Identification of 2D Interconnected Systems: An efficient Steepest-Descent approach.

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Abstract: In this paper, a new identification method for large 2D grids of interconnected systems is presented. The proposed algorithm minimizes the Output-Error of the lifted system by using a Steepest-Descent optimization method which exploit the Multilevel Sequentially-Semi Separable (MSSS) structure of the involved matrices. Furthermore, it is shown that the computational complexity of the proposed approach is linear with respect to the number of subsystems in the grid. Finally, a numerical example is presented in order to show the effectiveness of the proposed algorithm.

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1. INTRODUCTION

In recent years, there has been an increasing interest in identification and modeling of interconnected systems. This has been motivated for the large number of applications where interconnected systems appear, such as: car platoons, Horowitz and Varaiya (2000), structural beams, Scholte and D’Andrea (2003), irrigation networks, Cantoni et al. (2007), turbulent wavefront reconstruction, Fraanje et al. (2010), satellite formations, Massioni (2010) between others. In the case of Prediction-Error methods (PEM), an identification approach for Linear Parameter Varying (LPV) spatially interconnected systems in input-output form based on Least Squares (LS) estimation is presented in Mukhtar et al. (2010). This approach has also been extended to Box-Jenkins models in Mukhtar et al. (2011). In Dankers (2015), the theory of closed-loop system identification is extended to networks of dynamical systems with known topology. Moreover, Weerts et al. (2016) studies the conditions for uniquely identify the topology of the network when some nodes are not contaminated by noise. One of the drawbacks of the methods mentioned before is the computational complexity. Therefore, new tools need to be developed for the identification of large networks of interconnected systems with general topology. In the same line, Torres et al. (2015) proposes an efficient Output-Error identification method for 1D interconnected systems. The method exploits Sequentially-Semi Separable (SSS) matrix operations to speed up the optimization routines. Moreover, in Yu and Verhaegen (2017) a subspace identification approach for 1D heterogeneous interconnected systems that exploit local input-output data is presented. Even though these methods are efficient in terms of computational complexity they focus in 1D spatially interconnected systems and not in complex networks.

In order to cope with this limitation, in this paper, an efficient Output-Error identification method for 2D grids of interconnected systems is introduced. The algorithm exploits the Multilevel Sequentially Semi Separable (MSSS) structure, Qiu et al. (2015), of the matrices involved to perform an efficient Steepest-Descent optimization in linear complexity with respect to the number of subsystems in the grid. This is achieved by using basic operations with MSSS matrices. The main contribution of this paper is to show that all the matrices involved in the computations (including partial derivatives) have MSSS structure.

In this context, the outline of our paper is as follows: In Section 2, the problem formulation and the statement of the identification problem are presented. The complete algorithm is exposed in Section 3 while in Section 4 an example is provided in order to show the potential of the proposed algorithm. Finally, we draw our conclusions.
We consider a 2D grid of $M \times N$ interconnected systems as depicted in Fig. 1. Every sub-system is described by the following set of equations:

$$\Sigma_{i,j} : \begin{bmatrix} x_k^{(i,j)} \\ y_k^{(i,j)} \\ z_k^{(i,j)} \\ \eta_k^{(i,j)} \\ S_{k}^{(i,j)} \end{bmatrix} = \begin{bmatrix} A_{i,j} & B_{i,j} & C_{i,j} & D_{i,j} & \eta_k^{(i,j)} \\ C_{i,j} & D_{i,j} & W_{i,j}^{(i,j)} & 0 & 0 \\ C_{i,j} & D_{i,j} & W_{i,j}^{(i,j)} & 0 & 0 \\ C_{i,j} & D_{i,j} & W_{i,j}^{(i,j)} & 0 & 0 \\ S_{k}^{(i,j)} & \end{bmatrix} \begin{bmatrix} x_k^{(i,j)} \\ y_k^{(i,j)} \\ z_k^{(i,j)} \\ \eta_k^{(i,j)} \\ S_{k}^{(i,j)} \end{bmatrix},$$

where $x_k^{(i,j)} \in \mathbb{R}^n$ is the state vector, $w_k^{(i,j)} \in \mathbb{R}^r$ is the input vector, $y_k^{(i,j)} \in \mathbb{R}^q$ is the output vector, $\eta_k^{(i,j)} \in \mathbb{R}^q$, $w_k^{(i,j)} \in \mathbb{R}^q$, and $S_k^{(i,j)} \in \mathbb{R}^q$ are the interconnection variables. The dimensions of the matrices in (1) are consistent with the sizes of the corresponding vectors.

If we consider $N \times M$ sub-systems with the following boundary conditions: $c_k^{(0,j)} = 0$, $w_k^{(N+1,j)} = 0$, $\eta_k^{(0,0)} = 0$, and $S_k^{(i,M+1)} = 0$, for all $i = 1, 2, \ldots, M$, we can model the complete 2D grid by performing consecutive substitutions in (1):

$$\Sigma : \begin{bmatrix} \bar{x}_{k+1}^{(i,j)} \\ \bar{y}_k^{(i,j)} \end{bmatrix} = \begin{bmatrix} \bar{A} & \bar{B} \\ \bar{C} & \bar{D} \end{bmatrix} \begin{bmatrix} \bar{x}_k^{(i,j)} \\ \bar{y}_k^{(i,j)} \end{bmatrix},$$

where $\bar{x}_k^{(i,j)} \in \mathbb{R}^{NMn}$ is the interconnection state vector that contains all the local states stacked together. Similar definitions hold for $\bar{x}_{k+1}^{(i,j)}, \bar{y}_k^{(i,j)}, \bar{A}, \bar{B}, \bar{C},$ and $\bar{D}$. On the other hand, $\bar{A} \in \mathbb{R}^{NMn \times NMn}$, $\bar{B} \in \mathbb{R}^{NMn \times NMr}$, $\bar{C} \in \mathbb{R}^{NMn \times NMn}$, and $\bar{D} \in \mathbb{R}^{NMn \times NMr}$ are Multilevel Sequentially Semi-Separable (MSSS) matrices

characterized by the special structure depicted in Fig. 2.b). Notice from Fig. 2.b) that a MSSS matrix is a Sequentially Semi-Separable (SSS) matrix whose generators are SSS matrices themselves (see eq. (4)-(10)). Moreover, an SSS matrix has the structure depicted in Fig. 2.a). This matrix can be obtained by sequentially multiplying the generators appearing in eq. (4) as shown in Fig. 2.a) and by placing the matrices $A_{i,j}$ in the main diagonal.

$$\bar{A} = MSSS(B_N^{(i)}, W_N^{(i)}, C_N^{(i)}, A^{(i)}, B_s^{(i)}, W_s^{(i)}, C_s^{(i)}), \quad (3)$$

$$\bar{A}^{(i)} = SSS(B^{(i)}_c, W^{(i)}_c, C^{(i)}_c, A^{(i)}, B^{(i)}_w, W^{(i)}_w, C^{(i)}_w), \quad (4)$$

$$\bar{B}^{(i)}_N = SSS(0, 0, 0, B^{(i)}_N, 0, 0, 0), \quad (5)$$

$$\bar{W}^{(i)}_N = SSS(0, 0, 0, W^{(i)}_N, 0, 0, 0), \quad (6)$$

$$\bar{C}^{(i)}_N = SSS(0, 0, 0, C^{(i)}_N, 0, 0, 0), \quad (7)$$

$$\bar{B}^{(i)}_s = SSS(0, 0, 0, B^{(i)}_s, 0, 0, 0), \quad (8)$$

$$\bar{W}^{(i)}_s = SSS(0, 0, 0, W^{(i)}_s, 0, 0, 0), \quad (9)$$

$$\bar{C}^{(i)}_s = SSS(0, 0, 0, C^{(i)}_s, 0, 0, 0), \quad (10)$$

Therefore, for the rest of the lifted matrices we have a similar definition but exchanging the corresponding generators. Thus for obtaining $\bar{B}$ we replace matrix $A$ by $B$ and $C$ by $D$ in eq. (3)-(10). For $\bar{C}$ we replace matrix $A$ by $C$ and $B$ by $F$ and finally for $\bar{D}$ we replace matrices $A$ and $C$ by $D$ and $B$ by $F$.

Therefore, the identification problem can now be formulated as follows: Given the input-output sequences:

$$\{\bar{u}_k^{(i,j)}, \bar{y}_k^{(i,j)}\}_{k=1}^{(i,M+1),j=1}^{N},$$

estimate the distributed system matrices $A_{i,j}$, $B_{i,j}$, $C_{i,j}$, $D_{i,j}$, $B^{(i)}_c$, $W^{(i)}_c$, $C^{(i)}_c$, $D^{(i)}_c$, $B^{(i)}_w$, $W^{(i)}_w$, $C^{(i)}_w$, $D^{(i)}_w$, $F^{(i)}_e$, $F^{(i)}_w$, $F^{(i)}_s$, and $F^{(i)}_c$. 

Without loss of generality, we assume the subsystems have the same number of states, inputs, outputs and interconnection variables.
leads to the derivation of the following objective function:

\[
J = \left( K - \frac{\partial \theta}{\partial y} \right)^T J K \left( \frac{\partial \theta}{\partial y} \right) \quad \forall k \in \{1, \ldots, N\}
\]

and the adjoint state-space equation given by:

\[
\hat{X}_{k-1}(\theta) = \hat{A} \hat{X}_k(\theta) + C^T(\theta) \epsilon_k(\theta)
\]

where \( \hat{X}_k(\theta) \) and \( \epsilon_k(\theta) \) are obtained by simulating the model given in (2). Notice that in this case two simulations are needed: the backward simulation (18) and the forward simulation of (2). For the derivation of the equations presented above the interested reader is referred to Verhaegen and Verdult (2007).

The main contribution of this paper is to show that all the matrices involved in the computations have MSSS structure (denoted by \( \Phi \)). Furthermore, the partial derivatives involved in the formulas can also be written as MSSS matrices:

\[
\frac{\partial}{\partial \theta} [I_p \otimes \Phi_k(\theta)] = \text{MSSS}(B^{(i,j)}_N, W^{(i,j)}_N, \mathcal{C}^{(i,j)}, T^{(i,j)}_s, W^{(i,j)}_s, \mathcal{C}^{(i,j)}_s),
\]

\[
\frac{\partial}{\partial \theta} [I_p \otimes \Phi_k(\theta)] = \text{MSSS}(B^{(i,j)}_N, W^{(i,j)}_N, \mathcal{C}^{(i,j)}, T^{(i,j)}_s, W^{(i,j)}_s, \mathcal{C}^{(i,j)}_s),
\]

\[
\frac{\partial}{\partial \theta} [I_p \otimes \Phi_k(\theta)] = \text{MSSS}(F^{(i,j)}_r, \mathcal{C}^{(i,j)}, \mathcal{C}^{(i,j)}_s, T^{(i,j)}_s, W^{(i,j)}_s, \mathcal{C}^{(i,j)}_s),
\]

\[
\frac{\partial}{\partial \theta} [I_p \otimes \Phi_k(\theta)] = \text{MSSS}(F^{(i,j)}_r, \mathcal{C}^{(i,j)}, \mathcal{C}^{(i,j)}_s, T^{(i,j)}_s, W^{(i,j)}_s, \mathcal{C}^{(i,j)}_s),
\]

\[
\forall i \in \{1, \ldots, M\}
\]

where the generator matrices in caligraphic style are depicted in Fig. 3. The corresponding matrix recurrences \( \gamma \) and \( \delta \) used to compute the generators are given in Appendix A.

### 3.2 Computational Complexity

According to the exposed before, matrix-vector product and matrix transpose MSSS operations are required for the two simulations and the computation of the Jacobian in every iteration. Furthermore, both operations can be performed in linear complexity with respect to the number of subsystems in the grid \( O(NM) \). (Houtzager and Rice (2009)). Moreover, the identification problem described above requires \( p = ((n^2 + nr + 8nq + rn + rl + 4ql + 4q^2)NM \) parameters to be estimated. This is clearly much better than the complexity of the full parameterization of the state space model (2), that produces \( (n^2 + nr + n\ell + nr + q\ell + qr + 4q^2)NM \) parameters. (assuming that \( n, r, \ell, q << N, M \)).

In addition, the number of parameters to be stored for the partial derivatives also increase linearly with respect to the number of sub-systems in the grid: \( (2n^2 + 2n\ell + 12q)(n^2 + nr + 8nq + rn + rl + 4qr + 4ql + 4q^2)NM \) (see Fig. 3). In the case of Quasi-Newton methods, the implementation also
Fig. 3. Generator matrices used to compute the partial derivatives ($\otimes$ is the Kronecker product)

results in linear complexity with respect to the number of subsystems in the grid as the approximation of the Hessian can also be represented as a SSS matrix. In addition, faster convergence to the optimum should be expected.

4. SIMULATION EXAMPLE

The proposed identification algorithm is tested on the heat equation discretized in a 2D domain with Dirichlet boundary conditions:

\[
\begin{bmatrix}
\begin{bmatrix}
\begin{bmatrix}
A & B & C \\
D & E & F \\
G & H & I
\end{bmatrix}
\end{bmatrix}
\end{bmatrix}^T
\]

where 
\[
A = 0.006(\Gamma_{i,j} + \Gamma_{0,i-1,j}),
\]
\[
b = 0.006(\Gamma_{i,j} + \Gamma_{i+1,j}),
\]
\[
c = 0.006(\Gamma_{i,j} + \Gamma_{i-1,j}),
\]
\[
d = 0.006(\Gamma_{0,i,j} + \Gamma_{i+1,j}),
\]

\[
e = (c_{i,j} + d_{i,j}) - (c_{i,j} + f_{i,j}) - 0.5,
\]
\[
f = 0.3 \times 1 - 0.06 \text{ and } \Gamma_{i,j} \sim \mathcal{U}(0,5) \text{ are uniformly distributed parameters. The input signals are normally distributed random variables with covariance matrix equal to } \text{cov}(u_k) = I_k. \text{ In addition, a Gaussian noise is added to the outputs with a prescribed Signal-to-Noise Ratio (SNR).}

In the following subsections, the prediction capabilities of the identified model, the computational complexity of the algorithm and the effect of the noise level on the prediction error are investigated.

4.1 Variance Accounted For as function of the number of iterations

In this case, the Variance-Accounted-For (VAF)\(^2\) is studied as function of the number of iterations. A 2D grid of \(15 \times 15 = 225\) sub-systems is considered. A set of 20 experiments is performed where every trial is a random realization of the parameters of every sub-system in the grid. In addition, the outputs are corrupted with a Gaussian noise of SNR\(_{dB} = 40dB\). The mean of the VAF is calculated over 20 trials and depicted in Fig. A.1. It is observed that the VAF converges to the 100% as the number of iterations increases.

4.2 Computational Complexity

In this section, the computation time per iteration of the proposed algorithm is studied. The computation time is recorded for 2D grids of increasing number of sub-systems and for 20 Monte-Carlo simulations per each grid. The results are presented in Fig. A.2.

In fact, it is observed that the computation time increases linearly as the number of sub-systems grows. This result confirms our hypothesis that the computations are performed on linear time with respect to the number of sub-systems in the grid \(\mathcal{O}(NM)\).

\(^{2}\) the VAF is defined as: \(VAF = 100 \cdot \max\{0, (1 - \frac{\text{var}(\hat{y}_k - y_k)}{\text{var}(y_k)})\}\)\%
4.3 The effect of noise

Finally, the effect of noise is investigated on the VAF index. The outputs are contaminated with Gaussian noise of different SNR characteristics and with different number of samples. Moreover, different realization of the subsystems in the grid along 20 Monte-Carlo simulations are considered. The results for two different data sets are depicted in Fig. A.3: A) noisy data used for training the model and B) noise-free data used for validation.

In this scenario, two effects can be observed. If the number of samples remain constant at \( K = 30 \), the VAF diminishes when the power of the noise increases (SNR decreases). This means that the prediction performance is deteriorated when the influence of noise is predominant. On the other hand, if the noise power is set constant and high at \( \text{SNR}_{IB} = 10dB \), the prediction capabilities are enhanced (VAF is increased) if the amount of samples used on the identification is increased (\( K = 90 \)), i.e., the use of longer sequences in the training procedure attenuates the noise overfit improving the results on the validation data.

5. CONCLUSION

In this paper, an efficient Output-Error identification method for 2D interconnected systems based on the Steepest-Descent gradient method was presented. It was shown that for 2D interconnected systems the whole algorithm can be expressed within the MSSS-approach which resulted in linear time computations with respect to the number of sub-systems in the grid. In addition, the results were strengthened by a simulation example showing the linearity and effectiveness of the proposed method. Convergence of the VAF to the 100% in finite number of iterations was exposed as well. Even though the prediction capabilities of the resulting model were deteriorated by output data contaminated with increasing noise power (overfit), this negative effects can be attenuated by considering data sequences with larger number of samples.

Appendix A

A.1 Matrices \( \gamma \)

Given the matrices \( W_N \in \mathbb{R}^{q \times q} \) and \( C_N \in \mathbb{R}^{q \times n} \) we define the matrix sequence \( \gamma_N^{(i,j)} \) as follows:

\[
\gamma_N^{(i,j)} = \gamma_N^{(i-1,j)} W_N^{(i-1,j)T} + x_N^{T(i-1,j)} C_N^{(i-1,j)T},
\]

\[
\gamma_N^{(1,j)} = 0, \quad \forall j \in \{1, \ldots, N\}
\]

On the other hand, given \( W_s \in \mathbb{R}^{q \times q} \) and \( C_s \in \mathbb{R}^{q \times n} \) we have:

\[
\gamma_s^{(i,j)} = \gamma_s^{(i,j)} W_s^{(i,j)T} + x_s^{T(i,j)} C_s^{(i,j)T},
\]

\[
\gamma_s^{(M,j)} = 0, \quad \forall j \in \{1, \ldots, N\}
\]

Finally, given \( W_c \in \mathbb{R}^{q \times q} \), \( W_c \in \mathbb{R}^{q \times q} \), \( C_c \in \mathbb{R}^{q \times n} \) and \( C_c \in \mathbb{R}^{q \times n} \) we define:

\[
\gamma_c^{(i,j)} = \gamma_c^{(i,j-1)} W_c^{(i,j-1)T} + x_c^{T(i,j)} C_c^{(i,j-1)T},
\]

\[
\gamma_c^{(1,1)} = 0, \quad \forall i \in \{1, \ldots, M\}
\]
and:

\[ \gamma_w^{(i,j-1)} = \gamma_w^{(i,j)} W_w^{(i,j)} + x_{(i,j)}^T C_w^{(i,j)} T, \]

\[ \nexists i \in \{1, \cdots, M\}, \forall j \in \{N, \cdots, 2\} \]

\[ \gamma_w^{(i,N)} = 0. \]

\[ \forall i \in \{1, \cdots, M\} \]

**A.2 Matrices δ**

Moreover, given the matrices \( W_N \in \mathbb{R}^{q \times q} \) and \( D_N \in \mathbb{R}^{q \times r} \)
we define the matrix sequence \( \delta_N^{(i,j)} \) as follows:

\[ \delta_N^{(i,j)} = \delta_N^{(i-1,j)} W_N^{(i-1,j)} + u_{(i-1,j)}^T D_N^{(i-1,j)} T, \]

\[ \nexists i \in \{2, \cdots, M\}, \forall j \in \{1, \cdots, N\} \]

\[ \delta_N^{(M,j)} = 0. \]

\[ \forall j \in \{1, \cdots, N\} \]

On the other hand, given \( W_s \in \mathbb{R}^{q \times q} \) and \( D_s \in \mathbb{R}^{q \times r} \)
we have:

\[ \delta_s^{(i,j-1)} = \delta_s^{(i,j)} W_s^{(i,j)} + u_{(i,j)}^T D_s^{(i,j)} T, \]

\[ \nexists i \in \{M, \cdots, 2\}, \forall j \in \{1, \cdots, N\} \]

\[ \delta_s^{(M,j)} = 0. \]

\[ \forall j \in \{1, \cdots, N\} \]

Finally, given \( W_c \in \mathbb{R}^{q \times q} \), \( W_w \in \mathbb{R}^{q \times q} \), \( D_c \in \mathbb{R}^{q \times r} \)
and \( D_w \in \mathbb{R}^{q \times r} \) we define:

\[ \delta_c^{(i,j)} = \delta_c^{(i,j-1)} W_c^{(i,j-1)} + u_{(i,j-1)}^T D_c^{(i,j-1)} T, \]

\[ \nexists i \in \{1, \cdots, M\}, \forall j \in \{2, \cdots, N\} \]

\[ \delta_c^{(i,1)} = 0. \]

\[ \forall i \in \{1, \cdots, M\} \]

and:

\[ \delta_w^{(i,j-1)} = \delta_w^{(i,j)} W_w^{(i,j)} + u_{(i,j)}^T D_w^{(i,j)} T, \]

\[ \nexists i \in \{1, \cdots, M\}, \forall j \in \{N, \cdots, 2\} \]

\[ \delta_w^{(i,N)} = 0. \]

\[ \forall i \in \{1, \cdots, M\} \]

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