

Identification of MIMO linear models: introduction to subspace methods

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State space identification from impulse response data

Ho-Kalman realisation theory



Consider the finite dimensional, linear time-invariant (LTI) state space model:

$$\begin{array}{rcl} x(t+1) &=& Ax(t) + Bu(t) \\ y(t) &=& Cx(t) + Du(t) \end{array}$$

Realisation: the problem of computing [A,B,C,D] or an equivalent realisation for the system, from the impulse response (Markov parameters) of the system:

$$\begin{cases} h(0) = D\\ h(t) = CA^{t-1}B, \quad t > 0 \end{cases}$$

Ho-Kalman realisation theory (cont.d)



A few definitions:

• Extended observability matrix:

$$\Gamma_i = \begin{bmatrix} C^T & (CA)^T & (CA^2)^T & \dots & (CA^{i-1})^T \end{bmatrix}^T$$

• Extended controllability matrix:

$$\Delta_i = \begin{bmatrix} B & AB & A^2B & \dots & A^{i-1}B \end{bmatrix}$$



Hankel matrix

$$\mathbf{u}_{-}(t) = \begin{bmatrix} \mathbf{u}^{T}(t-1) & \mathbf{u}^{T}(t-2) & \dots & \mathbf{u}^{T}(t-j) \end{bmatrix}^{T}$$
$$\mathbf{y}_{+}(t) = \begin{bmatrix} \mathbf{y}^{T}(t) & \mathbf{y}^{T}(t+1) & \dots & \mathbf{y}^{T}(t+i-1) \end{bmatrix}^{T}$$

$$H_{i,j} = \begin{bmatrix} h(1) & h(2) & h(3) & \dots & h(j) \\ h(2) & h(3) & h(4) & \dots & h(j+1) \\ h(3) & h(4) & h(5) & \dots & h(j+2) \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ h(i) & h(i+1) & h(i+2) & \dots & h(j+i-1) \end{bmatrix}$$

$$y_+(t) = H_{ij}u_-(t)$$

Ho-Kalman realisation theory (cont.d)



Properties of the Hankel matrix:

- H_{i,j}, i,j , n, has rank n iff h(t) admits an n_{th} order [A,B,C,D] realisation;
- H_{i,j} can be equivalently written as

$$H_{ij} = \Gamma_i \Delta_j$$

Ho-Kalman realisation theory (cont.d)



The realisation can be constructed as follows:

- Let D=h(0);
- Construct the Hankel matrix H_{i,j} from h(1), h(2), ...;
- Factor the Hankel matrix to get Γ_i and Δ_i ;
- Let C=first l rows of Γ_i ;
- Let B=first m columns of Δ_i ;
- Compute A exploiting *shift invariance*, i.e., solving

$$\Gamma_{\uparrow}A = \begin{bmatrix} C \\ CA \\ CA^{2} \\ \vdots \\ CA^{i-2} \end{bmatrix} A = \begin{bmatrix} CA \\ CA^{2} \\ CA \\ \vdots \\ CA^{i-1} \end{bmatrix} = \Gamma_{\downarrow}$$



What if *noisy* measurements of h(t) are available? $\tilde{h}(t) = h(t) + w(t)$

Idea:

- Construct the noisy Hankel matrix \hat H_{i,i}
- Factor the matrix using the SVD:

$$\tilde{H}_{ij} = \begin{bmatrix} \mathbf{U}_s & \mathbf{U}_0 \end{bmatrix} \begin{bmatrix} \boldsymbol{\Sigma}_s & \mathbf{0} \\ O & \boldsymbol{\Sigma}_0 \end{bmatrix} \begin{bmatrix} \mathbf{V}_s^T \\ \mathbf{V}_0^T \end{bmatrix}$$

Estimate H_{i,j} as the best rank n approximation:

$$\widehat{H}_{ij} = \left(\mathbf{U}_s \boldsymbol{\Sigma}_s^{1/2}\right) \left(\boldsymbol{\Sigma}_s^{1/2} \mathbf{V}_s^T\right) \quad \rightarrow \quad \widehat{\boldsymbol{\Gamma}}_i = \mathbf{U}_s \boldsymbol{\Sigma}_s^{1/2}, \quad \widehat{\boldsymbol{\Delta}}_j = \boldsymbol{\Sigma}_s^{1/2} \mathbf{V}_s^T$$



14 🖵 10⁻³ 12 10 8 Impulse response e 2 n -2 ⊾ 0 100 200 500 700 800 300 400 600 samples

Model for a Peltier cell (n=4, i=20)

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Subspace Model Identification: deterministic case

The data equation



Note that we can write the following equation (i > n)

$$\begin{bmatrix} y(t) \\ y(t+1) \\ y(t+2) \\ \vdots \\ y(t+i-1) \end{bmatrix} = \begin{bmatrix} C \\ CA \\ CA^2 \\ \dots \\ CA^{i-1} \end{bmatrix} x(t) + \begin{bmatrix} D & 0 & 0 & \dots & 0 \\ CB & D & 0 & \dots & 0 \\ CAB & CB & D & \dots & 0 \\ \vdots \\ CA^{i-1}B & \dots & CB & D \end{bmatrix} \begin{bmatrix} u(t) \\ u(t+1) \\ u(t+2) \\ \vdots \\ u(t+i-1) \end{bmatrix}$$

which describes the system over a window of finite length.

The data equation (cont.d)



Repeating for various initial times we get the *data* equation

$$Y_{t,i,j} = \Gamma_i X_{t,j} + H_i U_{t,i,j}$$

where $Y_{t,i,j}$, $U_{t,i,j}$ are Hankel matrices:

$$Y_{t,i,j} = \begin{bmatrix} y(t) & \cdots & y(t+j-1) \\ y(t+1) & \cdots & y(t+j) \\ \vdots & \ddots & \vdots \\ y(t+i-1) & \cdots & y(t+i+j-2) \end{bmatrix}$$

and $X_{t,j}$ is defined as $X_{t,j} = \begin{bmatrix} x(t) & x(t+1) & \cdots & x(t+j-1) \end{bmatrix}$



The MOESP algorithm (Verhaegen and Dewilde 1991):

- 1. Construct projection $\Pi^{?}$ such that $U_{t,i,j} \Pi^{?} = 0$
- 2. Project data equation using $\Pi^{?}~$ to recover column space of Γ_{i}

$$Y_{t,i,j} \Pi^{\perp} = \Gamma X_{t,j} \Pi^{\perp}$$

- 3. Construct a basis for the column space of $\Gamma_{\rm I}$ and estimate A and C.
- 4. Solve LS problem for estimation of B and D.

Computing the projection $\Pi^{?}$



We look for $\Pi^{?}$ such that $U_{t,i,j} \Pi^{?} = 0$.

The solution is given by

$$\Pi^{\perp} = I - U_{t,i,j}^{T} \left(U_{t,i,j} U_{t,i,j}^{T} \right)^{-1} U_{t,i,j}$$

since in fact

$$U_{t,i,j} \Pi^{\perp} = U_{t,i,j} - U_{t,i,j} U_{t,i,j}^{T} \left(U_{t,i,j} U_{t,i,j}^{T} \right)^{-1} U_{t,i,j} = 0$$

Note that constructing Π ? requires $(U_{t,i,j}U_{t,i,j}^T)$ to be nonsingular.

Implementation of the projection



The projection $\Pi^{?}$ can be computed and implemented via the RQ factorisation:

$$\begin{bmatrix} U_{t,i,j} \\ Y_{t,i,j} \end{bmatrix} = \begin{bmatrix} R_{11} & 0 \\ R_{21} & R_{22} \end{bmatrix} \begin{bmatrix} Q_1 \\ Q_2 \end{bmatrix} = RQ, \quad QQ^T = \begin{bmatrix} Q_1Q_1^T & Q_1Q_2^T \\ Q_2Q_1^T & Q_2Q_2^T \end{bmatrix} = I$$

which can be written as

$$U_{t,i,j} = R_{11}Q_1$$

$$Y_{t,i,j} = \Gamma_i X_{t,j} + H_i U_{t,i,j} = R_{21}Q_1 + R_{22}Q_2$$

and therefore

$$R_{22} = \Gamma_i X_{t,j} Q_2^T$$





Therefore, considering the equation

$$Y_{t,i,j} = \Gamma_i X_{t,j} + H_i U_{t,i,j} = R_{21}Q_1 + R_{22}Q_2$$

and right-multiplying by Q_2^{T} one gets

$$R_{22} = \Gamma_i X_{t,j} Q_2^T$$

so R_{22} , of dimension (il £ il) and computed from data only, contains information on Γ_i .

Under what conditions range(R_{22})=range(Γ_i)?

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A rank condition



Theorem 1: if u(t) is such that

$$\operatorname{rank}\left(\begin{bmatrix} X_{t,j} \\ U_{t,i,j} \end{bmatrix} \right) = n + im$$

then

range(
$$R_{22}$$
) = range(Γ_i).

Problem: this is not yet an identifiability condition,since it depends on the state.However, it implies the following.

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Theorem 2 (Jansson 1997): if the input u is persistently exciting of order n+i, then

$$\lim_{N \to \infty} \frac{1}{N} \begin{bmatrix} X_{t,j} \\ U_{t,i,j} \end{bmatrix} \begin{bmatrix} X_{t,j}^T & U_{t,i,j}^T \end{bmatrix} > 0$$

(i.e., the rank condition of Theorem 1 holds).



Rank reduction of estimated column space of Γ_i performed via singular value decomposition of R₂₂. Under p.e. assumptions, rank(R₂₂=n), so

$$R_{22} = \begin{bmatrix} U_n & U_n^{\perp} \end{bmatrix} \mathbf{\Sigma} V^T =$$
$$= \begin{bmatrix} U_n & U_n^{\perp} \end{bmatrix} \begin{bmatrix} \mathbf{\Sigma}_n & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix} V^T \quad \Rightarrow \quad \widehat{\mathbf{\Gamma}}_i = U_n$$

The inspection of the singular values provides information about model order.

Estimation of A and C



Let C=first l rows of computed Γ_i ;

Compute A exploiting *shift invariance*, i.e., solving the system of linear equations

$$\Gamma_{\uparrow}A = \begin{bmatrix} C \\ CA \\ CA^2 \\ \vdots \\ CA^{i-2} \end{bmatrix} A = \begin{bmatrix} CA \\ CA^2 \\ CA \\ \vdots \\ CA^{i-1} \end{bmatrix} = \Gamma_{\downarrow}$$

A simple example (Van Der Veen et al. 1993)



Consider the LTI system ($|\alpha| < 1$)

$$\begin{array}{rcl} x(t+1) &=& \alpha x(t) + \alpha u(t) \\ y(t) &=& x(t) + u(t) \end{array}$$

and apply the input sequence (x(1)=0)

$$u = [1 \ 2 \ 1 \ 1]^T$$

that gives the corresponding output sequence

$$y = [1 \ 2 + \alpha \ 1 + 2\alpha + \alpha^2 \ 1 + \alpha + 2\alpha^2 + \alpha^3]^T$$

A simple example (cont.d)



Choosing i=2 and j=3 we can construct the compound matrix

$$\begin{bmatrix} U_{t,i,j} \\ Y_{t,i,j} \end{bmatrix} = \begin{bmatrix} 1 & 2 & 1 \\ 2 & 1 & 1 \\ 1 & 2+\alpha & 1+2\alpha+\alpha^2 \\ 2+\alpha & 1+2\alpha+\alpha^2 & 1+\alpha+2\alpha^2+\alpha^3 \end{bmatrix} = RQ$$

and computing the RQ factorisation we get

$$R = \begin{bmatrix} 1 & 2 & 0 \\ 2 & 1 & 0 \\ 1 & 2+\alpha & 5\alpha+3\alpha^2 \\ 2+\alpha & 1+2\alpha+\alpha^2 & 5\alpha^2+3\alpha^3 \end{bmatrix}$$

A simple example (cont.d)



We can now factor R_{22} as :

$$R_{22} = \begin{bmatrix} 5\alpha + 3\alpha^2 \\ 5\alpha^2 + 3\alpha^3 \end{bmatrix} = \begin{bmatrix} 1 \\ \alpha \end{bmatrix} \begin{bmatrix} 5\alpha + 3\alpha^2 \end{bmatrix}$$

So

$$\widehat{\Gamma}_2 = \begin{bmatrix} 1 \\ \alpha \end{bmatrix}$$

and finally

$$\hat{C} = 1, \quad \hat{A} = \alpha$$

A numerical example



Consider the order 2 system

$$\begin{array}{rcl} x_1(t+1) &=& 0.3x_1(t) + x_2(t) + u(t) \\ x_2(t+1) &=& 0.7x_2(t) + u(t) \\ y(t) &=& x_1(t) \end{array}$$

and measure the response to a 500 samples realisation of white gaussian noise.

I/O data







 U_{tij} and Y_{tij} are constructed with i=10 and j=490, so R_{22} is 10 \pm 10. Its singular values are given by



Estimated A and C



Numerical results of the estimation procedure:

$$A = \begin{bmatrix} 0.8033 & 0.5950 \\ -0.0874 & 0.1967 \end{bmatrix}, \quad C = \begin{bmatrix} -0.5897 & 0.7799 \end{bmatrix}$$

Note that

- The computed A and C are in a different state space basis from the original system;
- They are equivalent to the original A and C;
- Question: what determines the basis of the estimated matrices?

MATLAB code for the estimation of A and C



```
function [A,C]=omoesp(u,y,i,j,n);
```

```
sy=size(y);su=size(u);
datalen=min([max(sy) max(su)]);
m=min(su); l=min(sy);
```

```
H=[];
for ii=1:i
H=[H u(ii:ii+j-1,:)];
end
```

```
for ii=1:i
H=[H y(ii:ii+j-1,:)];
end
```

R=triu(qr(H))';

R22=R(m*i+1:(m+l)*i,m*i+1:(m+l)*i);

[U,S,Vt]=svd(R22);

Un=U(:,1:n);

C=Un(1:l,:);

A=Un(1:l*(i-1),:)\Un(l+1:l*i,:);



The output of the identified model is given by:

$$\widehat{y}(t) = Du(t) + \sum_{r=0}^{t-1} \mathbf{C} \mathbf{A}^{t-r-1} Bu(r)$$

we aim at writing the above as a linear regression in the elements of B and D:

$$\hat{y}(t) = \phi_D^T(t)vec(\mathbf{D}) + \phi_B^T(t)vec(\mathbf{B})$$

where for X 2 $R^{(m \ \ n)}$

 $vec(X) = \begin{bmatrix} x_{11} & \dots & x_{m1} & x_{12} & \dots & x_{m2} & x_{1n} & \dots & x_{mn} \end{bmatrix}^T$

For this, we need to introduce Kronecker products.



Let A 2 $R^{(m \pounds n)}$ and B 2 $R^{(r \pounds s)}$, then the (mr £ ns) matrix

$$A \otimes B = \begin{bmatrix} a_{11}B & a_{12}B & \dots & a_{1n}B \\ a_{21}B & a_{22}B & \dots & a_{2n}B \\ \vdots & & & \vdots \\ a_{m1}B & a_{m2}B & \dots & a_{mn}B \end{bmatrix}$$

is called the Kronecker product of A and B.

vec operation and Kronecker product



There is a connection between Kronecker products and the vec operation.

Let A 2 $R^{(m \text{ f } n)}$, B 2 $R^{(n \text{ f } o)}$, C 2 $R^{(o \text{ f } p)}$, then

$$vec(ABC) = (C^T \otimes A) vec(B)$$



Using Kronecker products the output of the identified model t-1

$$\hat{y}(t) = Du(t) + \sum_{r=0}^{\infty} \mathbf{C} \mathbf{A}^{t-r-1} Bu(r)$$

can be written as

$$\hat{y}(t) = [u(t)^T \otimes \mathbf{I}_l] vec(\mathbf{D}) + (\sum_{r=0}^{t-1} u(r)^T \otimes \mathbf{C} \mathbf{A}^{t-r-1}) vec(\mathbf{B})$$

so that B and D can be obtained from:

$$B, D = \arg\min_{B,D} \sum_{k=0}^{s} \left[y(t) - [u(t)^T \otimes \mathbf{I}_l] vec(\mathbf{D}) - (\sum_{r=0}^{t-1} u(r)^T \otimes \mathbf{CA}^{t-r-1}) vec(\mathbf{B}) \right]^2$$

which is clearly a least squares problem in B and D.



Subspace Model Identification: output error case

SMI: output error case



Consider the finite dimensional, linear time-invariant (LTI) state space model:

$$x(t+1) = Ax(t) + Bu(t)$$

with the measurement equation

$$y(t) = Cx(t) + Du(t) + v(t)$$

where v is a zero-mean, white measurement noise, uncorrelated with u.

We want to analyse the effect of v on the identification algorithm we studied in the deterministic case.

The data equation with measurement error

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When adding measurement noise the data equation becomes

$$Y_{t,i,j} = \Gamma_i X_{t,j} + H_i U_{t,i,j} + V_{t,i,j}$$

where $\boldsymbol{V}_{t,i,j}$ is defined as

$$V_{t,i,j} = \begin{bmatrix} v(t) & \cdots & v(t+j-1) \\ v(t+1) & \cdots & v(t+j) \\ \vdots & \ddots & \vdots \\ v(t+i-1) & \cdots & v(t+i+j-2) \end{bmatrix}$$



As in the deterministic case, we:

- Construct projection $\Pi^{?}$ such that $U_{t,i,j} \Pi^{?} = 0$
- Project data equation using $\Pi^{?}~$ to recover column space of Γ_{i}

$$Y_{t,i,j} \Pi^{\perp} = \Gamma X_{t,j} \Pi^{\perp} + V_{t,i,j} \Pi^{\perp}$$

• Using the RQ factorisation we obtain now

$$R_{22} = \Gamma_i X_{t,j} Q_2^T + V_{t,i,j} Q_2^T$$

Asymptotic properties of R_{22}



Can we use R_{22} to estimate the observability subspace?

Theorem 3: if v ' WN(0, σ^2 I) and u is p.e. of order n+i, then

$$\lim_{N \to \infty} \frac{1}{N} R_{22} R_{22}^T = \Gamma_i M \Gamma_i^T + \sigma^2 I_{il}, \quad M = \lim_{N \to \infty} \frac{1}{N} X_{t,j} \Pi^\perp X_{t,j}^T$$

and

$$\lim_{N \to \infty} \frac{1}{N} R_{22} R_{22}^T = \begin{bmatrix} U_n & U_n^{\perp} \end{bmatrix} \begin{bmatrix} \Sigma_n + \sigma^2 I_n & 0 \\ 0 & \sigma^2 I_{il-n} \end{bmatrix} U^T$$



Consider again the order n=2 system

$$\begin{aligned} x_1(t+1) &= 0.3x_1(t) + x_2(t) + u(t) \\ x_2(t+1) &= 0.7x_2(t) + u(t) \\ y(t) &= x_1(t) + v(t) \end{aligned}$$

and measure the response to a 500 samples realisation of white gaussian noise, subject to v'(0,0.01).

We repeat the identification 1000 times, with different realisations of the noise v to assess the *average* effect of measurement noise.

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Construction of R_{22} and SVD



 R_{22} is constructed with i=10 and j=490. Its singular values are given by



Estimated eigenvalues of A







Subspace Model Identification: the general case

Identification and control of rotary wing aircraft

SMI: the general case



Consider the finite dimensional, linear time-invariant (LTI) state space model:

$$x(t+1) = Ax(t) + Bu(t) + w(t)$$

with the measurement equation

$$y(t) = Cx(t) + Du(t) + v(t)$$

with w and v zero-mean white noises, uncorrelated with u.

Does the orthogonal projection algorithm still work?

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Example



Consider the n=1 system A=0.7; B=1; C=1; D=0; and compare the performance of the SMI algorithm with and without process noise w:



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When process noise is present, the data equation becomes

$$Y_{t,i,j} = \Gamma_i X_{t,j} + H_i U_{t,i,j} + E_i W_{t,i,j} + V_{t,i,j}$$

and therefore the residual is not white anymore and the results we have seen so far do not hold.

The problem can be solved by introducing Instrumental Variables.

Instrumental variable (IV) algorithms



Assume that a matrix Z (Instrumental Variable) can be found such that

$$rank\left(\lim_{N\to\infty}\frac{1}{N}(X_{t,j}\Pi^{\perp})Z^{T}\right) = n \quad \lim_{N\to\infty}\frac{1}{N}(E_{i}W_{t,i,j}+V_{t,i,j})Z^{T} = 0$$

Then the column space of Γ_i can be estimated from

$$Y_{t,i,j} \Pi^{\perp} Z^T = \Gamma X_{t,j} \Pi^{\perp} Z^T + (E_i W_{t,i,j} + V_{t,i,j}) \Pi^{\perp} Z^T$$



The term $Y_{t,i,j}\Pi^{?}Z^{\mathsf{T}}$ can be computed from the RQ factorisation

$$\begin{bmatrix} U_{t,i,j} \\ Z \\ Y_{t,i,j} \end{bmatrix} = \begin{bmatrix} R_{11} & 0 & 0 \\ R_{21} & R_{22} & 0 \\ R_{31} & R_{32} & R_{33} \end{bmatrix} \begin{bmatrix} Q_1 \\ Q_2 \\ Q_3 \end{bmatrix}$$

And it holds that

$$Y_{t,i,j} \mathbf{\Pi}^{\perp} Z^T = R_{32} R_{22}^T$$

and therefore

range (
$$\Gamma_i$$
) = range $\left(\lim_{N \to \infty} \frac{1}{N} R_{32} R_{22}^T\right)$



Possible choice of IVs (MOESP-PO, Verhaegen 1994):

- Consider the available I/O data set and split it in two parts (past and future), the second shifted ahead of i samples with respect to the first;
- Write two separate data equations for past and future data:

 $Y_1 = \Gamma_i X_1 + H_i U_1 + E_i W_1 + V_1$ $Y_2 = \Gamma_i X_2 + H_i U_2 + E_i W_2 + V_2$

Use past data as IVs in the future data equation;



 Using Past Inputs and Outputs as IVs one can compute the RQ factorisation

$$\begin{bmatrix} U_2 \\ Z \\ Y_2 \end{bmatrix} = \begin{bmatrix} U_1 \\ V_1 \\ Y_1 \end{bmatrix}_{Y_2} = \begin{bmatrix} R_{11} & 0 & 0 & 0 \\ R_{21} & R_{22} & 0 & 0 \\ R_{31} & R_{32} & R_{33} & 0 \\ R_{41} & R_{42} & R_{43} & R_{44} \end{bmatrix} \begin{bmatrix} Q_1 \\ Q_2 \\ Q_3 \\ Q_4 \end{bmatrix}$$
range $\left(\lim_{N \to \infty} \frac{1}{\sqrt{N}} \begin{bmatrix} R_{42} & R_{43} \end{bmatrix}\right) = \operatorname{range}(\Gamma_i)$

• Rank reduction of estimated column space of $\Gamma_{\rm i}$ performed via a singular value decomposition.

Persistency of excitation conditions



- An input u which is p.e. of order n+2i will "almost always" lead to a consistent estimate of A and C.
- The theory for the IV algorithm is not complete yet...

MATLAB code for the estimation of A and C



```
function [A,C]=moesppo(u,y,i,j,n);
```

```
sy=size(y);su=size(u);
datalen=min([max(sy) max(su)]);
m=min(su); l=min(sy);
Up=[];Uf=[];Yp=[];Yf=[];
```

```
for ii=1:i
    Up=[Up u(ii:ii+j1,:)];
    Yp=[Yp y(ii:ii+j1,:)];
end
```

```
for ii=i+1:2*i
Uf=[Uf u(ii:ii+j1,:)];
Yf=[Yf y(ii:ii+j1,:)];
end
```

R=triu(qr([Uf Up Yp Yf]))';

```
R4243=R((2*m+l)*i+1:2*(m+l)*i,m*i+
1:(2*m+l)*i);
```

[U,S,Vt]=svd(R4243);

Un=U(:,1:n);

C=Un(1:l,:);

A=Un(1:l*(i1),:)``Un(l+1:l*i,:);

(Some) extensions of SMI algorithms



- Recursive versions of all the presented algorithms;
- Identification of linear models in *continuous time*:

$$\dot{x}(t) = Ax(t) + Bu(t) + w(t)$$

$$y(t) = Cx(t) + Du(t) + v(t)$$

 Identification of classes of nonlinear models, including, e.g., Wiener models:





- Choice of parameter i:
 - The choice of i affects the variance of the estimates;
 - No general guidelines except for condition i >> n;
- Asymptotic variance of the estimated [A,B,C,D] matrices:
 - Analytical expressions for the variance of the estimates exist;
 - Expressions too complicated to be of practical use!
 - The estimates are asymptotically Gaussian;
 - No results available for efficiency;

SMI vs Prediction Error Methods



Advantages:

- SMI algorithms work equally well for SISO and MIMO problems;
- They are very reliable from the numerical point of view;

Disadvantages:

- SMI algorithms are not "optimal" in any sense;
- Very difficult to use them for structured problems;

Available software tools



- Functions for SMI in the System Identification Toolbox for Matlab;
- Dedicated SMI Toolbox, again based on Matlab;
- Fast code in C and Fortran available in the Slicot library.