; \theta)$$

$$y_k = Cx_k + \mathcal{N}_g(x_k, u_k; \theta)$$

ullet State fully observed \Rightarrow

$$x_{k+1} = \mathcal{N}_f(x_k, u_k; \ \theta)$$
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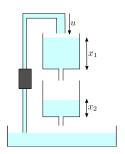
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Physics-inspired model structures

Two-tank system, input=flow u in upper tank, output=lower tank level x_2 .

- The system has two states: x_1 and x_2
- The state x_1 does not depend on x_2
- The state x₂ does not depend directly on u
- The state x_2 is observed



The observations above lead to the neural physics-inspired model structure

$$\dot{x}_1 = \mathcal{N}_1(x_1, u)$$

$$\dot{x}_2 = \mathcal{N}_2(x_1, x_2)$$

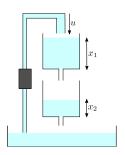
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In principle, simulation error minimization is a valid strategy:

$$\theta^o = \arg\min_{\theta, \hat{x}_0} \sum_{k=0}^{N-1} \|\hat{y}_k(\theta, \hat{x}_0) - y_k\|^2$$

where

$$\hat{x}_{k+1} = \mathcal{N}_f(\hat{x}_k, u_k; \ \theta)$$
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for
$$k = 0, 1, ..., N - 1$$
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However, it is not convenient from a computational perspective:

- Simulation is not parallelizable. Several neural network evaluations have to be performed sequentially.
- Back-propagation cost increases with the sequence length

In this work, we minimize instead the simulation error over batches of q subsequences, each one of length $m \ll N$.

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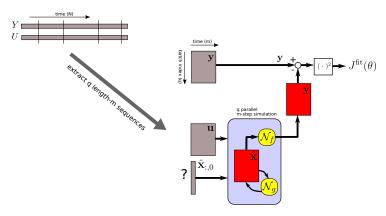
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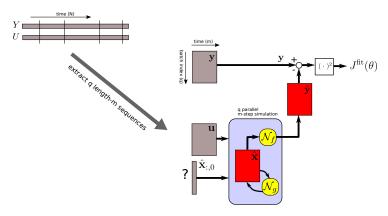
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for each iteration of gradient-based optimization:



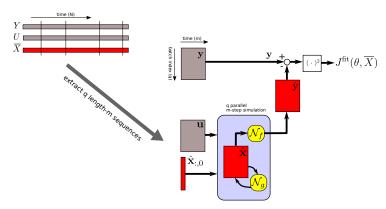
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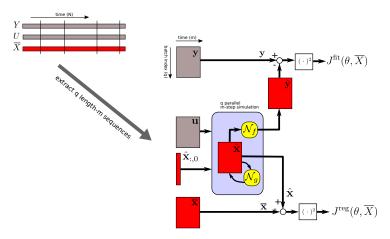
Problem: how do we choose $\hat{x}_{:,0}$, the initial state for each batch? We do not know it, but we need it to initialize all simulations.

We consider the unknown state sequence \overline{X} as an optimization variable. We sample from \overline{X} to obtain the initial state for simulation in each batch.



 $J^{ ext{fit}}$ is now a function of both θ and \overline{X} . We optimize w.r.t. both!

The hidden state sequence \overline{X} should also satisfy the identified dynamics! We enforce this by adding a regularization term in the cost function.



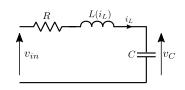
We minimize a weighted sum of J^{fit} and J^{reg} w.r.t. both θ and \overline{X} .

Simulation example

RLC circuit

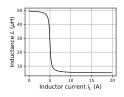
We consider a nonliner RLC circuit:

$$\begin{bmatrix} \dot{v}_{\textit{C}} \\ \dot{i}_{\textit{L}} \end{bmatrix} = \begin{bmatrix} 0 & \frac{1}{\textit{C}} \\ \frac{-1}{\textit{L}(i_{\textit{L}})} & \frac{-\textit{R}}{\textit{L}(i_{\textit{L}})} \end{bmatrix} \begin{bmatrix} v_{\textit{C}} \\ i_{\textit{L}} \end{bmatrix} + \begin{bmatrix} 0 \\ \frac{1}{\textit{L}(i_{\textit{L}})} \end{bmatrix} \textit{v}_{\textit{in}}$$



with nonlinear inductance $L(i_L)$

$$\textit{L(i}_{\textit{L}}) = \textit{L}_{\textit{0}} \bigg[\bigg(\frac{0.9}{\pi} \text{arctan} \big(-5 \big(|\textit{i}_{\textit{L}}| -5 \big) + 0.5 \bigg) + 0.1 \bigg]$$



Input: voltage v_{in} . Output: voltage v_C , current i_L . SNR=20

Neural model structure: fully observed state

$$x_{k+1} = \mathcal{N}_f(x_k, u_k; \theta)$$
$$y_k = x_k$$

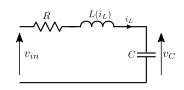


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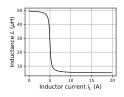
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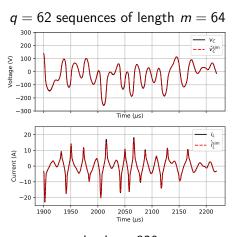
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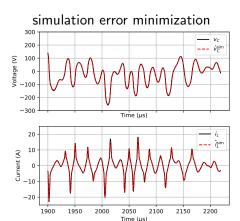


Numerical example

RLC circuit

Results in simulation on the test dataset. Training with:





train time: 320 s

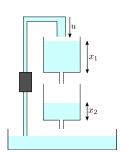
train time: 7000 s

Numerical example

Cascaded Tank System

Dataset with real measurements from www.nonlinearbenchmark.org





Neural model structure: physics-inspired

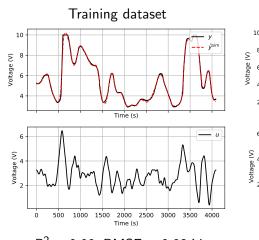
$$\dot{x}_1 = \mathcal{N}_1(x_1, u)
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y = x_2$$

The dependency of \mathcal{N}_2 on u models water overflow from upper tank.

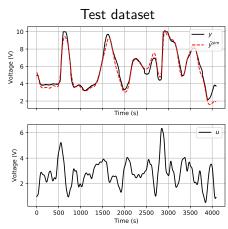
Numerical example

Cascaded Tank System

Training with m = 128, q = 64. Results on:



 $R^2 = 0.99$, RMSE = 0.08 V



 $R^2 = 0.97$, RMSE = 0.33 V

Conclusions

We have presented tailor-made neural structures for system identification embedding a priori knowledge.

We have shown how to parallelize the training using batches of short-size subsequences, and taking into account the effect of the initial condition.

Current/Future work

- Extension to the continuous-time setting
- Learning of Partial Differential Equations

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Thank you. Questions?

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