Latent Dynamic Networked System Identification with High-Dimensional Networked Data

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Abstract— Networked dynamic systems are ubiquitous in various domains, such as industrial processes, social networks, and biological systems. These systems produce high-dimensional data that reflect the complex interactions among the network nodes with rich sensor measurements. In this paper, we propose a novel algorithm for latent dynamic networked system identification that leverages the network structure and performs dimension reduction for each node via dynamic latent variables (DLVs). The algorithm assumes that the DLVs of each node have an auto-regressive model with exogenous input and interactions from other nodes. The DLVs of each node are extracted to capture the most predictable latent variables in the high dimensional data, while the residual factors are not predictable. The advantage of the proposed framework is demonstrated on an industrial process network for system identification and dynamic data analytics.

I. INTRODUCTION

Networked dynamic systems have become increasingly prevalent across various fields, from industrial processes to social networks and biological systems. These systems are composed of interconnected nodes with their own dynamics and dynamic interactions with other nodes [1], [2]. The nodes in the networked systems are often equipped with a rich set of sensors that yield high-dimensional time series data. Recent works have dealt with the estimation and system identification problems of these networked models with tremendous progress [3], [4], [5]. However, in dealing with high-dimensional networked systems data, little has been done to extract low-dimensional latent dynamic networks from the high-dimensional networked data [6].

To reduce dimension in data with static collinearity, statistical techniques such as principal component analysis (PCA), partial least squares (PLS), and canonical correlation analysis (CCA) are commonly used in data analytics [7], [8]. These methods effectively extract low-dimensional latent variables (LV) with statistical objectives. However, these methods do not focus on extracting latent dynamic information from data.

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To deal with the issue of co-moving dynamics or collinear dynamics in high-dimensional data, dynamic latent variables (DLVs) and dynamic factor models (DFMs) have been developed. Box and Tiao [9] introduced a technique to decompose multivariate time series into dynamic and static latent factors. Following this work, DFMs have been extensively used in financial time series to extract reduceddimensional dynamics [10]. Besides, compressive sensing offers another perspective for extracting reduced-dimensional dynamic LVs (DLVs) with principles of signal processing and optimization [11].

The aforementioned methods may not emphasize the predictability of the extracted DLVs. To this end, various reduced-dimensional DLVs methods with a focus on predictability have been proposed. For instance, the dynamic inner PCA (DiPCA) [12] algorithm, the dynamic inner PLS (DiPLS) [13] model, and the dynamic inner CCA (DiCCA) [14] method all extract the most predictable DLVs based on the univariate auto-regressive (AR) assumption. Recent works include the latent vector AR (LaVAR) model with fully interactive DLVs [15] and the latent state space representation with a CCA objective (LaSS-CCA) [16]. These methods also extract predictable DLVs in descending order and have been demonstrated to reduce the dimensionality of time-series data effectively. However, these methods do not consider the networked structure in the data, lacking the ability to capture the complex communications among the nodes of the networked system.

To address the challenges mentioned above, we propose a new framework for latent dynamic networked systems modeling with the following advantages:

- 1) To handle co-moving or collinear dynamics, low dimensional dynamics are captured on each individual node system to extract DLVs with a network topology;
- 2) The proposed framework extends single-node DLV methods to networked dynamic systems to analyze highdimensional networked time series data.
- 3) The proposed networked systems adopt vector autoregressive models with exogenous input (VARX) for each node, and dynamic connections to other nodes follow a fully interactive topology;
- 4) The identified network model is a networked latent VARX model (Net-LaVARX), which is readily suitable for networked system identification.

We demonstrate the effectiveness of the proposed framework by applying it to an industrial process network.

The remainder of the paper is organized as follows. In Section II, the proposed framework for latent dynamic networked systems is presented. The Net-LaVARX algorithm is developed in Section III with its learning procedures. In Section IV, a dynamic case study is detailed to verify the proposed method. The conclusions are given in Section V.

II. NETWORKED LATENT DYNAMIC SYSTEMS

A. Networked latent vector AR model with exogenous input (Net-LaVARX)

As described in the introduction section, high dimensional data usually have dynamics with a reduced dimension for each node of a networked system, giving rise to the need to develop networked latent dynamic modeling methods. For convenience, the notation is listed in Table I. Consider a networked dynamic system of M nodes, where Node $i \in [M]$ has a measurement vector $y_k^i \in \mathbb{R}^{p_i}$. Let $v_k^i \in \mathbb{R}^{\ell_i}$
represent the reduced dimensional dynamic latent vector represent the reduced dimensional dynamic latent vector (DLV) for Node i, whose elements lie independently in an ℓ_i -dimensional subspace. Following [15] for a single node system, the dynamics of y_k^i are related to v_k^i with an outer model,

$$
\mathbf{y}_k^i = \mathbf{P}_i \mathbf{v}_k^i + \bar{\mathbf{P}}_i \bar{\boldsymbol{\varepsilon}}_k^i,\tag{1}
$$

where $\bar{\varepsilon}_k^i \in \Re^{p_i-\ell_i}$ denotes the static noise, and $\mathbf{P}_i \in \Re^{p_i \times \ell_i}$ and $\overline{\mathbf{P}}_i \stackrel{\sim}{\in} \Re^{p_i \times (p_i - \ell_i)}$ are the respective loadings for the DLVs and static latent factors.

It is noted that the auto-correlated DLVs v_k^i and static noise $\bar{\varepsilon}_k^i$ complement each other to make up the whole signal series y_k^i . We give the following definition for a networked latent vector auto-regressive model with exogenous input.

Definition 1: A networked latent vector auto-regressive model with exogenous input (Net-LaVARX) system is described by (1) with

1) $[\mathbf{P}_i \quad \bar{\mathbf{P}}_i] \in \Re^{p_i \times p_i}$ form a non-singular matrix,

2) the DLVs v_k^i of Node *i* have auto-regressive terms with exogenous input as

$$
\boldsymbol{v}_{k}^{i} = \sum_{h=1}^{s_{i}} \left(\mathbf{A}_{h}^{i} \boldsymbol{v}_{k-h}^{i} + \mathbf{B}_{h}^{i} \boldsymbol{u}_{k-h}^{i} \right) + \sum_{j \in \mathcal{N}_{i}} \sum_{h=1}^{s_{i}} \mathbf{C}_{h}^{ij} \boldsymbol{v}_{k-h}^{j} + \varepsilon_{k}^{i},
$$
\n(2)

where the notation is given in Table I, and

Fig. 1: A diagram of a networked latent dynamic systems with three nodes.

3) the noise terms ε_k^i , $\bar{\varepsilon}_k^j$ are serially uncorrelated and mutually independent random sequences for $i \in \llbracket M \rrbracket$.

With the definition, if $\ell_i = p_i$, the model is a standard full-dimensional networked dynamic system, and $\bar{\varepsilon}_k^i$ is null. On the other hand, if $\ell_i = 0$, the model is composed of static nodes only. Without the cross terms \mathbf{C}_h^{ij} , the above model is a standard latent VARX model for Node i. The networked dynamic system is illustrated in Fig. 1, where dashed arrows represent dynamic interactions between nodes.

B. Extracting Net-LaVARX DLVs via oblique projections

The ultimate objective of the Net-LaVARX modeling is to develop algorithms to extract the DLVs v_k^i and estimate model parameters of (1) and (2). To do this, we design a projection to eliminate the column space of \overline{P}_i and focus on that of P_i . We give the following lemma to state the result.

Lemma 1: Let $\bar{P}_{i}^{\perp} \in \Re^{p_i \times \ell_i}$ have full column rank and $\bar{\mathbf{P}}_i^{\intercal} \bar{\mathbf{P}}_i^{\perp} = \mathbf{0}$. Then

$$
\boldsymbol{v}_k^i = \mathbf{R}_i^{\mathsf{T}} \boldsymbol{y}_k^i, \tag{3}
$$

where $\mathbf{R}_i = \bar{\mathbf{P}}_i^{\perp} (\mathbf{P}_i^{\mathsf{T}} \bar{\mathbf{P}}_i^{\perp})^{-1}$, $\mathbf{R}_i^{\mathsf{T}} \mathbf{P}_i = \mathbf{I}$, and $\mathbf{P}_i \mathbf{R}_i^{\mathsf{T}}$ i ^t is an oblique projection matrix, i.e., $(\mathbf{P}_i \mathbf{R}_i^{\mathsf{T}})$ $(\mathbf{F}_i)^2 = \mathbf{P}_i \mathbf{R}_i^{\intercal}$ $\frac{\intercal}{i}$.

The proof of this lemma is given in the appendix of the paper. This lemma shows that we can extract the DLVs via an oblique projection. Further, using (3), the projection

$$
\boldsymbol{y}_{k}^{i,\text{proj}} = \mathbf{P}_{i} \mathbf{R}_{i}^{\mathsf{T}} \boldsymbol{y}_{k}^{i} = \mathbf{P}_{i} \boldsymbol{v}_{k}^{i}
$$
(4)

is a reconstruction of y_k^i .

While separating the dynamic variations from static ones in (1) is feasible, P_i cannot be uniquely identified due to the bilinear decomposition in (1). Assuming v_k^{*i} is the true DLV with loadings \mathbf{P}_i^* , we have

$$
\boldsymbol{y}_k^i = \mathbf{P}_i^* \boldsymbol{v}_k^{*i} + \mathbf{\bar{P}}_i \bar{\boldsymbol{\varepsilon}}_k^i.
$$

For any nonsingular $\mathbf{M} \in \Re^{\ell_i \times \ell_i}$, we have

$$
\mathbf{P}_i^* \boldsymbol{v}_k^{*i} = (\mathbf{P}_i^*\mathbf{M}^{-1})(\mathbf{M}\boldsymbol{v}_k^{*i})
$$

It is obvious that $v_k^i = M v_k^{*i}$ is an equivalent dynamic latent vector. Therefore, the true loadings matrix can only

be identified up to the same range space, since $P_i^*M^{-1}$ and \mathbf{P}_{i}^{*} share the same range space. For this reason, the subspace of the DLVs is uniquely identified. There are extra degrees of freedom to make the DLVs in v_k^i in a pre-specified order, e.g., a descending order of predictability with constrained magnitudes [14], [15].

Substituting (3), (2) into (1) gives the model of measurement variables as

$$
\mathbf{y}_{k}^{i} = \sum_{h=1}^{s_{i}} (\mathbf{P}_{i} \mathbf{A}_{h}^{i} \mathbf{R}_{i}^{\mathsf{T}} \mathbf{y}_{k-h}^{i} + \mathbf{P}_{i} \mathbf{B}_{h}^{i} \mathbf{u}_{k-h}^{i})
$$

$$
+ \sum_{j \in \mathcal{N}_{i}} \sum_{h=1}^{s_{i}} \mathbf{P}_{i} \mathbf{C}_{h}^{ij} \mathbf{R}_{i}^{\mathsf{T}} \mathbf{y}_{k-h}^{j} + \mathbf{e}_{k}^{i}, \tag{5}
$$

which is a networked VARX model with reduced-rank autoregressive matrices $P_i A_h^i R_i^{\mathsf{T}}$ i^{T} , where

$$
\boldsymbol{e}^i_k = \mathbf{P}_i \boldsymbol{\varepsilon}^i_k + \mathbf{\bar{P}}_i \bar{\boldsymbol{\varepsilon}}^i_k
$$

is the equivalent noise sequence of the model.

Since $\varepsilon_k^i \in \Re^{\ell_i}$ and $e_k^i \in \Re^{p_i}$ are the random noises that are serially independent, the best predictions of v_k^i and y_k^i are, respectively

$$
\hat{\mathbf{v}}_k^i = \sum_{h=1}^{s_i} (\mathbf{A}_h^i \mathbf{v}_{k-h}^i + \mathbf{B}_h^i \mathbf{u}_{k-h}^i) + \sum_{j \in \mathcal{N}_i} \sum_{h=1}^{s_i} \mathbf{C}_h^{ij} \mathbf{v}_{k-h}^j, \tag{6}
$$

$$
\hat{\mathbf{y}}_k^i = \mathbf{P}_i \hat{\mathbf{v}}_k^i \tag{7}
$$

with the auto-regressive coefficient matrix $A_h^i \in \Re^{\ell_i \times \ell_i}$, the exogenous input coefficient matrix $\mathbf{B}_h^i \in \mathbb{R}^{\ell_i \times m_i}$, and the regression coefficient matrix $\mathbf{C}_h^{ij} \in \Re^{\ell_i \times \ell_j}$.

C. Net-LaVARX with a CCA objective

Taking account of the interactions between the networked nodes, we aim to find the projection weight matrices \mathbf{R}_i for $i \in [\![M]\!]$ that make the DLVs v_k^i the most predictable
based on the latent dynamics, with $\ell \geq n$, $\forall i \in [\![M]\!]$ for the based on the latent dynamics, with $\ell_i < p_i$, $\forall i \in [\![M]\!]$ for the dimension reduction purposes. Let $e = \max\{e_i, i \in [\![M]\!]$ dimension reduction purposes. Let $s = \max\{s_i, i \in \llbracket M \rrbracket\}$ and consider that we have $s + N$ samples of training data $\{\{\mathbf{y}_k^i, \mathbf{u}_k^i\}_{k=1}^{s+N}, i \in [\![M]\!] \}$. Similar to [15] for a single node
system, we use the lest N samples to form the following system, we use the last N samples to form the following Net-LaVARX CCA objective

$$
\min_{\{\mathbf{R}_i, \mathbf{A}_h^i, \mathbf{B}_h^i, \mathbf{C}_h^{ij}\}} J = \sum_{i=1}^M \sum_{k=s+1}^{s+N} \|\mathbf{v}_k^i - \hat{\mathbf{v}}_k^i\|^2
$$
(8)

$$
\text{s.t.} \quad \sum_{k=s+1}^{s+N} \mathbf{v}_k^i (\mathbf{v}_k^i)^\mathsf{T} = \mathbf{I}, \quad i \in [\![M]\!], \quad (9)
$$

where the first s samples are saved for initialization.

III. LEARNING PROCEDURE OF DYNAMIC LATENT NETWORK SYSTEMS

A. Net-LaVARX objective with training data

With $s + N$ samples of training data $\{\{y_k^i, u_k^i\}_{k=1}^{s+N}, i \in$ $[M]$, we denote the matrices of output and input data as

$$
\mathbf{Y}^{i} = \begin{bmatrix} \mathbf{y}_{1}^{i} & \mathbf{y}_{2}^{i} & \cdots & \mathbf{y}_{s+N}^{i} \end{bmatrix}^{\mathsf{T}} \in \mathbb{R}^{(s+N)\times p_{i}},
$$

$$
\mathbf{U}^{i} = \begin{bmatrix} \mathbf{u}_{1}^{i} & \mathbf{u}_{2}^{i} & \cdots & \mathbf{u}_{s+N}^{i} \end{bmatrix}^{\mathsf{T}} \in \mathbb{R}^{(s+N)\times m_{i}}.
$$
 (10)

The time-shifted matrices with N samples are formed as

$$
\mathbf{Y}_{j}^{i} = \begin{bmatrix} \mathbf{y}_{j+1}^{i} & \mathbf{y}_{j+2}^{i} & \cdots & \mathbf{y}_{j+N}^{i} \end{bmatrix}^{\mathsf{T}} \in \mathbb{R}^{N \times p_{i}}
$$

\n
$$
\mathbf{U}_{j}^{i} = \begin{bmatrix} \mathbf{u}_{j+1}^{i} & \mathbf{u}_{j+2}^{i} & \cdots & \mathbf{u}_{j+N}^{i} \end{bmatrix}^{\mathsf{T}} \in \mathbb{R}^{N \times m_{i}} \quad (11)
$$

\n
$$
\mathbf{V}_{j}^{i} = \begin{bmatrix} \mathbf{v}_{j+1}^{i} & \mathbf{v}_{j+2}^{i} & \cdots & \mathbf{v}_{j+N}^{i} \end{bmatrix}^{\mathsf{T}} \in \mathbb{R}^{N \times \ell_{i}},
$$

for $j = 0, 1, ..., s$, where $\mathbf{V}_j^i = \mathbf{Y}_j^i \mathbf{R}_i$. For the case of $j = s$, we have

$$
\mathbf{V}_s^i = \left[\begin{array}{cccc} \boldsymbol{v}_{s+1}^i & \boldsymbol{v}_{s+2}^i & \cdots & \boldsymbol{v}_{s+N}^i \end{array} \right]^\intercal \in \Re^{N \times \ell_i}
$$

The objective (8) is reformulated as

$$
\min_{\{\mathbf{R}_i, \mathbf{A}_h^i, \mathbf{B}_h^i, \mathbf{C}_h^{ij}\}} J = \sum_{i=1}^M \|\mathbf{V}_s^i - \hat{\mathbf{V}}_s^i\|_F^2
$$
(12)

s.t. $(\mathbf{V}_s^i)^\mathsf{T} \mathbf{V}_s^i = \mathbf{I}, \quad i \in [\![M]\!]$ (13)

where $\|\cdot\|_F$ denotes the Frobenius norm and $\hat{\mathbf{V}}_s^i$ depends on $\{\mathbf R_i, \mathbf A_h^i, \mathbf B_h^i, \mathbf C_h^{ij}\}\$ which can be obtained from (6) as

$$
\hat{\mathbf{V}}_{s}^{i} = \sum_{h=1}^{s_{i}} \left(\mathbf{V}_{s-h}^{i} \mathbf{A}_{h}^{i \mathsf{T}} + \mathbf{U}_{s-h}^{i} \mathbf{B}_{h}^{i \mathsf{T}} \right) + \sum_{j \in \mathcal{N}_{i}} \sum_{h=1}^{s_{i}} \mathbf{V}_{s-h}^{j} \mathbf{C}_{h}^{ij \mathsf{T}}
$$
\n
$$
= \mathbb{V}_{s}^{i} \mathbb{A}^{i} + \mathbb{U}_{s}^{i} \mathbb{B}^{i} + \sum_{j \in \mathcal{N}_{i}} \mathbb{V}_{s}^{j} \mathbb{C}^{ij}, \qquad (14)
$$

where the augmented matrices are given as follows

$$
\begin{aligned}\n\mathbb{V}_{s}^{j} &= \left[\begin{array}{cccc} \mathbf{V}_{s-1}^{j} & \mathbf{V}_{s-2}^{j} & \cdots & \mathbf{V}_{s-s_i}^{j} \end{array} \right], \quad j \in \llbracket M \rrbracket \\
\mathbb{U}_{s}^{i} &= \left[\begin{array}{cccc} \mathbf{U}_{s-1}^{i} & \mathbf{U}_{s-2}^{i} & \cdots & \mathbf{U}_{s-s_i}^{i} \end{array} \right], \\
\mathbb{A}^{i} &= \left[\begin{array}{cccc} \mathbf{A}_{1}^{i} & \mathbf{A}_{2}^{i} & \cdots & \mathbf{A}_{s_i}^{i} \end{array} \right]^{T}, \\
\mathbb{B}^{i} &= \left[\begin{array}{cccc} \mathbf{B}_{1}^{i} & \mathbf{B}_{2}^{i} & \cdots & \mathbf{B}_{s_i}^{i} \end{array} \right]^{T}, \\
\mathbb{C}^{ij} &= \left[\begin{array}{cccc} \mathbf{C}_{1}^{ij} & \mathbf{C}_{2}^{ij} & \cdots & \mathbf{C}_{s_i}^{ij} \end{array} \right]^{T}, \quad j \in \mathcal{N}_i\n\end{aligned}
$$

for $i \in M$. To further simplify the notations, we denote

$$
\mathbb{Z}^i = \begin{bmatrix} \mathbb{V}^i & \mathbb{U}^i & [\mathbb{V}^j]_{j \in \mathcal{N}_i} \end{bmatrix},
$$

where $[\mathbb{V}^j]_{j \in \mathcal{N}_i}$ is arranged row-wise with all \mathbb{V}^j for $j \in \mathcal{N}_i$. Similarly, we collect all DLV model parameters as

$$
\mathbb{Q}^i = \begin{bmatrix} \mathbb{A}^i \\ \mathbb{B}^i \\ [\mathbb{C}^{ij}]_{j \in \mathcal{N}_i} \end{bmatrix},
$$

where $[\mathbb{C}^{ij}]_{j \in \mathcal{N}_i}$ is arranged column-wise with all \mathbb{C}^{ij} for $j \in \mathcal{N}_i$. The objective (12) is rewritten as

$$
\min_{\{\mathbf{R}_i, \mathbb{Q}^i\}} \quad J = \sum_{i=1}^M \|\mathbf{V}_s^i - \hat{\mathbf{V}}_s^i\|_F^2 = \sum_{i=1}^M \|\mathbf{V}_s^i - \mathbb{Z}^i \mathbb{Q}^i\|_F^2
$$
\n
$$
\text{s.t.} \quad (\mathbf{V}_s^i)^\mathsf{T} \mathbf{V}_s^i = \mathbf{I}, \quad i \in [\![M]\!]. \tag{15}
$$

The above equality constraints guarantee that the objective is equivalent to the sum of canonical correlations of all nodes [15].

B. Solution of the Net-LaVARX models

Due to the bilinear decomposition of the Net-LaVARX models, an alternating iterative solution is expected, as shown in early work [15]. Below we give an iterative solution to update ${\bf R}_i\}_{i=1}^M$. Given ${\bf R}_i\}_{i=1}^M$, ${\bf \{\mathbb{V}_s^i\}}_{i=1}^M$ can be calculated. We have the following solution for the i -th component by minimizing (15) ,

$$
\mathbb{Q}^i = \left(\mathbb{Z}^i\right)^+ \mathbf{V}_s^i,\tag{16}
$$

where $(\cdot)^+$ refers to the Moore–Penrose pseudo-inverse. Then we have

$$
\hat{\mathbf{V}}_s^i = \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathbf{V}_s^i.
$$
 (17)

The objective (15) can be rewritten as M

$$
\min_{\{\mathbf{R}_i\}} \quad J = \sum_{i=1}^M \left\| \mathbf{V}_s^i - \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathbf{V}_s^i \right\|_F^2
$$
\n
$$
= \sum_{i=1}^M \left\| \mathbf{V}_s^i \right\|_F^2 - \sum_{i=1}^M \left\| \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathbf{V}_s^i \right\|_F^2
$$
\n
$$
= \sum_{i=1}^M \ell_i - \sum_{i=1}^M \left\| \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathbf{Y}_s^i \mathbf{R}_i \right\|_F^2
$$
\ns.t.
$$
(\mathbf{V}_s^i)^{\mathsf{T}} \mathbf{V}_s^i = \mathbf{I}_{\ell_i}, \quad i \in \llbracket M \rrbracket
$$

In the iterative scheme, \mathbb{Z}^i depends on the previous values of $\{R_i\}_{i=1}^M$. Therefore, the above objective is equivalent to maximizing the following M separate objectives,

$$
\max_{\mathbf{R}_i} J_i = \left\| \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathbf{Y}_s^i \mathbf{R}_i \right\|_F^2
$$
\n
$$
\text{s.t.} \quad (\mathbf{V}_s^i)^\mathsf{T} \mathbf{V}_s^i = \mathbf{R}_i^\mathsf{T} (\mathbf{Y}_s^i)^\mathsf{T} \mathbf{Y}_s^i \mathbf{R}_i = \mathbf{I}_{\ell_i} \tag{19}
$$

for $i \in [M]$. To make it easy to satisfy the equality constraints, we perform the following singular value decomposition (SVD)

$$
\mathbf{Y}_s^i = \mathcal{U}_s^i \mathcal{D}_s^i (\mathcal{V}_s^i)^\mathsf{T}
$$

where \mathcal{D}_s^i retains all non-zero singular values of \mathbf{Y}_s^i which makes $\mathcal{U}_s^i = \mathbf{Y}_s^i \mathcal{V}_s^i(\mathcal{D}_s^i)^{-1}$. The DLV score matrix is equivalently obtained by linearly combining \mathcal{U}_s^i as

$$
\mathbf{V}_s^i = \mathcal{U}_s^i \mathbf{W}_i = \mathbf{Y}_s^i \mathcal{V}_s^i (\mathcal{D}_s^i)^{-1} \mathbf{W}_i = \mathbf{Y}_s^i \mathbf{R}_i
$$

which makes

$$
\mathbf{R}_i = \mathcal{V}_s^i (\mathcal{D}_s^i)^{-1} \mathbf{W}_i \tag{20}
$$

Therefore, solving W_i is equivalent to solving for R_i .

With the above SVD the objective (19) is transformed into the following objective

$$
\begin{aligned}\n\max_{\mathbf{W}_{i}} \quad & J_{i} = \left\| \mathbb{Z}^{i}(\mathbb{Z}^{i})^{+} \mathcal{U}_{s}^{i} \mathbf{W}_{i} \right\|_{F}^{2} \\
&= \text{tr} \left\{ \mathbf{W}_{i}^{\mathsf{T}}(\mathcal{U}_{s}^{i})^{+} \mathbb{Z}^{i}(\mathbb{Z}^{i})^{+} \mathcal{U}_{s}^{i} \mathbf{W}_{i} \right\} \qquad (21) \\
\text{s.t.} \quad & (\mathbf{V}_{s}^{i})^{\mathsf{T}} \mathbf{V}_{s}^{i} = \mathbf{W}_{i}^{\mathsf{T}} \mathbf{W}_{i} = \mathbf{I}_{\ell_{i}}.\n\end{aligned}
$$

It is clear that the above objective with $\mathbf{W}_i^{\mathsf{T}} \mathbf{W}_i = \mathbf{I}_{\ell_i}$ is achieved only if \mathbf{W}_i are the eigenvectors of $(\mathcal{U}_s^i)^\top \mathbb{Z}^i (\mathbb{Z}^i)^\top \mathcal{U}_s^i$ corresponding to the ℓ_i maximum eigenvalues. Therefore, performing eigen-decomposition

$$
(\mathcal{U}_s^i)^{\mathsf{T}} \mathbb{Z}^i (\mathbb{Z}^i)^+ \mathcal{U}_s^i = \bar{\mathbf{W}}_i \Lambda_i \bar{\mathbf{W}}_i^{\mathsf{T}},\tag{22}
$$

the optimal solution is $\mathbf{W}_i = \bar{\mathbf{W}}_i(:, 1 : \ell_i)$ for $i \in [\![M]\!]$. The complete iterative Net-LaVARX algorithm is summarized in Algorithm 1. Note that in the algorithm, the DLV scores are scaled to unit variance as a post-processing step. The loadings matrix is solved by

$$
\mathbf{P}_{i} = \arg\min_{\mathbf{V}_{s}} \|\mathbf{Y}_{s}^{i} - \mathbf{V}_{s}^{i} \mathbf{P}_{i}^{\mathsf{T}}\|_{F}^{2}
$$

= $(\mathbf{Y}_{s}^{i})^{\mathsf{T}} \mathbf{V}_{s}^{i} = \mathcal{V}_{s}^{i} \mathcal{D}_{s}^{i} \mathbf{W}^{i}.$ (23)

Algorithm 1 Pseudo-code for Net-LaVARX-CCA

Input: Scale output Y^i and inputs U^i to zero mean and unit variance; specify the number of DLVs ℓ_i and order s_i for $i \in \llbracket M \rrbracket$;

Output: Weight matrices \mathbf{R}_i , loadings matrices \mathbf{P}_i , and VARX coefficients \mathbb{Q}^i , for $i \in [M];$
Perform SVD: $\mathbf{V}^i = \mathcal{U}^i \mathcal{D}^i (\mathcal{Y}^i)^\mathsf{T}$ Perform SVD: $\mathbf{Y}_s^i = U_s^i D_s^i (V_s^i)^T$, $i \in [M]$;

Initialize $\mathbf{W}^i - \mathbf{I}$ (c) $1 \cdot \ell$) where x_i is the Initialize $\mathbf{W}^i = \mathbf{I}_{r_i}(:, 1 : \ell_i)$, where r_i is the rank of \mathcal{D}_s^i ; Form $\mathbb{U}_s^i = \left[\begin{array}{cc} \mathbf{U}_{s-1}^i & \mathbf{U}_{s-2}^i & \cdots & \mathbf{U}_{s-s_i}^i \end{array} \right], i \in [M];$
while not converge it iter \leq mer iter do. while not converge $\&$ iter \lt max iter do for $i = 1 : M$ do Update *i*-th DLV node $V_s^i = U_s^i W^i$; Update the corresponding \mathbb{V}_s^i with \mathbf{V}_s^i ; Concatenate augmented DLVs and inputs for \mathbb{Z}_s^i ; Perform eigen-decomposition on $(U_s^i)^T \mathbb{Z}_s^i (\mathbb{Z}_s^i)^+ U_s^i$ and store the ℓ_i leading eigenvectors in \mathbf{W}^i ; end for Convergence detection on W^i ;

end while

Calculate $R_i =$ √ $\overline{N-1}\mathcal{V}_{s}^{i}(\mathcal{D}_{s}^{i})^{-1}\mathbf{W}^{i};$ Calculate DLV scores $V^i = Y^i R^i$; Calculate DLV scores $\mathbf{V} = \mathbf{r} \cdot \mathbf{R}$;
Calculate loading matrix $\mathbf{P}_i = \mathcal{V}_s^i \mathcal{D}_s^i \mathbf{W}^i / \sqrt{2}$ $\overline{N-1};$ Calculate VARX coefficients \mathbb{Q}^i , for $i \in [M]$.

IV. CASE STUDY

In this section, we present an application of the proposed net-LaVARX model on the Tennessee Eastman process (TEP) [17] with specially simulated dynamic data under decentralized control [18]. The goal is to demonstrate the effectiveness of the dynamic latent networked model in predicting the behavior of the TEP systems and to compare the variable partition strategies with different network partitioning of the process.

The current simulation represents an improved version of the original TEP problem [19]. We record data from 12 manipulated variables and 22 measured process variables. Unlike the heavily utilized simulated data with a 3-minute interval for steady-state modeling [20], the dataset used in this study is sampled every 36 seconds to reflect latent dynamics in the data. Nine manipulated variables, namely $XMV(1)-XMV(8)$ and $XMV(10)-XMV(11)$, are treated as

the exogenous inputs, while XMEAS(1)-XMEAS(22) are chosen as the process measurements.

The central units in TEP consist of a reactor, a stripper, and a vapor-liquid separator (Separ.) column with a condenser (Cond.) and a compressor (Comp.). By examining the process flow chart in [17] and decentralized control in [18], we respectively model the TEP process as networked dynamic systems with two partition scenarios, which are in Table II. Partition-I includes the node of the reactor with the condenser, the node of the separator with the compressor, and the node of the stripper. Partition-II moves the compressor to the node of the separator. For comparison, we also tested a model that treats all measurements and inputs as a whole, i.e., the LaVARX model, without considering the networked structure, denoted as *Monolithic*.

The Net-LaVARX models are built on each partition of the TEP process. For the Net-LaVARX model, the hyperparameters include the number of DLVs ℓ_i and the AR orders s_i . To select the optimal number of hyper-parameters, the grid search method is utilized, where the first 60% of the samples are used to train the model, and the next 15% is used for validation purposes. The last 25% of the data is reserved for testing the model generalization performance. The performance metrics adopted are the coefficient of determination $(R²)$, the correlations (Corr) coefficients, the root mean squared error (RMSE), and the mean absolute error (MAE) between the predicted and actual measurement values. The optimal number of hyper-parameters ℓ_i^* and s_i^* with performance metrics on the validation set are depicted in Table III. It is noticed that the sums of the optimal number of ℓ^* and s^* for Partition-I and Partition-II are comparable to that of the Monolithic (LaVARX) model, but their prediction performance is rather different. Precisely, the Net-LaVARX models with Partition-I and Partition-II attain the respective best RMSE and MAE, and the best R^2 and Corr on the validation set. It also verifies the feasibility of dividing the TEP into a networked system model with three nodes according to its central columns, and the effectiveness in improving the prediction performance.

Once the optimal ℓ_i^* and s_i^* are decided, we train the final Net-LaVARX models again using the entire training set (first 75% of the samples). Then, we evaluate the performance of the final model on the remaining 25% of the samples, with the same metrics given in Table IV. The results show that networked systems with both partitions outperform the performance of the monolithic model, as evidenced in Table III and Table IV.

To visualize the interactions between the latent nodes in these networked dynamic systems, the cross-correlations among the DLV scores of each node are calculated. With the threshold of 0.1 on the correlations, the DLVs network graph for Partition-II is generated and visualized in Fig. 2, where the nodes are marked with distinct colors. Each node contains the optimal number of DLVs. For easy visualization, the arcs from the leading two DLVs are colored, whereas the remaining ones are shown in light grey. The arcs indicate the connectivity between these DLVs, and the line width of each arc corresponds to the correlation magnitude between the connected DLVs. It is noticed that the thick arcs of different colors only appear between the leading 2 DLVs among these three nodes. This phenomenon coincides with our model objective that the predictable DLVs in each node are ranked in descending order. The leading DLVs of all nodes have consistently strong interactions among them, which dominate the underlying dynamics of the networked system. The minor DLVs have limited predictability and low correlations among them. As a particular case, DLV N3.5 in Fig. 2 of the stripper node has no significant correlations with other nodes.

In addition, the sizes of dots for the node and DLVs in Fig. 2 represent the corresponding R^2 values, which correspond to predictability. It can be seen that the leading DLVs tend to have larger sizes of dots, but this is not absolute. For instance, although N2.2 has a weak intragroup prediction ability, it has considerable connectivity with N1. The above observations shed light on interpreting latent dynamic networked systems for complex plants.

V. CONCLUSIONS

In this paper, a new framework for latent dynamic networked system modeling is successfully developed to extract low-dimensional network dynamics from high-dimensional networked data. The new latent networked models can handle co-moving or collinear dynamics with dimension reduction

TABLE IV: The performance metrics with the optimal hyper-

parameters on the test data.

Fig. 2: The graph visualization of the networked system nodes according to cross-correlations among the DLV scores in Partition-II. N1.x marked in blue is the reactor node; N2.x in red is the separator node; N3.x in green is the stripper node.

and take the process network topology into account. The Net-LaVARX models applied to the dynamic TEP dataset show that it is beneficial to partition complex plants into networked systems by employing the process topology. The results demonstrate the potential of the proposed latent dynamic networked modeling for system identification. Future work will focus on the computational complexity and robustness analysis of the networked system identification framework.

APPENDIX

A. Proof of Lemma 1

We first need to show that $(\bar{P}_i^{\perp})^{\dagger}P_i$ is invertible. Since $[\mathbf{P}_i \quad \bar{\mathbf{P}}_i]$ in (1) is non-singular,

$$
(\bar{\mathbf{P}}_i^\perp)^\intercal [\mathbf{P}_i \quad \bar{\mathbf{P}}_i] = [(\bar{\mathbf{P}}_i^\perp)^\intercal \mathbf{P}_i \quad \mathbf{0}_{\ell_i \times (p_i - \ell_i)}]
$$

has a rank of ℓ_i , i.e., $(\bar{\mathbf{P}}_i^{\perp})^{\intercal} \mathbf{P}_i$ has full rank. Therefore, $(\bar{\mathbf{P}}_i^{\perp})$ ^T \mathbf{P}_i is invertible.

It is straightforward to show that

$$
\mathbf{R}_i^{\mathsf{T}} \mathbf{P}_i = ((\bar{\mathbf{P}}_i^{\perp})^{\mathsf{T}} \mathbf{P}_i)^{-1} (\bar{\mathbf{P}}_i^{\perp})^{\mathsf{T}} \mathbf{P}_i = \mathbf{I}.
$$

It follows that

$$
(\mathbf{P}_i\mathbf{R}_i^{\mathsf{T}})^2 = \mathbf{P}_i(\mathbf{R}_i^{\mathsf{T}}\mathbf{P}_i)\mathbf{R}_i^{\mathsf{T}} = \mathbf{P}_i\mathbf{R}_i^{\mathsf{T}},
$$

which shows that $P_i R_i^{\mathsf{T}}$ \overline{i} is an oblique projection [21]. Moreover, it follows from $\bar{P}_i^{\mathsf{T}} \bar{P}_i^{\perp} = 0$ that

$$
\mathbf{R}_i^\mathsf{T} \bar{\mathbf{P}}_i = ((\bar{\mathbf{P}}_i^\perp)^\mathsf{T} \mathbf{P}_i)^{-1} (\bar{\mathbf{P}}_i^\perp)^\mathsf{T} \bar{\mathbf{P}}_i = \mathbf{0}.
$$

Thus, we have

$$
\mathbf{R}_i^{\intercal}\boldsymbol{y}_k^i = \mathbf{R}_i^{\intercal}\mathbf{P}_i\boldsymbol{v}_k^i + \mathbf{R}_i^{\intercal}\bar{\mathbf{P}}_i\bar{\boldsymbol{\varepsilon}}_k^i = \boldsymbol{v}_k^i + \boldsymbol{0} = \boldsymbol{v}_k^i.
$$

This completes the proof.

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