Frequency domain subspace-based identification of discrete-time power spectra from uniformly spaced measurements

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A subspace algorithm proposed by Van Overche etal. (1997)) to identify state space models from given uniformly spaced power spectrum measurements.

- When spectrum samples are on a uniform grid of frequencies, the inverse discrete-Fourier transform (IDFT) of the given power spectrum can be expressed in terms of the state-space parameters of the spectral factor.
- A realization theory based on McKelvey *etal*.:1996 is then devised to obtain the system matrices from this IDFT.
- Two methods ensuring positivity of the identified spectrum are proposed.

Consider the $I \times I$ dimensional square discrete time system:

$$x_{k+1} = Ax_k + Bu_k,$$

$$y_k = Cx_k + Du_k$$
(1)

with $A \in \mathbf{R}^{n \times n}$, $B \in \mathbf{R}^{n \times l}$, $C \in \mathbf{R}^{l \times n}$ and $D \in \mathbf{R}^{l \times l}$ nonsingular. The vector sequences u_k , $y_k \in \mathbf{R}^l$ are the input and output sequences, respectively.

- (1)–(2) stable and strictly minimum phase: all eigenvalues of A and $A BD^{-1}C$ strictly inside the unit circle.
- $\{A, B\}$ and $\{A, C\}$ controllable and observable.
- A non-singular.

The system (1)–(2) is thus a minimal stochastic system.

The transfer function of the system:

$$G(z)=D+C(zI_n-A)^{-1}B.$$

The power spectrum associated with (I)–(2):

$$S(z) = G(z)G^{T}(z^{-1}).$$
 (2)

The system (I)–(2) is the innovation form, unity variance, minimum phase spectral factor associated with S(z).

• (2) satisfies the positive realness condition

$$S(z) > 0, \qquad |z| = 1.$$

imposing positivity constraints on the spectrum samples as well as on the identified power spectrum.

Splitting of the power spectrum (Caines:1988)

$$S(z) = H(z) + H^T(z^{-1})$$

where

• The IDFT v_k of a given complex signal $V_k = V(e^{j2\pi k/2N})$, $k = O, \dots, 2N - 1$ is defined by

$$v_k = rac{1}{2N} \sum_{r=0}^{2N-1} V_r e^{j2\pi rk/2N}).$$

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The problem treated can now be stated as follows:

Given

N + 1 matrices $S_k \in \mathbf{C}^{l \times l}$ of the power spectrum S(z) evaluated at N + 1 equidistant points over the unit circle:

$$S_k = S(e^{j2\pi k/2N}), \qquad k = 0, \cdots N,$$

Find

- The system matrices A, G, C, Λ₀ describing the power spectrum.
- The system matrices *A*, *B*, *C*, *D* describing the spectral factor (I)–(2).

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Expand the N + 1 given points S_k to 2N points as follows:

$$S_{N+k} = S^*_{N-k}, \qquad k = 1, \cdots, N-1.$$

From now on, S_k denotes this signal of length 2N.

Theorem 1 (Inverse discrete Fourier transform). With $M = (I_n - A^{2N})^{-1}$, the IDFT $s_k \in \mathbf{R}^{l \times l}$ of the given power spectrum S_k is given by

$$\begin{split} s_0 &= \Lambda_0 + C A^{2N} M G + G^T (A^T)^{2N} M^T C^T, \\ s_k &= C A^{k-1} M G + G^T (A^T)^{2N-k-1} M^T C^T, \ 1 \leq k < 2N. \end{split}$$

• The proof of this theorem is based on McKelvey etal.:1996.

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The extended observability matrix $\Gamma_q \in \mathbf{R}^{lq \times n}$ and reversed extended observability matrix $\tilde{\Gamma}_r \in \mathbf{R}^{lr \times n}$ are defined as

$$\Gamma_q = \begin{pmatrix} C \\ \vdots \\ CA^{q-1} \end{pmatrix}, \qquad \tilde{\Gamma}_r = \Pi \Gamma_r$$

where $\Pi \in \mathbf{R}^{IN \times IN}$ is the permutation matrix given by

$$\Pi = \left(\begin{array}{ccc} 0 & \cdots & I_l \\ \vdots & \ddots & \vdots \\ I_l & \cdots & 0 \end{array}\right)$$

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The extended controllability matrix $\Delta_r \in \mathbf{R}^{n \times lr}$ and reversed extended controllability matrix $\tilde{\Delta}_q \in \mathbf{R}^{n \times lq}$ are defined as

$$\Delta_r = (G \cdots A^{r-1}), \qquad \tilde{\Delta}_q = \Delta_q \Pi.$$

Since (I)–(2) is minimal, the matrices Γ_q, Γ_r and Δ_r, Δ_q are respectively of full column and row rank *n*.

Let

$$S = \left(\begin{array}{ccc} s_1 & \cdots & s_r \\ \vdots & \ddots & \vdots \\ s_q & \cdots & s_{q+r-1} \end{array}\right)$$

with $q, r \ge 2n, r + q < 2N$.

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Theorem 2 The block-Hankel matrix S can be decomposed as

$$\mathcal{S} = \begin{pmatrix} \Gamma_q \ \tilde{\Delta}_q^T \end{pmatrix} \begin{pmatrix} M & 0 \\ 0 & M^T \end{pmatrix} \begin{pmatrix} \Delta_r \\ \tilde{\Gamma}_r^T \end{pmatrix}$$

which leads to the following results:

- $\operatorname{rank}(\mathcal{S}) = 2n$.
- The column space of S can be expressed in terms of the system matrices as column space S = column space (Γ_q Δ̃^T_q).
- The row space of S can be expressed in terms of the system matrices as

row space
$$S = \text{row space} \left(\begin{array}{c} \Delta_r \\ \tilde{\Gamma}_r^T \end{array} \right)$$

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- A is determined by a procedure similar to in Akçay and Türkay:2004 (SVD followed by eigen decomposition).
- G, C, and A are determined in the same step of Van Overschee *etal*.:1997 whereas G, C, and Δ₀ are estimated from S_k by an LS procedure in Akçay and Türkay:2004.

• With
$$M = (I_n - A^{2N})^{-1}$$
,

$$\Delta_o = s_0 - C A^{2N-1} M G - G^T (A^T)^{2N-1} M^T C^T.$$

- *B* and *D* are determined as in Akçay and Türkay:2004.
- Consistency follows from McKelvey *etal.*:1996.

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When S_k are not generated by a finite dimensional linear system or noise corrupted, there is no guarantee that the identified power spectrum, which is determined by A, G, C, and Λ_0 will be positive real.

- When the identified sequence is not positive real, the Riccati equation has no positive definite solution and the spectral factor cannot be computed.
- Two possible solutions to this problem are presented.
 - Both of these solutions start from given matrices *A* and *C*. The solutions then state how *G* and Λ_0 are determined through the solution of an optimization problem which guarantees a positive real identified power spectrum.

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Given *A* and *C*, a positive real identified spectrum can be guaranteed by solving the following optimization problem:

Given the known transfer matrix

$$L(z) = \left(C(zI_n - A)^{-1} I_l \right),$$

solve

$$\min_{Q,S,R} \sum_{k=0}^{2N-1} \|S_k - L(e^{j2\pi k/2N}) \begin{pmatrix} Q & S \\ S^T & R \end{pmatrix} L(e^{-j2\pi k/2N}) \|_F^2$$

constrained to

$$\left(\begin{array}{cc} Q & S \\ S^T & R \end{array}\right) \ge 0.$$

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 The system matrices G and Λ₀ can then be found by solving the set of equations:

$$P = APA^T + Q, \qquad (3)$$

$$G = APC^{T} + S, \qquad (4)$$

$$\Lambda_0 = CPC^T + R.$$
 (5)

- The constraint guarantees that the resulting identified quadruple A, G, C, Λ₀ leads to a positive real spectrum.
- This optimization problem can be converted to an LMI.
- The drawback: the LMI software is not suited for large N.

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Given *A* and *C*, a positive real identified spectrum can be guaranteed by solving the following NLS optimization:

$$\min_{B,D} \sum_{k=0}^{2N-1} \|S_k - L(e^{j2\pi k/2N}) \begin{pmatrix} B \\ D \end{pmatrix} (B^T D^T) L(e^{-j2\pi k/2N}) \|_F^2.$$

 To insure a minimum phase model, (3)–(5) can be solved for *G* and Λ₀, after which a new B and D (guaranteeing a minimum phase model) can be computed through the solution of the Riccati equation.

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Initial guesses of *B* and *D*

"Perturb Λ_0 to $\tilde{\Lambda}_0$ so that the the power spectrum associated with the resulting quadruple $\{A, G, C, \tilde{\Lambda}_0\}$ is positive real"

$$\tilde{\Lambda}_0 = \Lambda_0 + \tau I_I, \qquad \tau > 0.$$

 Taking *τ* large will trivially ensure positive realness of the power spectrum. However, we would like to keep *τ* as small as possible, which can be posed as an LMI:

$$\begin{array}{ll} \mbox{min} & \tau & \mbox{subject to} \ \tau > 0, \ \ P > 0, \\ \left(\begin{array}{cc} P - APA^T & G - APC^T \\ G^T - CPA^T & \Lambda_0 - CPC^T + \tau I_I \end{array} \right) > 0. \end{array}$$

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- A general perturbation model Λ₀ = Λ₀ + τ where the Frobenius norm of τ is to be minimized can be introduced.
- The Riccati equation associated with {A, G, C, Λ₀} can now be solved. This leads to matrices B⁰ and D⁰ which can serve as initial guesses for the NLS optimization problem.
- Both methods can be used with data sampled on a non-uniform grid of frequencies.
- The identification algorithm and the methods ensuring positivity of the spectrum are further studied in Hinnen *etal*.:2005.

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