

# Identification of linear dynamic networks with reduced-rank noise

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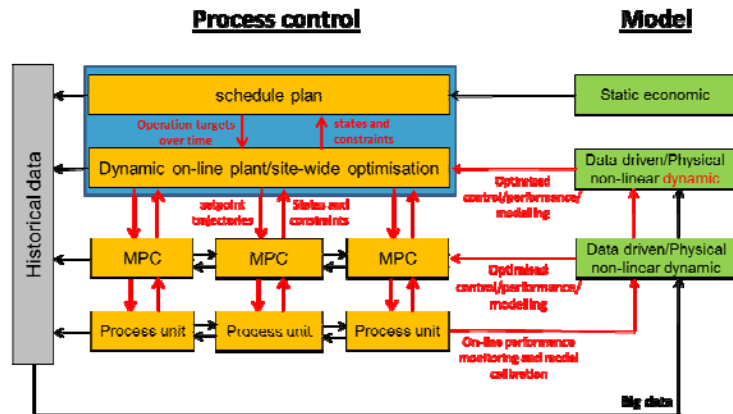
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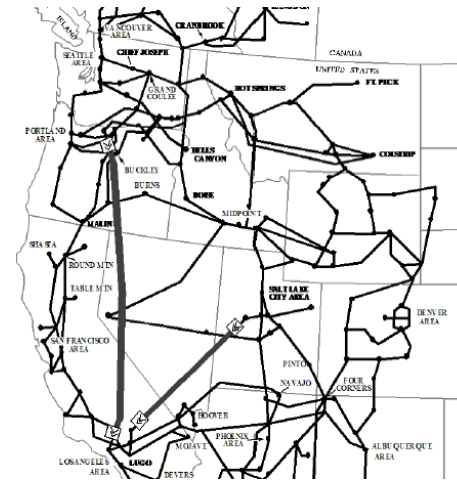
**Where innovation starts**

# Introduction – dynamic networks

## Decentralized process control

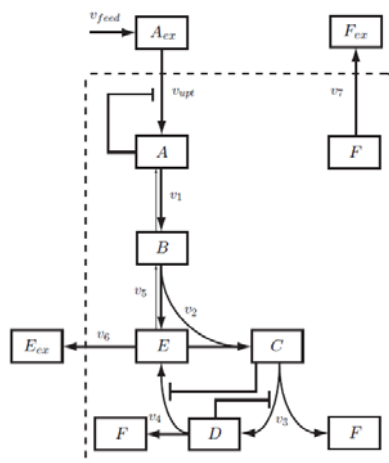


## Power grid



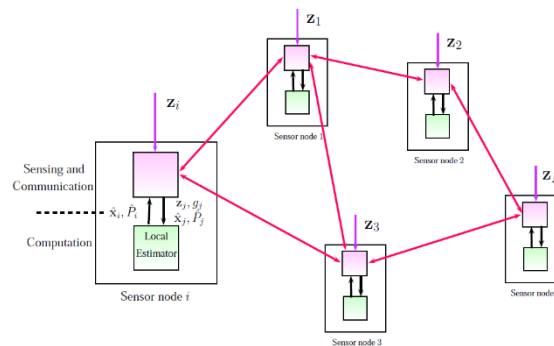
Pierre et al. (2012)

## Metabolic network



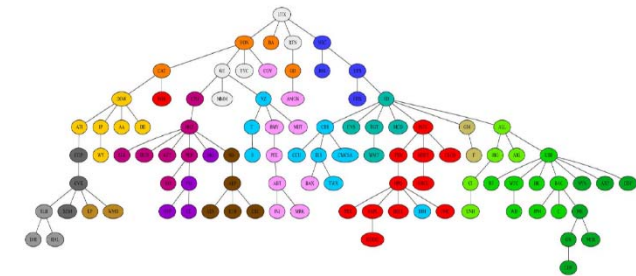
Hillen (2012)

## Distributed control (robotic networks)



Simonetto (2012)

## Stock market

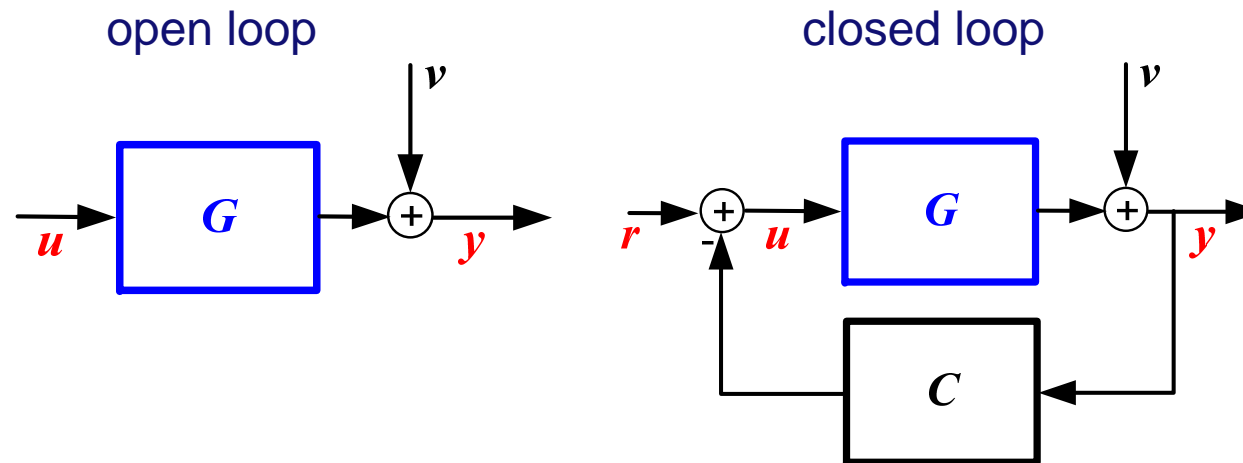


Materassi et al. (2010)

# Introduction – identification

2

The classical (multivariable) identification problems: [Ljung (1999)]

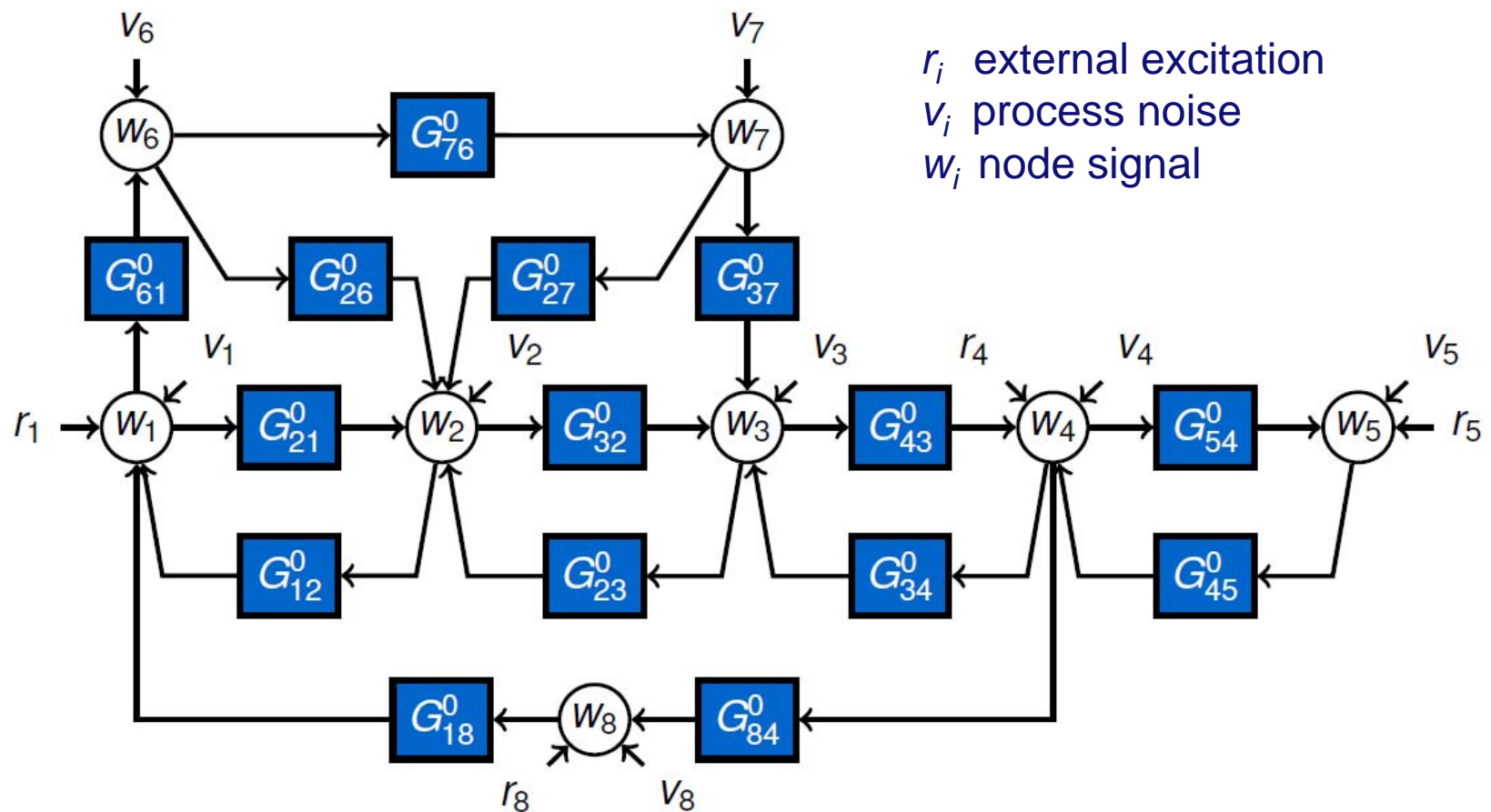


Identify a plant model  $\hat{G}$  on the basis of measured signals  $u$ ,  $y$  (and possibly  $r$ )

We have to move from a fixed and known configuration to deal with and exploit **structure** in the problem.

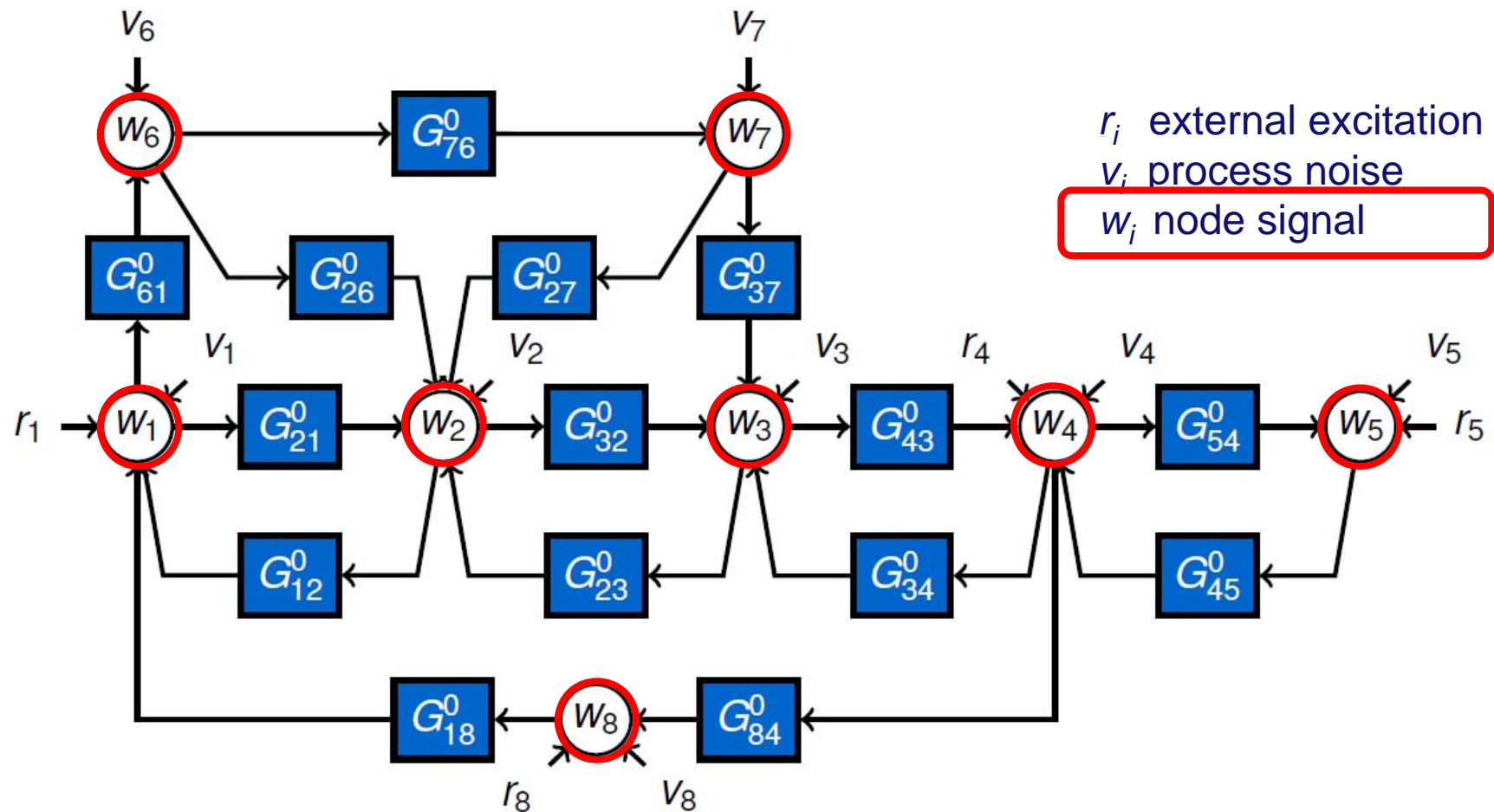
# Dynamic network: what is it?

3



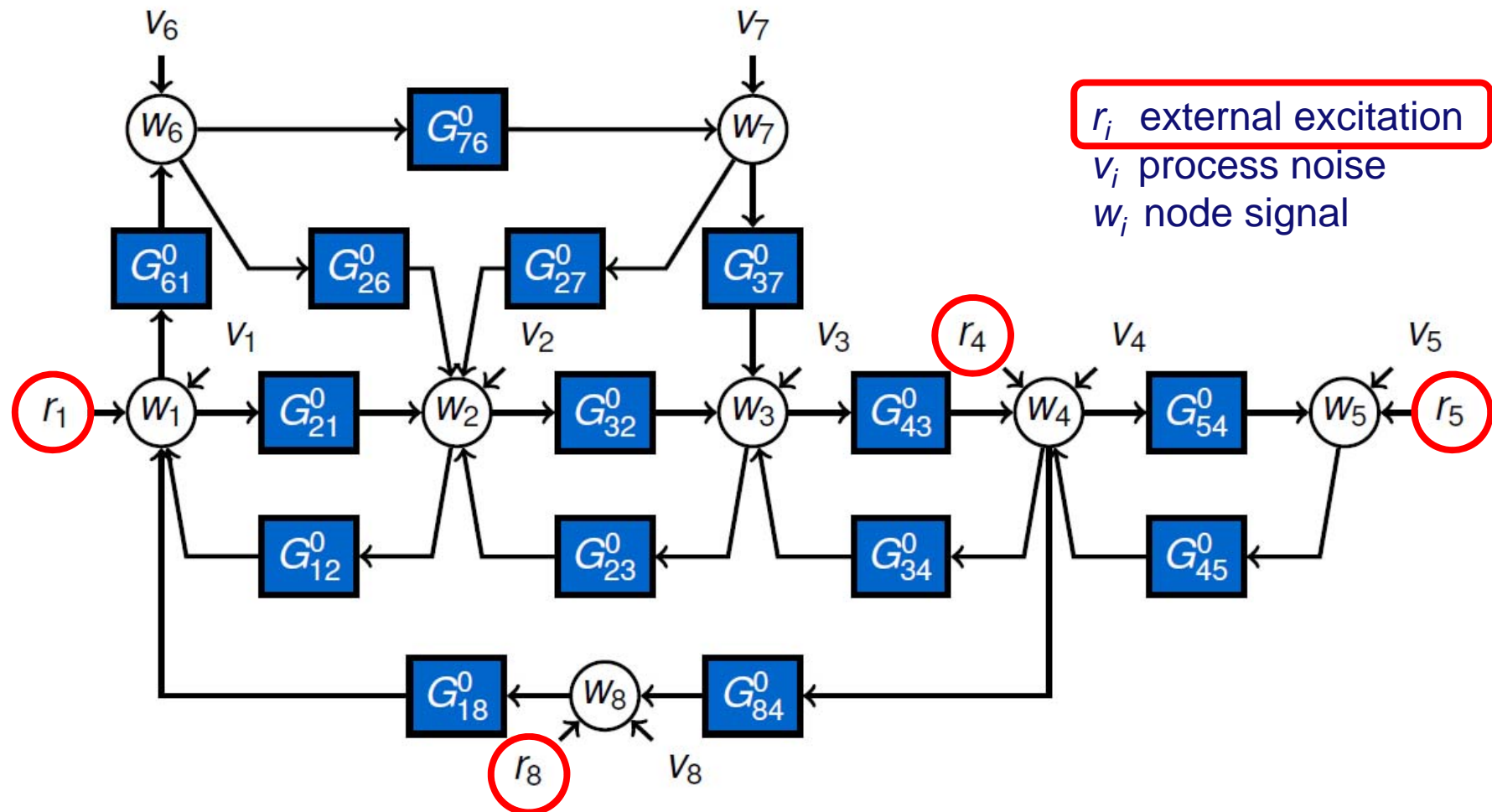
# Dynamic network: what is it?

4



# Dynamic network: what is it?

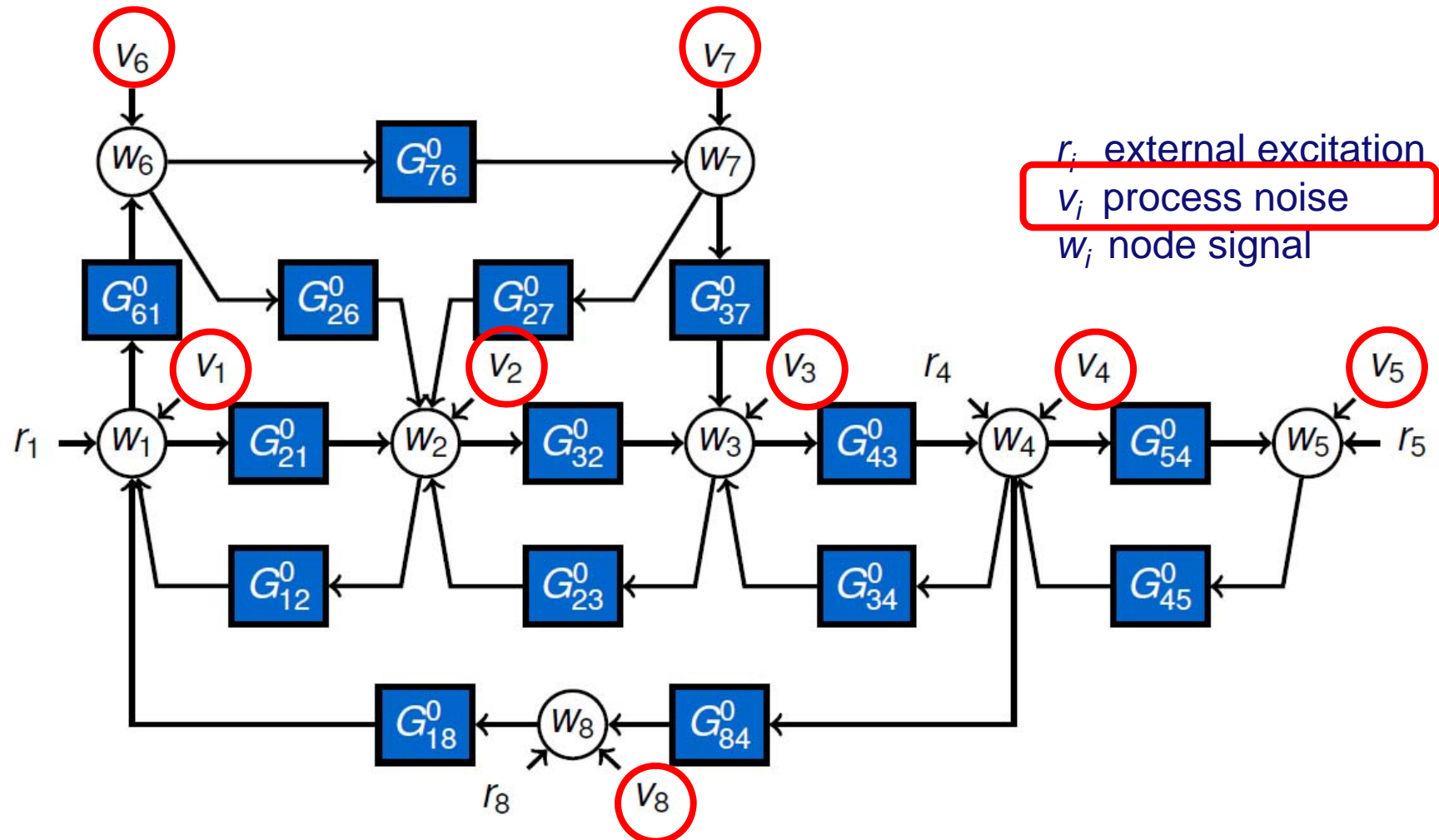
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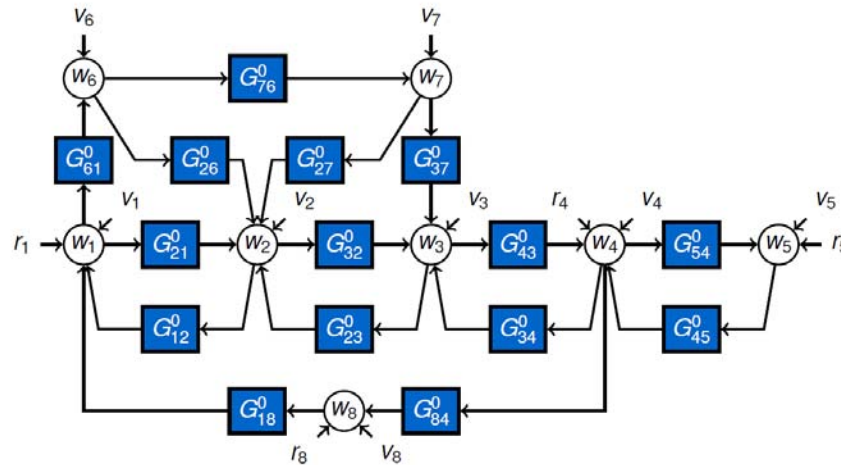
# Dynamic network: what is it?

6



# Introduction

7



$r_i$  external excitation

$v_i$  process noise

$w_i$  node signal

$$\mathbf{v}(t) = \begin{bmatrix} v_1(t) \\ \vdots \\ v_L(t) \end{bmatrix}$$

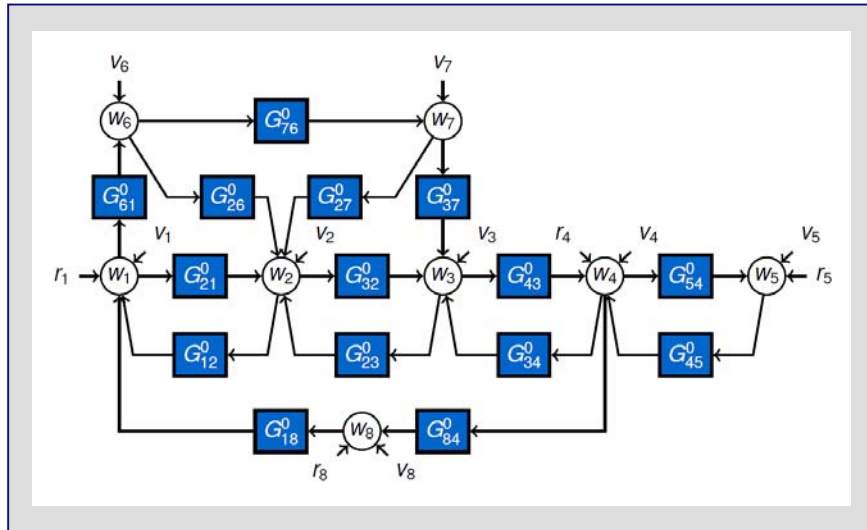
What are assumptions on process noises when identifying (parts of) a network?

- Independent white noise processes
- Vector stochastic process with full rank spectrum,  $\text{rank } \Phi_v(\omega) = L$  a.e. leading to a square noise model:  $\mathbf{v}(t) = \mathbf{H}(q)\mathbf{e}(t)$
- If  $\dim(\mathbf{e}) < L$  then we have “singular” or “reduced-rank” noise



# Network Setup

8



## Assumptions:

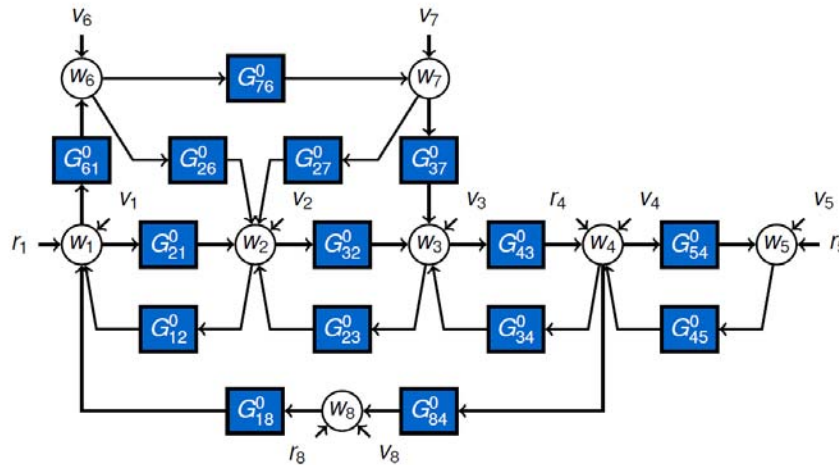
- Total of  $L$  nodes
- Network is well-posed and stable
- Modules may be unstable
- Node signals and excitation signals can be measured

$$\begin{bmatrix} w_1 \\ w_2 \\ \vdots \\ w_L \end{bmatrix} = \underbrace{\begin{bmatrix} 0 & G_{12}^0 & \cdots & G_{1L}^0 \\ G_{21}^0 & 0 & \cdots & G_{2L}^0 \\ \vdots & \cdots & \ddots & \vdots \\ G_{L1}^0 & G_{L2}^0 & \cdots & 0 \end{bmatrix}}_{G^0} \begin{bmatrix} w_1 \\ w_2 \\ \vdots \\ w_L \end{bmatrix} + R^0(q) \begin{bmatrix} r_1 \\ r_2 \\ \vdots \\ r_K \end{bmatrix} + \underbrace{H^0(q)}_v \begin{bmatrix} e_1 \\ e_2 \\ \vdots \\ e_p \end{bmatrix}$$

$$w = G^0 w + R^0 r + H^0 e$$

# Introduction

9



$r_i$  external excitation

$v_i$  process noise

$w_i$  node signal

$$\begin{bmatrix} v_1(t) \\ \vdots \\ v_L(t) \end{bmatrix} = H^0(q) \begin{bmatrix} e_1(t) \\ \vdots \\ e_p(t) \end{bmatrix}$$

**Main question:**

How to identify (parts of) a dynamic network,  
when the process noise is of reduced rank ( $p < L$ )?

- Modelling a reduced-rank stochastic process
- Multi-output identification in a dynamic network  
the joint-direct method with weighted LS
- Constrained LS and maximum likelihood estimation
- Variance-free estimation, minimum variance and the CRLB
- Simulation example

# Modelling reduced rank noise

11

## Assumption

The node signals  $w_j$  are ordered in such a way that the first  $p$  noise components  $v_j$ ,  $j = 1, \dots, p$  constitute a full rank process.

# Modelling reduced rank noise

12

A reduced-rank stochastic process  $v$  with dimension  $L$  and rank  $p$  can equivalently be described in two ways:

a)  $v(t) = \check{H}^0(q)\check{e}(t)$

With  $\check{H}^0 \in \mathbb{R}^{L \times L}(z)$ ,  $\check{e}(t) \in \mathbb{R}^L$  a white noise process,  
 $\check{H}^0$  stable, stably invertible, and monic, and  
 $cov(\check{e}) = \check{\Lambda}^0$  having rank  $p$

b)  $v(t) = H^0(q)e(t)$

With  $H^0 \in \mathbb{R}^{L \times p}(z)$ ,  $e(t) \in \mathbb{R}^p$  a white noise process,  
 $H^0 = \begin{bmatrix} H_a^0 \\ H_b^0 \end{bmatrix}$  with  $H_a^0$  square, stable, stably invertible, and monic,  
 $cov(e) = \Lambda^0$  having full rank  $p$

# Modelling reduced rank noise

13

Relations between descriptions:

$$v(t) = \check{H}^0(q)\check{e}(t) = \begin{bmatrix} H_a^0(q) & 0 \\ H_b^0(q) - \Gamma^0 & I \end{bmatrix} \begin{bmatrix} e \\ \Gamma^0 e \end{bmatrix}$$

with  $\Gamma^0 = \lim_{z \rightarrow \infty} H_b^0(z)$

while  $\check{\Lambda}^0 = \begin{bmatrix} I \\ \Gamma^0 \end{bmatrix} \Lambda^0 \begin{bmatrix} I \\ \Gamma^0 \end{bmatrix}^T$  and  $[\Gamma^0 \quad -I] \check{e}(t) = 0$

Both noise models  $\check{H}^0$  and  $H^0 = \begin{bmatrix} H_a^0 \\ H_b^0 \end{bmatrix}$  will be used.



# Joint-direct identification method

14

We follow a prediction error approach, by predicting **all** node variables:

$$\hat{w}(t|t-1) := \mathbb{E} \left\{ w(t) \mid w^{t-1}, r^t \right\}$$

Then: 
$$\hat{w}(t|t-1) = W_w^0(q)w(t) + W_r^0(q)r(t)$$

with: 
$$\begin{aligned} W_w^0(q) &= I - (\check{H}^0(q))^{-1}(I - G^0(q)), \\ W_r^0(q) &= (\check{H}^0(q))^{-1}R^0(q). \end{aligned}$$

being the unique predictor filters.

# Joint-direct identification method

15

The **network** is defined by:  $(G^0, R^0, H^0, \Lambda^0)$

a network model is denoted by:  $M = (G, R, H, \Lambda)$

and a **network model set** by:

$$\mathcal{M} = \{M(\theta) = (G(\theta), R(\theta), H(\theta), \Lambda(\theta)), \theta \in \Theta\}$$

Then the parametrized predictor:

$$\hat{w}(t|t-1) = W_w(q, \theta)w(t) + W_r(q, \theta)r(t)$$

leads to the prediction error:  $\varepsilon(t, \theta) = w(t) - \hat{w}(t|t-1; \theta)$

**Weighted LS criterion:**

$$\hat{\theta}_N^{WLS} = \arg \min_{\theta \in \Theta} \frac{1}{N} \sum_{t=1}^N \varepsilon^T(t, \theta) Q \varepsilon(t, \theta) \quad Q > 0$$

# Joint-direct identification method

16

**Weighted LS criterion:**

$$\hat{\theta}_N^{WLS} = \arg \min_{\theta \in \Theta} \frac{1}{N} \sum_{t=1}^N \varepsilon^T(t, \theta) Q \varepsilon(t, \theta) \quad Q > 0$$

**Properties:**

- Consistent estimate under regularity conditions,
- Provided model set large enough, appropriate excitation, global network identifiability,
- But for minimum variance an optimal  $Q$  has to be chosen

Typical choice, leading to minimum variance estimator for  $Q \in \mathbb{R}^{L \times L}$

$$Q = [\text{cov}(\check{e})]^{-1} = (\check{\Lambda}^0)^{-1}$$

but in our situation  $\check{\Lambda}^0$  is singular

# Constrained LS and Maximum Likelihood

17

The WLS estimator does not take account of the dependencies in the innovation:

$$\begin{bmatrix} \Gamma^0 & -I \end{bmatrix} \check{e}(t) = 0$$

or differently formulated:

$$\begin{bmatrix} \Gamma^0 & -I \end{bmatrix} \begin{bmatrix} \varepsilon_a(t, \theta_0) \\ \varepsilon_b(t, \theta_0) \end{bmatrix} = 0$$

This can be imposed, by restricting the parametrized model to satisfy:

$$\underbrace{\Gamma(\theta)\varepsilon_a(t, \theta) - \varepsilon_b(t, \theta)}_{:= Z(t, \theta)} = 0$$

We denote:

# Constrained LS and Maximum Likelihood

18

**Constrained LS criterion:**

$$\hat{\theta}_N^{CLS} = \arg \min_{\theta \in \Theta} \frac{1}{N} \sum_{t=1}^N \varepsilon_a^T(t, \theta) Q_a \varepsilon_a(t, \theta) \quad Q_a > 0$$

$$\text{subject to } \frac{1}{N} \sum_{t=1}^N Z^T(t, \theta) Z(t, \theta) = 0$$

**Properties:**

- Consistent estimate under similar conditions as WLS
- The choice  $Q_a = (\Lambda^0)^{-1}$

leads to minimum variance, and ML properties in case of Gaussian noise.

- For indendently parametrized  $\Lambda(\theta)$ , the cost function turns into a determinant function

# Constrained LS and Maximum Likelihood

19

## Implementation:

In practice, constraints could be unfeasible, e.g. in case  $\mathcal{S} \notin \mathcal{M}$

## Constraint relaxation:

$$\hat{\theta}_N^{rel} = \arg \min_{\theta} \frac{1}{N} \sum_{t=1}^N \left( \epsilon_a^T(t, \theta) Q_a \epsilon_a(t, \theta) + \lambda Z^T(t, \theta) Z(t, \theta) \right), \quad \lambda \in \mathbb{R}$$

with tuning parameter  $\lambda \in \mathbb{R}$

For  $\lambda > 0$  the consistency result remains true.

For  $\lambda \rightarrow \infty$  constraint satisfaction

The criterion is equivalent to WLS with

$$Q(\theta) = \begin{bmatrix} Q_a + \lambda \Gamma^T(\theta) \Gamma(\theta) & -\lambda \Gamma^T(\theta) \\ -\lambda \Gamma(\theta) & \lambda I \end{bmatrix}$$



# Minimum variance and CRLB

20

**Asymptotic criterion:**

$$\theta^* = \arg \min_{\theta \in \Theta} \bar{\mathbb{E}} \varepsilon^T(t, \theta) Q_a \varepsilon(t, \theta) \quad \text{subject to} \\ \bar{\mathbb{E}} \mathbf{Z}(t, \theta) \mathbf{Z}^T(t, \theta) = \mathbf{0}$$

When linearizing  $\mathbf{Z}(t, \theta)$  in the neighbourhood of the optimum:

$$\mathbf{Z}(t, \theta) \approx \mathbf{Z}(t, \theta^*) + \mathbf{A}(t)(\theta - \theta^*)$$

the constrained parameter space can be characterized by

$$\theta = \mathbf{S}\rho + \mathbf{C} \quad \rho \in \mathbb{R}^{n_\rho} \text{ of reduced dimension}$$

with  $\mathbf{S}, \mathbf{C}$  determined by:

$$\begin{cases} \mathbf{\Pi S} = \mathbf{0} \\ \mathbf{C} = -\mathbf{\Pi}^\dagger \mathbf{\Pi} \theta^* \end{cases} \quad \text{and } \mathbf{S} \text{ full rank, where } \bar{\mathbb{E}} \mathbf{A}^T(t) \mathbf{A}(t) = \mathbf{\Pi}^T \mathbf{\Pi} \\ \mathbf{\Pi}^\dagger \text{ right inverse}$$

