# MODEL STRUCTURES AND FITTING CRITERIA FOR SYSTEM IDENTIFICATION WITH NEURAL NETWORKS

Marco Forgione, Dario Piga

IDSIA Dalle Molle Institute for Artificial Intelligence SUPSI-USI, Lugano, Switzerland

14th IEEE International Conference on Application of Information and Communication Technologies

1/15

#### Motivations

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.

#### **Motivations**

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  $\Rightarrow$ 

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



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  $\Rightarrow$ 

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



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  $\Rightarrow$ 

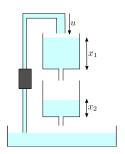
$$x_{k+1} = \mathcal{N}_f(x_k, u_k; \ \theta)$$
$$y_k = x_k$$



#### 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 x<sub>2</sub> does not depend directly on u
- The state  $x_2$  is observed



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

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

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

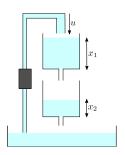
$$y = x_2$$



#### 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 x<sub>2</sub> does not depend directly on u
- The state  $x_2$  is observed



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

$$\dot{x}_1 = \mathcal{N}_1(x_1, u)$$
$$\dot{x}_2 = \mathcal{N}_2(x_1, x_2)$$
$$y = x_2$$



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

$$\hat{x}_{k+1} = \mathcal{N}_f(\hat{x}_k, u_k; \ \theta)$$
$$\hat{y}_k = \mathcal{N}_g(\hat{x}_k; \ \theta)$$

for 
$$k = 0, 1, ..., N - 1$$
.

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  $m \ll N$ .

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

$$\hat{x}_{k+1} = \mathcal{N}_f(\hat{x}_k, u_k; \ \theta)$$
$$\hat{y}_k = \mathcal{N}_g(\hat{x}_k; \ \theta)$$

for k = 0, 1, ..., N - 1.

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  $m \ll N$ .

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

$$\hat{x}_{k+1} = \mathcal{N}_f(\hat{x}_k, u_k; \ \theta)$$
$$\hat{y}_k = \mathcal{N}_g(\hat{x}_k; \ \theta)$$

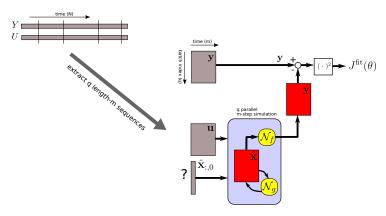
for k = 0, 1, ..., N - 1.

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

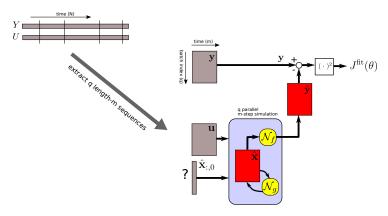
In this work, we minimize instead the simulation error over batches of q subsequences, each one of length  $m \ll N$ .

for each iteration of gradient-based optimization:



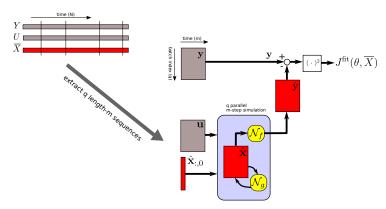
Problem: how do we choose  $\hat{x}_{:,0}$ , the initial state for each batch? We do not know it, but we need it to initialize all simulations.

for each iteration of gradient-based optimization:



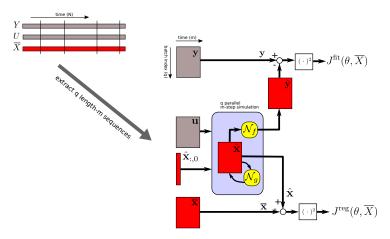
Problem: how do we choose  $\hat{x}_{:,0}$ , the initial state for each batch? We do not know it, but we need it to initialize all simulations.

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  $\theta$  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.



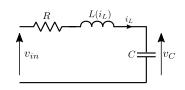
We minimize a weighted sum of  $J^{\mathrm{fit}}$  and  $J^{\mathrm{reg}}$  w.r.t. both  $\theta$  and  $\overline{X}$ .

## Simulation example

#### **RLC** circuit

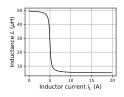
We consider a nonliner RLC circuit:

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



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

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

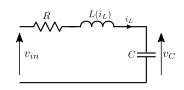


## Simulation example

#### RLC circuit

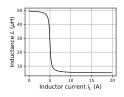
We consider a nonliner RLC circuit:

$$\begin{bmatrix} \dot{v}_C \\ \dot{i}_L \end{bmatrix} = \begin{bmatrix} 0 & \frac{1}{C} \\ \frac{-1}{L(i_L)} & \frac{-R}{L(i_L)} \end{bmatrix} \begin{bmatrix} v_C \\ \dot{i}_L \end{bmatrix} + \begin{bmatrix} 0 \\ \frac{1}{L(i_L)} \end{bmatrix} v_{in}$$



with nonlinear inductance  $L(i_L)$ 

$$\textit{L(i_L)} = \textit{L}_0 \bigg[ \bigg( \frac{0.9}{\pi} \text{arctan} \big( -5 \big( |\textit{i_L}| -5 \big) + 0.5 \bigg) + 0.1 \bigg]$$



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

Neural model structure: fully observed state

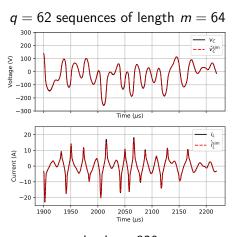
$$x_{k+1} = \mathcal{N}_f(x_k, u_k; \theta)$$
$$y_k = x_k$$

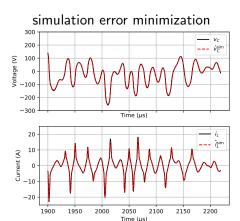


## Numerical example

#### RLC circuit

Results in simulation on the test dataset. Training with:





train time: 320 s

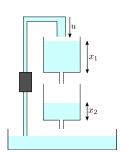
train time: 7000 s

## Numerical example

#### Cascaded Tank System

Dataset with real measurements from www.nonlinearbenchmark.org





Neural model structure: physics-inspired

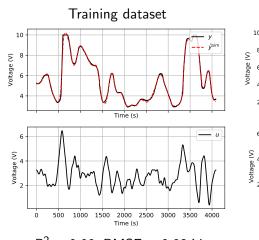
$$\dot{x}_1 = \mathcal{N}_1(x_1, u) 
\dot{x}_2 = \mathcal{N}_2(x_1, x_2, u) 
y = x_2$$

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

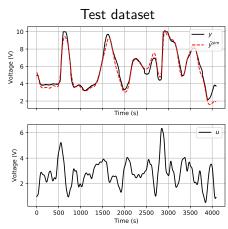
## Numerical example

#### 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

#### 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

# Thank you. Questions?

marco.forgione@idsia.ch