

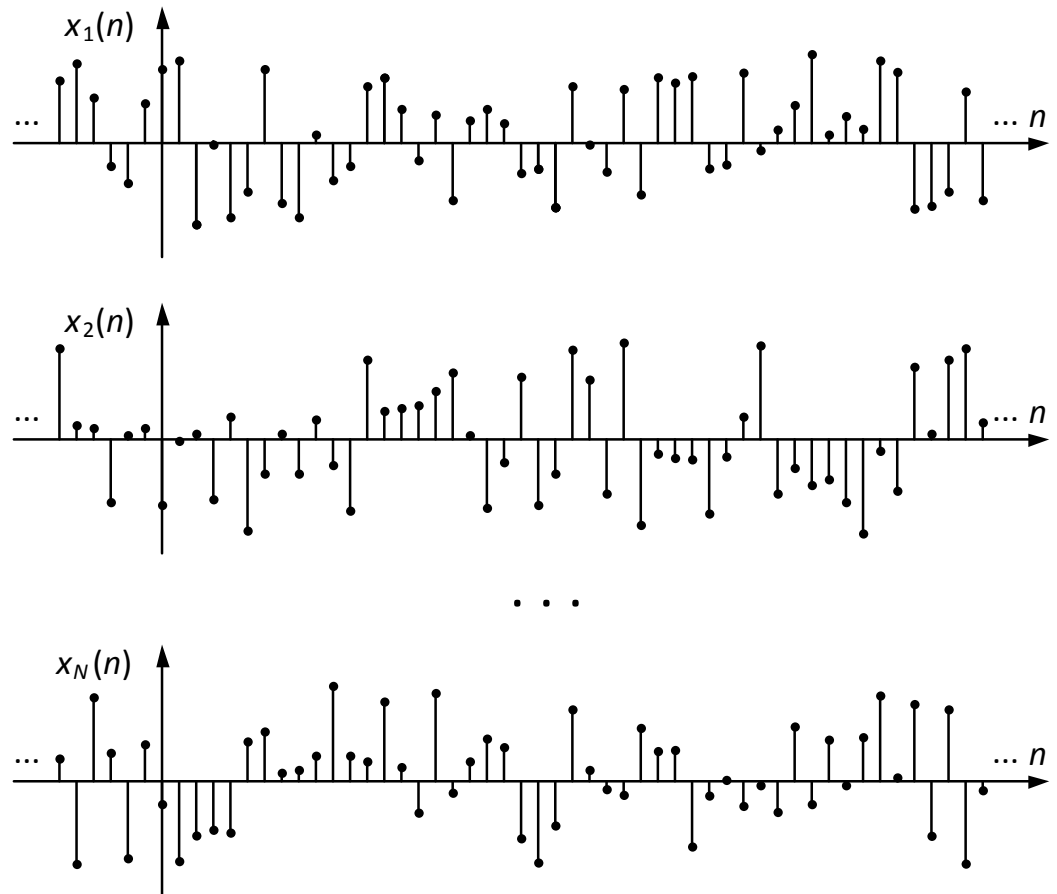
# **SPECTRAL ANALYSIS OF DISCRETE-TIME SIGNALS**

# Random discrete signals

- Exact values are unknown outside of the range in which the signal is observed
- Described through statistical parameters
  - This approach is sensible for many real-world signals, which occur as a consequence of mutual interaction of a number a complex causes
- A random signal is one specific realization of a *random process*
  - A *discrete-time random process*  $X_n$  represents a time series of random variables

# Statistical ensemble

- A set of all possible realizations of a random process
  - different signals but same statistical properties



# Random variables and random processes

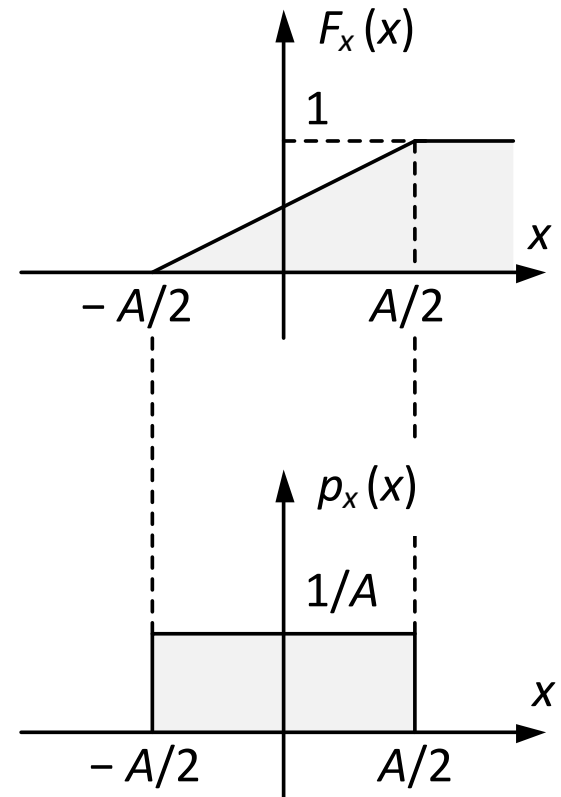
- A random variable can be described through its *cumulative probability distribution function* (CDF):

$$F_x(x) = P(X \leq x)$$

and if it is continuous, it can also be described through its *probability density function* (PDF):

$$p_x(x) = \frac{dF_x(x)}{dx}$$

- In this way we can also describe any random process at a particular moment



# Random processes

- A random process at a moment  $n$  can be described using a CDF dependent on  $n$ :

$$F_{X_n}(x, n) = P(X_n \leq x)$$

and if it is continuous, it can also be described using a PDF dependent on  $n$ :

$$p_{X_n}(x, n) = \frac{dF_{X_n}(x, n)}{dx}$$

- The values of  $X_n$  are generally mutually dependent, and to fully describe a random process we should know a joint CDF for any subset of moments:

$$F_{X_{n_1} X_{n_2} \dots X_{n_N}}(x_1, x_2, \dots, x_N, n_1, n_2, \dots, n_N) = P(X_{n_1} \leq x_1, X_{n_2} \leq x_2, \dots, X_{n_N} \leq x_N)$$

- In practice this function is most often unknown, and we describe random processes using statistical parameters such as mean and autocorrelation function

# Stationarity of a random process

- A random process is (*strict-sense*) *stationary* if its joint CDF does not change in time, i.e. if, for any integer  $k$ , the following holds:

$$F_{X_{n_1} X_{n_2} \dots X_{n_N}}(x_1, x_2, \dots, x_N, n_1, n_2, \dots, n_N) \\ = F_{X_{n_1+k} X_{n_2+k} \dots X_{n_N+k}}(x_1, x_2, \dots, x_N, n_1 + k, n_2 + k, \dots, n_N + k)$$

- There are many random processes for which this does not hold, and which go through various *states*, in which their statistical properties may be quite different

# Mean and autocorrelation function

- The *mean* of a random process is its mathematical expectation:

$$m_{X_n}(n) = E\{X_n\} = \int_{-\infty}^{\infty} x p_{X_n}(x, n) dx$$

and if it does not depend on  $n$ , the process is *stationary with respect to its mean*

- The *autocorrelation function* of a random process is the measure of the similarity of its values at the time instants  $n$  and  $n+k$ :

$$r_{X_n X_{n+k}}(n, n+k) = E\{X_n X_{n+k}\} = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} x_1 x_2 p_{X_n X_{n+k}}(x_1, x_2, n, n+k) dx_1 dx_2$$

and if it depends solely on  $k$ , the process is *stationary with respect to its autocorrelation function*

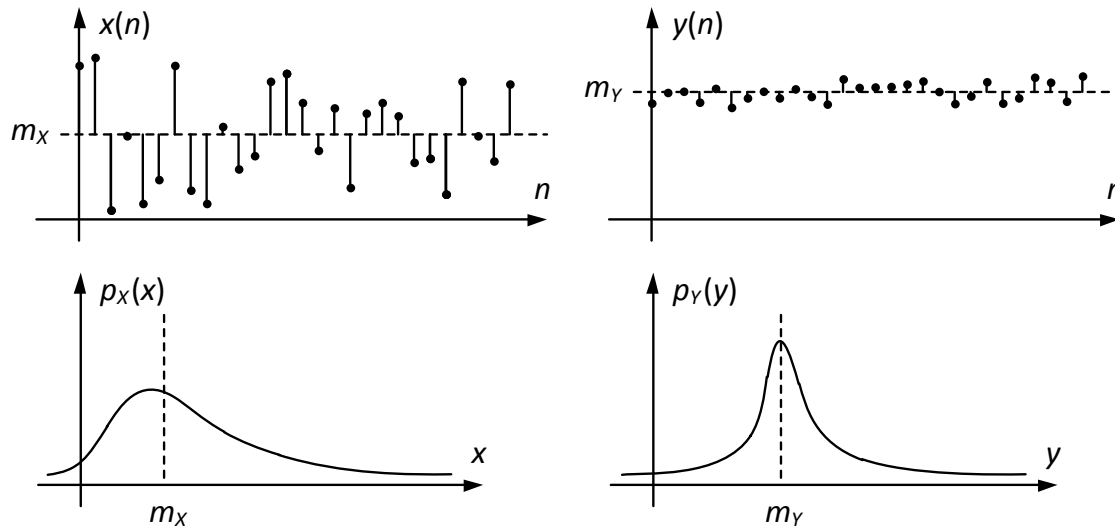
- A process stationary both in terms of its mean and its autocorrelation function is considered *wide-sense stationary*

# Variance

- The variance of a wide-sense stationary (WSS) random process is the measure of its expected deviation from its mean:

$$\sigma_X^2 = E\{(X_n - m_X)^2\} = \int_{-\infty}^{\infty} (x - m_X)^2 p_{X_n}(x, n) dx$$

$$\sigma_X^2 > \sigma_Y^2$$



- If  $m_X = 0$ , the variance is equal to the *average power* of the signal



# Estimation of statistical parameters

- We usually have only one member of the statistical ensemble  $x(n)$  at our disposal, from which we should infer the values of the statistical parameters of a WSS random process
- The mean of the discrete signal  $x(n)$  is:

$$\overline{x(n)} = \lim_{N \rightarrow \infty} \frac{1}{2N + 1} \sum_{n=-N}^N x(n)$$

and its autocorrelation is:

$$r_{xx}(k) = \lim_{N \rightarrow \infty} \frac{1}{2N + 1} \sum_{n=-N}^N x(n)x(n+k)$$

and if the process is *ergodic*, they are equal to the ensemble mean and autocorrelation function

- In practice we cannot calculate either of these parameters because we would need to know infinitely many values of signal samples and calculate infinite sums

# Examples of parameter estimation

## Mean estimation

- The standard estimate of the mean is:

$$\hat{m}_x = \frac{1}{N} \sum_{n=0}^{N-1} x(n)$$

- This estimate is *unbiased* (its expectation is equal to the actual mean  $m_x$ ):

$$E\{\hat{m}_x\} = \frac{1}{N} E\left\{\sum_{n=0}^{N-1} X_n\right\} = \frac{1}{N} \sum_{n=0}^{N-1} E\{X_n\} = m_x$$

- This estimate is also *consistent* (variance goes to 0 when  $N$  goes to infinity)
  - Consistency is a highly desirable property of an estimate because with a consistent estimate we can expect the same result from any member of the ensemble

# Examples of parameter estimation

## Estimation of the autocorrelation function

- The standard estimate of the autocorrelation function is:

$$\hat{r}_{xx}(k) = \frac{1}{N} \sum_{n=0}^{N-|k|-1} x(n)x(n+k)$$

- This estimate is *biased* (its expectation is not equal to the actual autocorrelation function  $r_{xx}(k)$ ):

$$E\{\hat{r}_{xx}(k)\} = \frac{1}{N} \sum_{n=0}^{N-|k|-1} E\{X_n X_{n+k}\} = \frac{1}{N} \sum_{n=0}^{N-|k|-1} r_{xx}(k) = \left(1 - \frac{|k|}{N}\right) r_{xx}(k)$$

but it is *asymptotically unbiased* (its expectation converges to  $r_{xx}(k)$  as  $N$  approaches infinity)

- This estimate is *consistent*

# Spectrum of a discrete-time random process

- Discrete-time random signals have infinite energy and thus have no DTFT, and we cannot calculate the energy spectral density (which is equal to the square of the magnitude spectrum)
- Discrete-time random signals have finite *average power*:

$$P_x = \lim_{N \rightarrow \infty} \frac{1}{2N + 1} \sum_{n=-N}^N |x(n)|^2$$

and it is possible to define their *power spectral density (power spectrum)*, by analogy with a corresponding definition for deterministic signals:

$$S_x(\omega) = \lim_{N \rightarrow \infty} E \left\{ \frac{1}{2N + 1} \left| \sum_{n=-N}^N x_n e^{-j\omega n} \right|^2 \right\}$$

# Spectrum of a discrete-time random process

- Power spectral density (PSD) is equal to the Fourier transform of the autocorrelation function:

$$S_X(\omega) = \sum_{k=-\infty}^{\infty} r_{XX}(k)e^{-j\omega k}$$

and it is a real, non-negative function, and in the case of real signals also even

- The estimation of power spectral density is one of the most frequent and most important tasks in digital signal processing
  - The problem is that we usually have only one member of the statistical ensemble at our disposal, and we usually know the values of only a limited number of its samples

# Methods for power spectrum estimation

## Non-parametric (classical) methods

- The spectrum is directly calculated by applying DTFT to either a finite-length segment of the input signal or its autocorrelation
- We do not introduce any assumptions as to the shape of the power spectrum

## Parametric methods

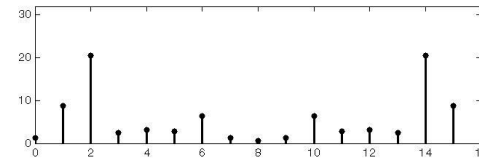
- We introduce a *model* of power spectral density function and the problem amounts to the estimation of the parameters of this model
- If an adequate model is used, these methods can give much better results than non-parametric ones, particularly if only a short segment of the input signal is available

# Non-parametric methods for estimating $S_X(\omega)$

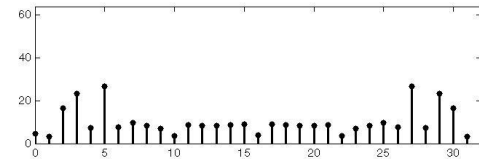
- The estimation of  $S_X(\omega)$  is more accurate if a longer segment of the input signal is available (as was the case with deterministic signals)

Example: estimation of the power spectrum of a signal composed from two sinusoids contaminated with noise

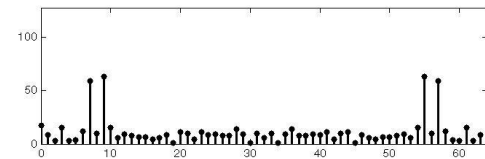
$N = 16$



$N = 32$



$N = 64$



- If  $N$  is too small, we can increase the density of DFT samples by zero padding, although this does not affect frequency resolution
  - Zero padding, as a means for increasing the density of DFT samples is of particular importance for non-stationary signals

# Periodogram

- Direct estimate based on the definition of  $S_x(\omega)$ , with the time range limited to the interval from 0 to  $N-1$ :

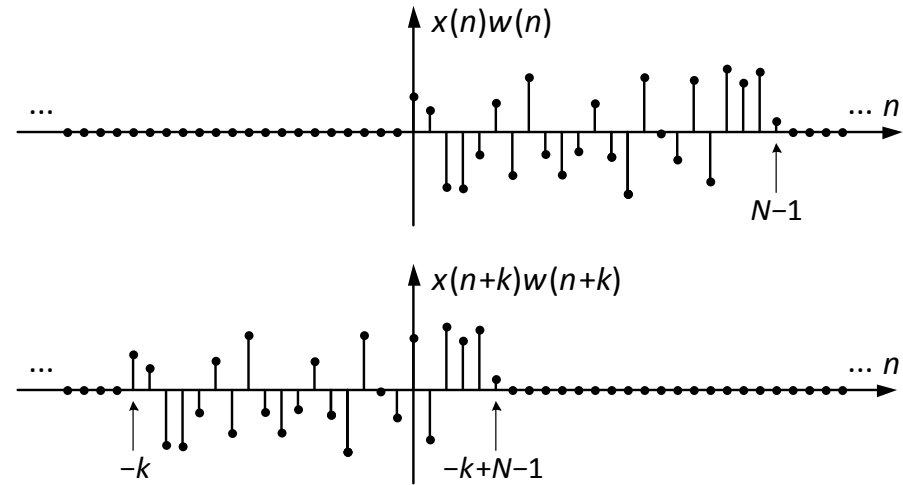
$$S_x(\omega) = \lim_{N \rightarrow \infty} E \left\{ \frac{1}{2N+1} \left| \sum_{n=-N}^N x_n e^{-j\omega n} \right|^2 \right\} \rightarrow \hat{S}_x(\omega) = \frac{1}{N} \left| \sum_{n=0}^{N-1} x(n) e^{-j\omega n} \right|^2$$

which is a real, non-negative function, and in the case of real signals also even, just as was the case for  $S_x(\omega)$



# Periodogram

$$\begin{aligned}
 \hat{S}_x(\omega) &= \frac{1}{N} |\bar{X}(\omega)|^2 = \frac{1}{N} \bar{X}(\omega) \bar{X}^*(\omega) \\
 &= \frac{1}{N} \sum_{n=0}^{N-1} x(n) e^{-j\omega n} \sum_{m=0}^{N-1} x(m) e^{j\omega m} \\
 &= \frac{1}{N} \sum_{n=0}^{N-1} \sum_{m=0}^{N-1} x(n) x(m) e^{-j\omega(m-n)} \\
 &= \sum_{k=-N+1}^{N-1} \left( \frac{1}{N} \sum_{n=0}^{N-|k|-1} x(n) x(n+k) \right) e^{-j\omega k} \\
 &= \sum_{k=-N+1}^{N-1} \hat{r}_{xx}(k) e^{-j\omega k}
 \end{aligned}$$



- Periodogram represents the Fourier transform of the standard estimate of the autocorrelation function of a random process

# Properties of the periodogram

$$\begin{aligned} E\{\hat{S}_x(\omega)\} &= E\left\{\sum_{k=-N+1}^{N-1} \hat{r}_{xx}(k) e^{-j\omega k}\right\} = \sum_{k=-N+1}^{N-1} E\{\hat{r}_{xx}(k)\} e^{-j\omega k} \\ &= \sum_{k=-N+1}^{N-1} \left(1 - \frac{|k|}{N}\right) r_{xx}(k) e^{-j\omega k} \end{aligned}$$

- The expectation of the periodogram  $\hat{S}_x(\omega)$  is not equal to the actual power spectrum  $S_x(\omega) = \text{DTFT}\{r_{xx}(k)\}$ , but to  $\text{DTFT}\{r_{xx}(k)w_B(k)\}$ , where  $w_B(k)$  is the triangular window function
  - Periodogram is a *biased* estimate of  $S_x(\omega)$ , but still *asymptotically unbiased*
- When estimating  $S_x(\omega)$  on the basis of a periodogram, the problem of *spectral leakage* occurs

$$E\{\hat{S}_x(\omega)\} = S_x(\omega) \circledast W_B(\omega) \quad W_B(\omega) = \frac{1}{N} \left( \frac{\sin N\omega/2}{\sin \omega/2} \right)^2$$

# Properties of the periodogram

- Variance of the periodogram  $\hat{S}_x(\omega)$  for large values of  $N$  in most cases of practical interest is proportional to the square of  $S_x(\omega)$

$$\lim_{N \rightarrow \infty} \text{Var} \left\{ \hat{S}_x(\omega) \right\} \sim S_x^2(\omega)$$

- Periodogram is an *inconsistent* estimate of  $S_x(\omega)$ , which means that it can greatly depend on the realization of the random process
- Due to its inconsistency, periodogram is not considered to be a high quality estimate of  $S_x(\omega)$
- An idea: to split the available segment of the input signal into  $K$  equal parts, to calculate the periodogram on each of them and to calculate the average of the results

# Periodogram averaging

- Let the available segment of the input signal  $x(n)$  of length  $N$  be divided into  $K$  equal parts of length  $M = N/K$ , with no overlap (basic version, proposed by Bartlett)
- Let  $\hat{S}_x^{(i)}(\omega)$  be the periodogram calculated on the  $i$ -th segment:

$$\hat{S}_x^{(i)}(\omega) = \frac{1}{M} \left| \sum_{n=0}^{M-1} x_i(n) e^{-j\omega n} \right|^2$$

- The averaged periodogram is obtained as the average of thus obtained  $K$  periodograms:

$$\hat{S}_x^B(\omega) = \frac{1}{K} \sum_{i=0}^{K-1} \hat{S}_x^{(i)}(\omega)$$

# Properties of the averaged periodogram

- Mathematical expectation of  $\hat{S}_x^B(\omega)$  is equal to the mathematical expectation of a single periodogram  $\hat{S}_x^{(i)}(\omega)$  :

$$E \left\{ \hat{S}_x^B(\omega) \right\} = \sum_{k=-M+1}^{M-1} \left( 1 - \frac{|k|}{M} \right) r_{xx}(k) e^{-j\omega k} = S_x(\omega) \otimes W_B(\omega)$$

and since the triangular window function is now  $K$  times shorter, the frequency resolution is also  $K$  times worse

- The variance is reduced by a factor of  $K$  with respect to the standard periodogram (the variance of the mean of  $K$  identically distributed random variables is  $K$  times smaller than the variance of a single one of them):

$$\lim_{N \rightarrow \infty} \text{Var} \left\{ \hat{S}_x(\omega) \right\} \sim \frac{1}{K} S_x^2(\omega)$$

- By averaging in this way, we have improved the consistency of the estimate at the expense of the frequency resolution

# Modifications of the averaged periodogram

- The basic idea of Bartlett can be modified in two ways (both proposed by Welch):
  - Consecutive segments can overlap
    - This increases the number of segments and decreases variance further
    - Overlap should not be too significant because that would undermine the assumption of independence between consecutive segments (if the overlap is above 50% the variance will abruptly stop decreasing)
  - Every segment  $x_i(n)$  can be shaped by applying a non-rectangular window function before calculating the periodogram
    - This reduces the correlation between consecutive segments, and also mitigates the problem of lower reliability of the estimate of autocorrelation function for large  $k$
- Without affecting the frequency resolution, the variance is now lower: for a large  $N$  it still behaves as  $S_x^2(\omega)/K$ , but  $K$  is now greater owing to the overlap

# Periodogram smoothing

- The idea is based on windowing the estimate of the autocorrelation function (as proposed by Blackman and Tukey):

- This reduces the influence of less reliable values of  $\hat{r}_{xx}(k)$ , obtained for large  $|k|$

$$\hat{S}_x^{BT}(\omega) = \text{FTD}\{\hat{r}_{xx}(k)w(k)\} = \sum_{k=-K+1}^{K-1} \hat{r}_{xx}(k)w(k)e^{-j\omega k}$$

- Mathematical expectation of a thus smoothed periodogram is:

$$\begin{aligned} E\{\hat{S}_x^{BT}(\omega)\} &= \text{FTD}\{E\{\hat{r}_{xx}(k)\}w(k)\} = \text{FTD}\{\hat{r}_{xx}(k)w_B(k)w(k)\} \\ &= S_x(\omega) \circledast W_B(\omega) \circledast W(\omega) = E\{\hat{S}_x(\omega)\} \circledast W(\omega) \end{aligned}$$

where  $w(k)$  is usually much shorter than  $w_B(k)$ , and thus the influence of convolution with  $W_B(\omega)$  is negligible

- The resulting spectral leakage depends on  $W(\omega)$ 
  - The choice of  $W(\omega)$  affects the trade-off between the variance and the frequency resolution

# Parametric methods for estimating $S_x(\omega)$

- We introduce a *model* of the random process, and the problem of estimating  $S_x(\omega)$  amounts to the problem of estimating model parameters
- We assume that the signal  $x(n)$  is the output of an LTI system with transfer function:

$$H(z) = \frac{B(z)}{A(z)} = \frac{\sum_{i=0}^q b_i z^{-i}}{1 + \sum_{i=1}^p a_i z^{-i}}$$

where the input is white noise  $s(n)$  of zero mean and variance (average power) equal to  $\sigma^2$

- The term *white* implies that the samples of noise are mutually uncorrelated, i.e.:

$$r_{ss}(k) = \sigma^2 \delta(k)$$

- This reduces the problem to the modeling of a *system*
- Power spectral density of the output signal is equal to:  $S_x(\omega) = \sigma^2 \frac{|B(\omega)|^2}{|A(\omega)|^2}$



# Parametric methods for estimating $S_X(\omega)$

- First we need to choose the values for  $p$  and  $q$ , and then to estimate  $a_i$  and  $b_i$ 
  - $q = 0$ : AR (autoregressive) models
  - $p = 0$ : MA (*moving average*) models
  - $q \neq 0, p \neq 0$ : ARMA models
- The choice is usually based on some previous knowledge of the random process
  - Owing to their simplicity, AR models are most commonly used (the others yield non-linear systems of equations)
  - An example of a problem where AR is the obvious solution is the signal which is composed of a sinusoid of an unknown frequency and noise
    - Poles of the model should match the unknown frequency of the sinusoid

$$\frac{X(z)}{S(z)} = \frac{\sum_{i=0}^q b_i z^{-i}}{1 + \sum_{i=1}^p a_i z^{-i}}$$

# Estimation of $S_x(\omega)$ for AR random processes

$$x(n) = -\sum_{i=1}^p a_i x(n-i) + s(n)$$

$$E\{x(n)x(n-k)\} = -\sum_{i=1}^p a_i E\{x(n-i)x(n-k)\} + E\{s(n)x(n-k)\}$$

$$r_{xx}(k) = -\sum_{i=1}^p a_i r_{xx}(k-i) + E\left\{s(n) \sum_{m=0}^{\infty} h(m) s(n-k-m)\right\}$$

$$= -\sum_{i=1}^p a_i r_{xx}(k-i) + \sum_{m=0}^{\infty} h(m) E\{s(n)s(n-k-m)\}$$

$$= -\sum_{i=1}^p a_i r_{xx}(k-i) + \sum_{m=0}^{\infty} h(m) \sigma^2 \delta(m+k)$$

$$= \begin{cases} -\sum_{i=1}^p a_i r_{xx}(k-i), & k > 0 \\ -\sum_{i=1}^p a_i r_{xx}(k-i) + \sigma^2, & k = 0 \end{cases}$$

we assume  $k \geq 0$ , and the values for negative  $k$  can be easily found because  $r_{xx}(k)$  is even

$h(m)$  exists only for  $m \geq 0$ , and thus for  $k \geq 0$  the only case when the right sum is non-zero is when  $k = m = 0$

we have used  $h(0) = 1$ , which follows from  $B(z) = 1$

# Estimation of $S_x(\omega)$ for AR random processes

- Coefficients  $a_i$  and  $\sigma^2$  are the solutions of the Yule-Walker equations:

$$\begin{bmatrix} r_{xx}(0) & r_{xx}(-1) & \dots & r_{xx}(-p) \\ r_{xx}(1) & r_{xx}(0) & \dots & r_{xx}(-p+1) \\ \vdots & \vdots & \ddots & \vdots \\ r_{xx}(p) & r_{xx}(p-1) & \dots & r_{xx}(0) \end{bmatrix} \begin{bmatrix} 1 \\ a_1 \\ \vdots \\ a_p \end{bmatrix} = - \begin{bmatrix} \sigma^2 \\ 0 \\ \vdots \\ 0 \end{bmatrix}$$

- The system matrix is a symmetrical Toeplitz matrix (each descending diagonal from left to right is constant), and thus the system can be solved efficiently, in  $N^2$  instead of  $N^3$  time, using the Levinson-Durbin algorithm
  - The same matrix is obtained if the problem is posed as a *linear prediction* problem, i.e. if the following is assumed:

$$\hat{x}(n) = - \sum_{i=1}^p a_i x(n-i)$$

- In practice the autocorrelation function  $r_{xx}(k)$  is not available and we only have its estimate, but the values of this estimate for  $k \leq p$  are still much more reliable than the values for relatively large  $k$