

## TTT4120 Digital Signal Processing Fall 2017

### Linear prediction

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## Lecture in course book\*

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- Proakis, Manolakis Digital Signal Processing, 4<sup>th</sup> Ed.
  - 14.3.1 Forward linear prediction
  - 14.3.2 The Yule-Walker method for AR model parameters
- A comprehensive overview of topics treated in the lecture, see “[Introduksjon til statistisk signalbehandling](#)” on Blackboard

\*Level of detail is defined by lectures and problem sets

## Contents and learning outcomes

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- How to find the AR parameters for a general process
- Linear prediction
- How many coefficients to choose? Model order estimation

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## Estimation in practice

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- Only access to finite-length realization,  $x[n]$ , of process  $X[n]$ 
  - True  $\gamma_{XX}[l]$  must be estimated from  $x[n] \Rightarrow \hat{\gamma}_{XX}[l]$
  - Parameter values computed using  $\hat{\gamma}_{XX}[l]$  becomes parameter estimates  $\{\hat{a}_k\} \Rightarrow$  Power spectrum estimate

$$\hat{\Gamma}_{XX}(f) = \frac{\hat{\sigma}_f^2}{|\hat{A}(f)|^2} = \frac{\hat{\sigma}_f^2}{|1 + \sum_{k=1}^p \hat{a}_k e^{-j2\pi f k}|^2}$$

$\begin{aligned} \gamma_{XX}[l] &\rightarrow \{a_k\} \rightarrow \Gamma_{XX}(f) \\ &\downarrow \text{(estimation)} \\ \hat{\gamma}_{XX}[l] &\rightarrow \{\hat{a}_k\} \rightarrow \hat{\Gamma}_{XX}(f) \end{aligned}$
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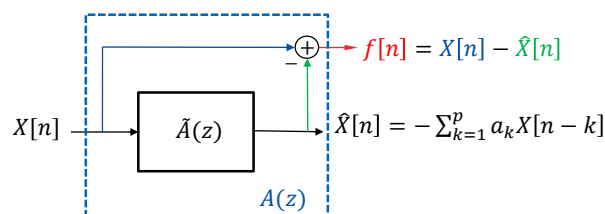
## Estimation in practice...

- In practice process  $X[n]$  may not be a true  $\text{AR}(p)$  process
  - How to choose parameters  $\{\hat{a}_k\}$  to closely model  $X[n]$  using an  $\text{AR}(p)$  process?
  - How do we measure closeness between model process and physical process?
- We will design  $p$ th-order linear predictor:
  - We observe/measure process  $X[n]$
  - Store  $p$  prior values of  $X[n]$ , i.e.,  $\{X[n-1], \dots, X[n-p]\}$
  - Make linear combination of past values to estimate of  $X[n]$

$$\hat{X}[n] = -\sum_{k=1}^p a_k X[n-k]$$

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## Linear prediction

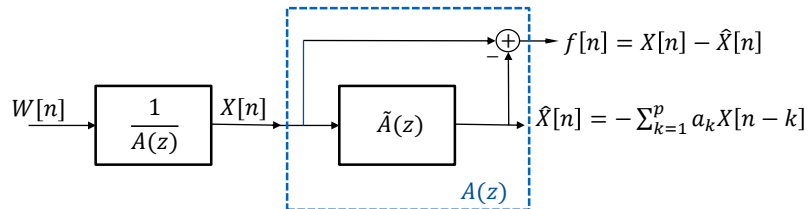


- Design  $a_k$  to match  $X[n]$  as good as possible in some sense
  - We can compute the prediction error
  - Error  $f[n]$  should be small
  - Find predictor coefficients that minimize mean-square error

$$\sigma_f^2 = E \left\{ (X[n] - \hat{X}[n])^2 \right\} = E \left\{ \left( X[n] + \sum_{k=1}^p a_k X[n-k] \right)^2 \right\}$$

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## Linear prediction...



- If  $X[n]$  is a true  $AR(p)$  process then  $f[n] = W[n]$  whenever the prediction coefficients  $a_k$  match those of the  $AR(p)$  process
- In practice this assumption leads to an approximation

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## Linear prediction...

- Elaborate the MSE

$$\begin{aligned}
 \sigma_f^2 &= E \left\{ (X[n] - \hat{X}[n])^2 \right\} = E \left\{ (X[n] + \sum_{k=1}^p a_k X[n-k])^2 \right\} \\
 &= E \left\{ X^2[n] + 2 \sum_{k=1}^p a_k X[n-k] X[n] + \sum_{l=1}^p \sum_{k=1}^p a_k a_l X[n-k] X[n-l] \right\} \\
 &= \gamma_{XX}[0] + 2 \sum_{k=1}^p a_k \gamma_{XX}[k] + \sum_{l=1}^p \sum_{k=1}^p a_k a_l \gamma_{XX}[l-k]
 \end{aligned}$$

- MSE is minimum if we choose  $a_k$  such that

$$\frac{d\sigma_f^2}{da_k} = 0, k = 1, 2, \dots, p$$

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## Linear prediction...

- Example: Find optimal predictor for  $p = 1$ , i.e.,  $\hat{X}[n] = -a_1 X[n-1]$

$$\begin{aligned}\sigma_f^2 &= E\{(X[n] - \hat{X}[n])^2\} = E\{(X[n] + a_1 X[n-1])^2\} \\ &= \gamma_{XX}[0] + 2a_1 \gamma_{XX}[1] + a_1^2 \gamma_{XX}[0] \\ &= \gamma_{XX}[0] - \frac{\gamma_{XX}^2[1]}{\gamma_{XX}[0]} + \gamma_{XX}[0] \left(a_1 + \frac{\gamma_{XX}[1]}{\gamma_{XX}[0]}\right)^2\end{aligned}$$

- Prediction error variance minimized for value  $a_1$  that gives  $\frac{d\sigma_f^2}{da_1} = 0$ :

$$\frac{d\sigma_f^2}{da_1} = 2\gamma_{XX}[1] + 2a_1 \gamma_{XX}[0] = 0 \Rightarrow a_1 = -\frac{\gamma_{XX}[1]}{\gamma_{XX}[0]}$$

- Resulting prediction variance:  $\sigma_f^2 = \gamma_{XX}[0] + a_1 \gamma_{XX}[1]$

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## Linear prediction...

- In vector notation:  $\sigma_f^2 = \gamma_{XX}[0] + 2\mathbf{a}^T \boldsymbol{\gamma}_{XX} + \mathbf{a}^T \boldsymbol{\Gamma}_{XX} \mathbf{a}$

$$\mathbf{a} = \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_p \end{bmatrix}, \boldsymbol{\Gamma}_{XX} = \begin{bmatrix} \gamma_{XX}[0] & \gamma_{XX}[1] & \dots & \gamma_{XX}[p-1] \\ \gamma_{XX}[1] & \gamma_{XX}[0] & \dots & \gamma_{XX}[p-2] \\ \vdots & \ddots & \ddots & \vdots \\ \gamma_{XX}[p-1] & \gamma_{XX}[p-2] & \dots & \gamma_{XX}[0] \end{bmatrix}, \boldsymbol{\gamma}_{XX} = \begin{bmatrix} \gamma_{XX}[1] \\ \gamma_{XX}[2] \\ \vdots \\ \gamma_{XX}[p] \end{bmatrix}$$

- Set the gradient  $\nabla_{\mathbf{a}} \sigma_f^2 = \mathbf{0}$ , i.e.,

$$\nabla_{\mathbf{a}} \sigma_f^2 = \begin{bmatrix} \frac{\partial \sigma_f^2}{\partial a_1} & \dots & \frac{\partial \sigma_f^2}{\partial a_p} \end{bmatrix}^T = [0 \quad \dots \quad 0]^T$$

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## Linear prediction...

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- $\nabla_{\mathbf{a}} \sigma_f^2 = \mathbf{0}$ :

$$\begin{aligned}\nabla_{\mathbf{a}} \sigma_f^2 &= 2\boldsymbol{\gamma}_{XX} + 2\boldsymbol{\Gamma}_{XX}\mathbf{a} = \mathbf{0} \\ \Rightarrow \mathbf{a} &= -\boldsymbol{\Gamma}_{XX}^{-1}\boldsymbol{\gamma}_{XX}\end{aligned}$$

- Minimum MSE:

$$\begin{aligned}\sigma_f^2 &= \gamma_{XX}[0] + 2\mathbf{a}^T\boldsymbol{\gamma}_{XX} + \mathbf{a}^T\boldsymbol{\Gamma}_{XX}\mathbf{a} \\ &= \gamma_{XX}[0] + 2\mathbf{a}^T\boldsymbol{\gamma}_{XX} - \mathbf{a}^T\boldsymbol{\gamma}_{XX} \\ &= \gamma_{XX}[0] + \mathbf{a}^T\boldsymbol{\gamma}_{XX} = \gamma_{XX}[0] + \sum_{k=1}^p a_k \gamma_{XX}[k]\end{aligned}$$

- Same solution as we had for a pure AR( $p$ ) process

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## Linear prediction...

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- Alternative approach by completing the square:

$$\begin{aligned}\sigma_f^2 &= \gamma_{XX}[0] + 2\mathbf{a}^T\boldsymbol{\gamma}_{XX} + \mathbf{a}^T\boldsymbol{\Gamma}_{XX}\mathbf{a} \\ &= \gamma_{XX}[0] - \boldsymbol{\gamma}_{XX}^T\boldsymbol{\Gamma}_{XX}^{-1}\boldsymbol{\gamma}_{XX} + (\mathbf{a} + \boldsymbol{\Gamma}_{XX}^{-1}\boldsymbol{\gamma}_{XX})^T\boldsymbol{\Gamma}_{XX}(\mathbf{a} + \boldsymbol{\Gamma}_{XX}^{-1}\boldsymbol{\gamma}_{XX})\end{aligned}$$

- The above holds true whenever  $\boldsymbol{\Gamma}_{XX}$  is positive definite, i.e.,

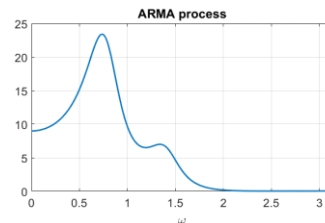
$$\mathbf{x}^T\boldsymbol{\Gamma}_{XX}\mathbf{x} > 0, \forall \mathbf{x} \neq \mathbf{0}$$

- Consequently,  $\sigma_f^2$  is minimized when last term equals zero

$$\mathbf{a} = -\boldsymbol{\Gamma}_{XX}^{-1}\boldsymbol{\gamma}_{XX}$$

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## Linear prediction...



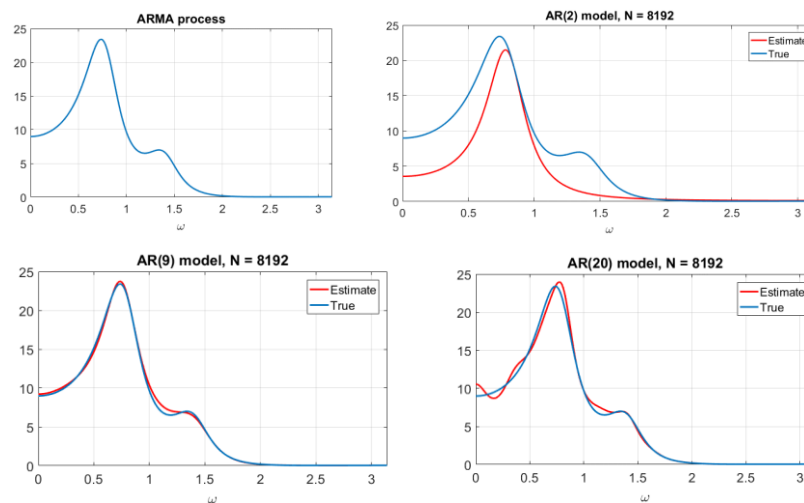
- Example: Estimate  $\Gamma_{XX}(f)$  from a realization of an  $N$ -point ARMA process,

$$X[n] = -\sum_{k=1}^p a_k X[n-k] + \sum_{k=0}^q b_k W[n-k], \quad W[n] \sim N(0, \sigma_w^2)$$

- Approximate with an AR( $p$ ) process and estimate model coefficients,  $\hat{a}_k$ , by minimizing prediction error variance,  $\sigma_f^2$ 
  - What model order should I use?

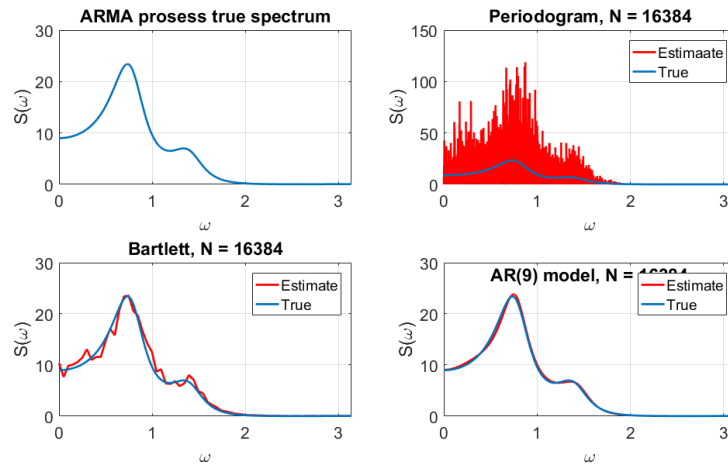
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## Linear prediction...



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## Linear prediction...



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## Determining model order $p$

- Model order not known when we shall model a physical process
- Proper choice of order  $p$  is necessary for good modelling capability
  - Too small  $p$  leads to smoothened spectrum
  - Too large  $p$  leads to spurious low-level peaks in the spectrum
- Prediction variance  $\sigma_f^2(p)$  could be an indicator
  - Monotonically decreasing with  $p$
  - Need to decide when changes are sufficiently small
  - Usually imprecise: in general no clear knee visible in plot  $\sigma_f^2(p)$

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## Determining model order $p$

- Different criteria that penalizes high model order  $p$ :

$$\text{FPE}(p) = \sigma_f^2(p) \frac{N + p + 1}{N - p - 1}$$

$$\text{MDL}(p) = N \log \sigma_f^2(p) + p \log N$$

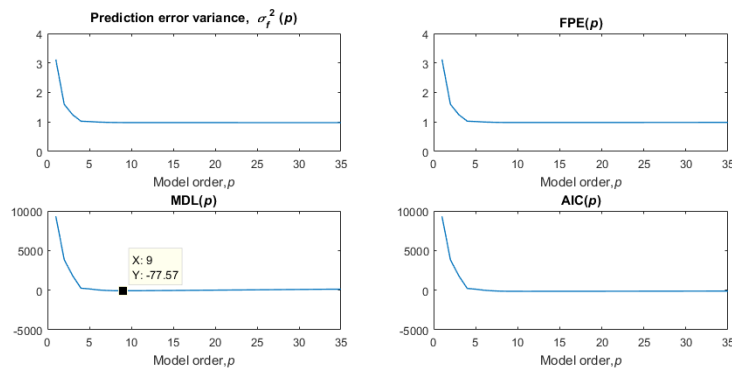
$$\text{AIC}(p) = N \log \sigma_f^2(p) + 2p$$

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## Determining model order $p...$

- Example: Estimate  $\Gamma_{XX}(f)$  from a realization of an ARMA process

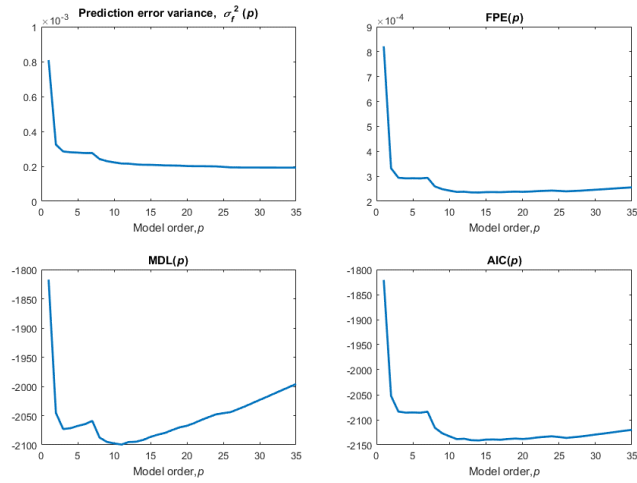
$$X[n] = -\sum_{k=1}^p a_k X[n-k] + \sum_{k=0}^q b_k W[n-k], W[n] \sim N(0, \sigma_w^2)$$



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## Determining model order $p$ ...

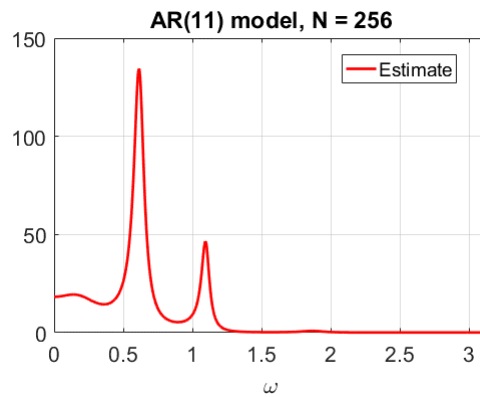
- Example: Vowel 'æ',  $N = 256$ :



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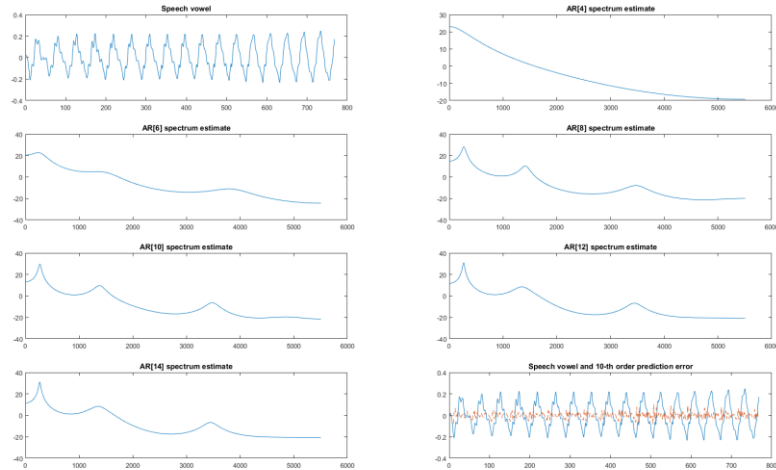
## Determining model order $p$ ...

- Example: Vowel 'æ',  $N = 256$ :



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## Determining model order $p...$



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## Final notes on estimation in practice

- All methods looked at so far assume
  - Random processes to be stationary and ergodic
  - Random processes are autoregressive (AR)
- In practice, all physical processes of interest are nonstationary
  - Short-time stationarity: process varies slowly and within a certain time window, statistical properties are constant
  - Assume stationarity over  $M$  times and we need  $N < M$  points
- Other methods for finding estimates
  - Usually lead to similar performance. Main differences are in the performance with few data points

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## Summary

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- Today we discussed:
  - Linear prediction
  - Model order
- Next time:
  - FIR filter design

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