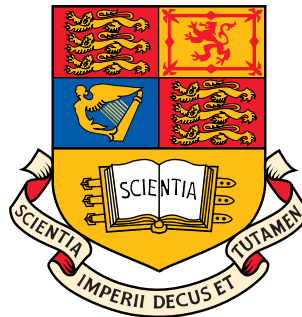

Advanced Signal Processing

Lecture 1: Random Variables

Prof Danilo Mandic
room 813, ext: 46271



Department of Electrical and Electronic Engineering
Imperial College London, UK

d.mandic@imperial.ac.uk, URL: www.commsp.ee.ic.ac.uk/~mandic

Introduction \rightsquigarrow Recap

Discrete Random Signals:

discrete vs. **digital** \leftrightarrow **quantisation**

- $\{x[n]\}_{n=0:N-1}$ is a sequence of indexed random variables $x[0], x[1], \dots, x[N-1]$, and the symbol ' $[\cdot]$ ' indicates the random nature of signal x (*every sample is random too!*)
- The sequence is discrete with respect to sample index n (*discrete time or some other physical variable, such as spatial index in arrays of sensors*)
- $x[n]$ real or complex with discrete or continuous values

NB: signals can be continuous or discrete in *time* as well as *amplitude*.

Digital signal = discrete in time and amplitude

Discrete-time signal = discrete in time, amplitude either discrete or continuous

Standardisation and normalisation

(e.g. to be invariant of amplifier gain or the quality of sensor contact)

Some real-world applications require data of specific mean and variance, yet measured variables are usually of different natures and magnitudes. We refer to **standardisation** as the process of converting the data to an arbitrary mean $\bar{\mu}$ and variance $\bar{\sigma}^2$, and to **normalisation** as the particular case $\bar{\mu} = 0$, $\bar{\sigma}^2 = 1$. In practice, **raw data** $\{x[n]\}_{n=0:N-1}$ are normalised by subtracting the sample mean, μ , and dividing by the sample std. dev., σ

- **Compute statistics:** $\mu = \frac{1}{N} \sum_{n=0}^{N-1} x[n]$, $\sigma^2 = \frac{1}{N} \sum_{n=0}^{N-1} (x[n] - \mu)^2$
- **Centred data:** $x^c = x - \mu$
- **Centred and scaled data (normalised):** $x^{cs} = \frac{x^c}{\sigma}$ ($\mu = 0, \sigma = 1$)

Normalised data can be **standardised** to any mean $\bar{\mu}$ and variance $\bar{\sigma}^2$ by

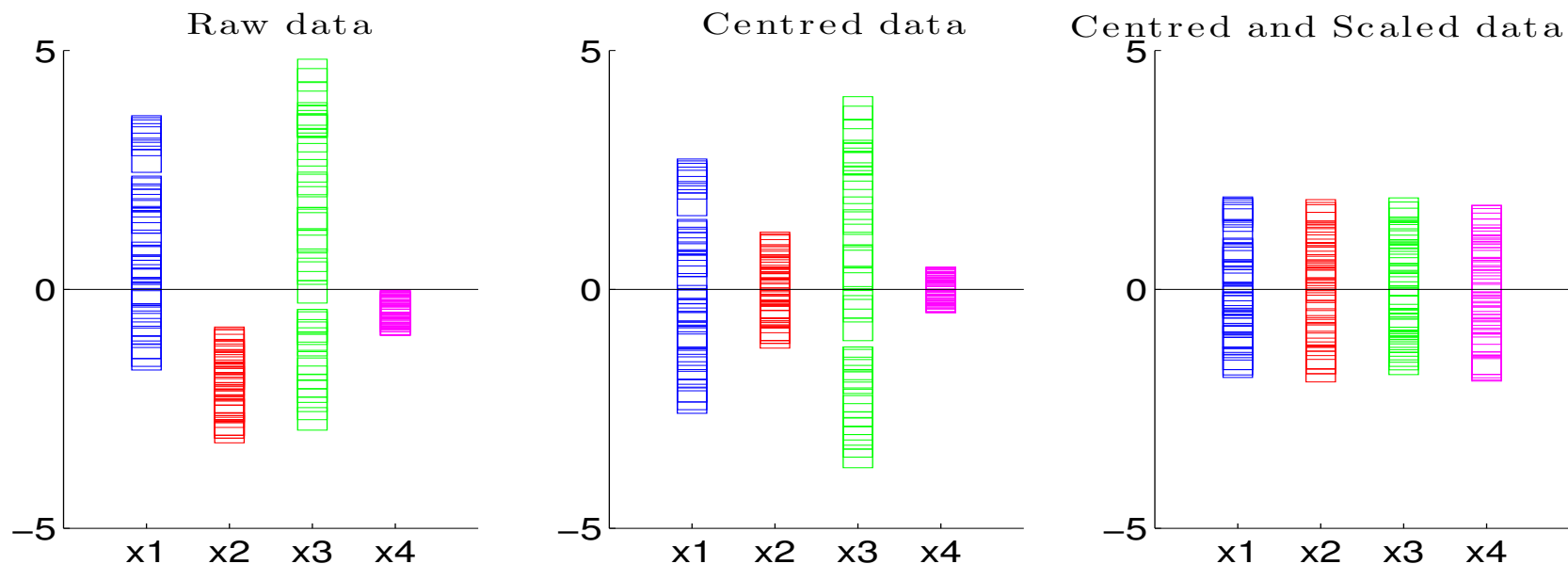
$$x^{st} = \frac{x^{cs} - \bar{\mu}}{\bar{\sigma}}$$

Standardize to zero mean and range $[-1, 1]$ $\leadsto x(n) = \frac{x(n) - (x_{max} + x_{min})/2}{(x_{max} - x_{min})/2}$

Standardisation: Example

The bars denote the amplitudes of the samples of signals x_1-x_4

For the raw measurements: $\{x_1[n], x_2[n], x_3[n], x_4[n]\}_{n=1:N}$



- Standardisation allows for a *coherent and aligned* handling of different variables, as the amplitude plays a role in regression algorithms.
- Furthermore, input variable selection can be performed by assigning smaller or larger weighting to samples (confidence intervals).

How do we describe a signal? (statistically)

Probability distribution functions – very convenient!

- **Cumulative Density Function (CDF)** – probability of a random variable falling within a given range.

$$F_X(x[n]) = \text{Probability}(X[n] \leq x[n]) \quad (1)$$

$X[n]$ – random quantity, $x[n]$ – particular fixed value.

- **Probability Density Function (pdf)** – relative likelihood for a random variable to occur at a given point in the observation space.

$$p(x[n]) = \frac{\partial F_X(x[n])}{\partial x[n]} \quad \Leftrightarrow \quad F(x) = \int_{-\infty}^x p(X) dX \quad (2)$$

NB: For random signals, for two time instants n_1 and n_2 , the pdf of $x[n_1]$ need not be identical to that of $x[n_2]$. (e.g. $\sin(n) + w(n)$)

Statistical distributions: Uniform distribution

Important: Recall that probability densities sum up to unity

$$\int_{-\infty}^{\infty} p(x[n]) dx[n] = 1$$

and that the connection between pdf and its *cumulative density function* CDF is

$$F(x[n]) = \int_{-\infty}^{x[n]} p(z) dz, \quad \text{also} \quad \lim_{x[n] \rightarrow \infty} F(x[n]) = 1$$

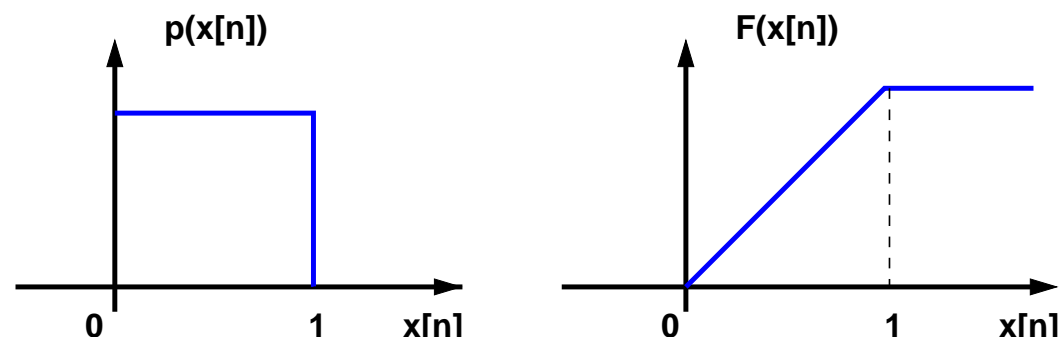
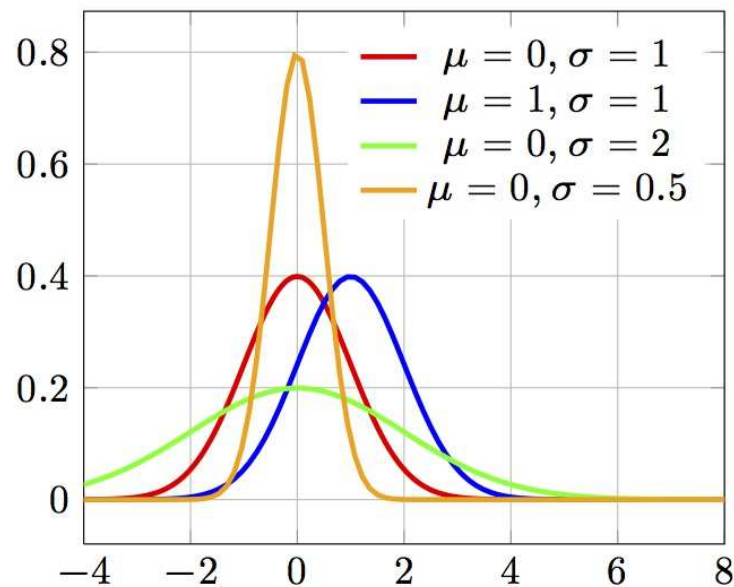


Figure: pdf and CDF for uniform distribution. In MATLAB – function rand

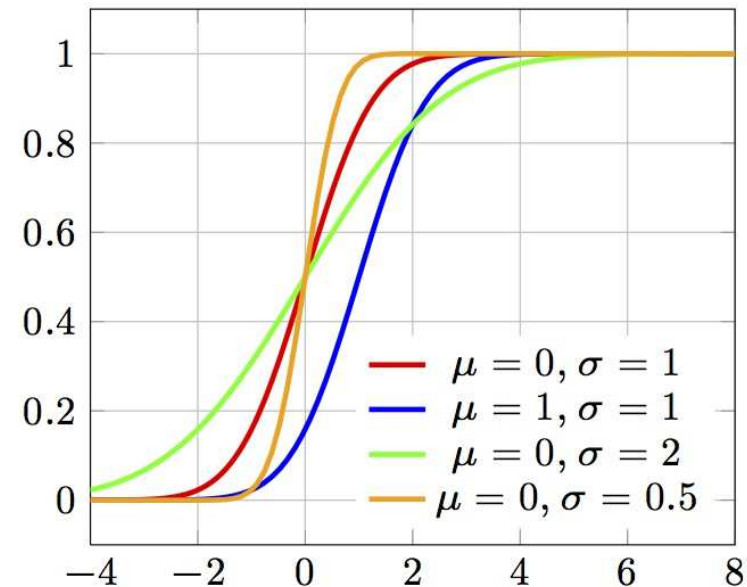
Gaussian probability and cumulative density functions

How does the variance σ^2 influence the shape of CDP and pdf?

Gaussian pdf



Gaussian CDF



$$p(x; \mu, \sigma) = \frac{1}{\sigma\sqrt{2\pi}} \exp\left(-\frac{(x - \mu)^2}{2\sigma^2}\right)$$

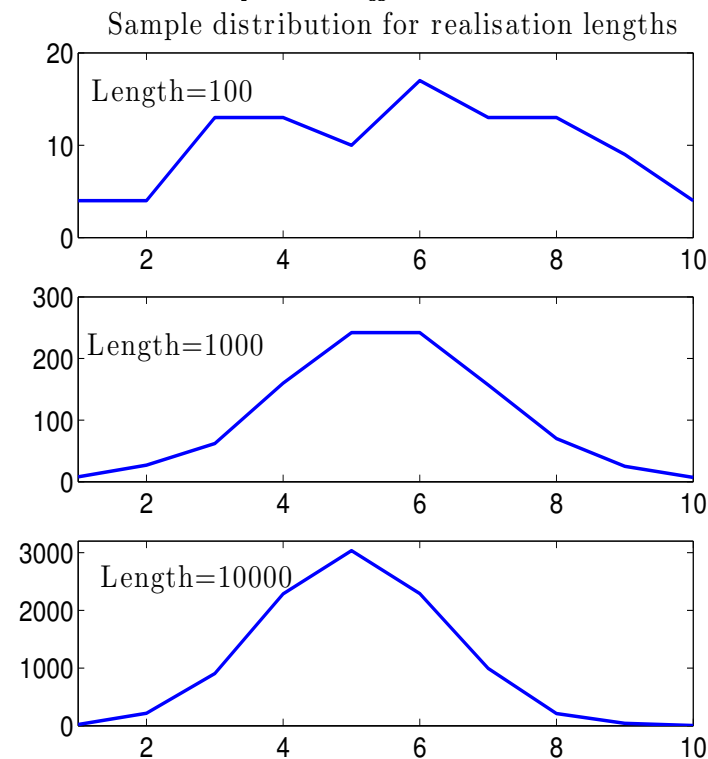
$$P(x; \mu, \sigma) = \frac{1}{2} \left[1 + \operatorname{erf}\left(\frac{x - \mu}{\sqrt{2}\sigma}\right) \right]$$

The **standard Gaussian** distribution ($\mu = 0, \sigma = 1$) is given by $p(x) = \frac{1}{\sqrt{2\pi}} \exp\left(-\frac{x^2}{2}\right)$

Statistical distributions: Gaussian \rightarrow randn in Matlab

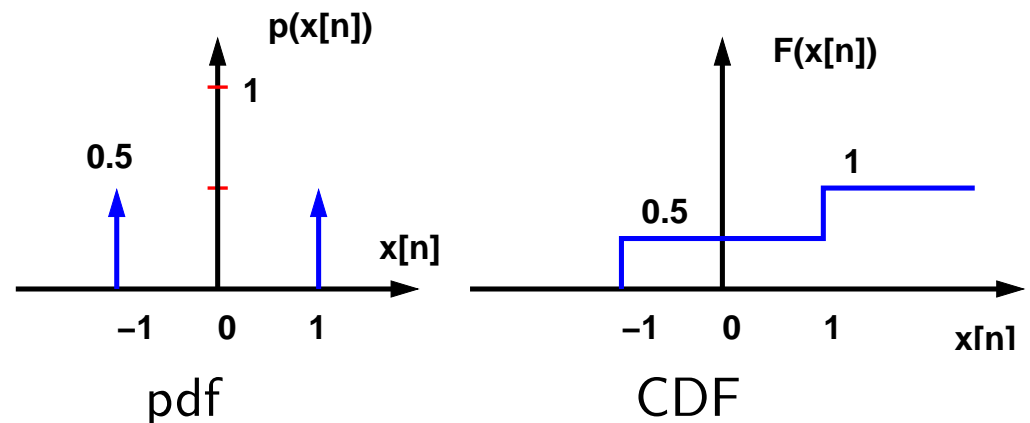
Very convenient (mathematical tractability) - especially in terms of **log-likelihood** $\log p(x[n])$

$$p(x[n]) = \frac{1}{\sqrt{2\pi\sigma_x^2}} e^{-\frac{(x[n] - \mu_x)^2}{2\sigma_x^2}} \Rightarrow \log p(x[n]) = -\frac{(x[n] - \mu_x)^2}{2\sigma_x^2} - \frac{1}{2} \log(2\pi\sigma_x^2)$$



$$x[n] \sim \mathcal{N}(\mu_x, \sigma_x^2) \quad \mu_x \rightarrow \text{mean}, \sigma_x^2 \rightarrow \text{variance}$$

Bipolar distribution



Sample densities for varying N

Multi-dimensionality versus multi-variability

Univariate vs. **Multivariate** vs. **Multidimensional**

- Single input single output (SISO) e.g. single-sensor system
- Multiple input multiple output (MIMO) (arrays of transmitters and receivers) *can measure one source with many sensors*
- Multidimensional processes (3D inertial body motion sensors, radar, vector fields, wind anemometers) – *intrinsically multidimensional*

Example: Multivariate function with single output (MISO)

$$stockvalue = f(stocks, oilprice, GNP, month, \dots)$$

⇒ Complete probabilistic description of $\{x[n]\}$ is given by its pdf

$$p(x[n_1], \dots, x[n_k]) \quad \text{for all } k \text{ and } n_1, \dots, n_k.$$

Much research is being directed towards the reconstruction of the process history from observations of one variable only (Takens)

Joint distributions of delayed samples (temporal)

Joint distribution (bivariate CDF)

$$F(x[n_1], x[n_2]) = \text{Prob}(X[n_1] \leq x[n_1], X[n_2] \leq x[n_2])$$

and its pdf

$$p(x[n_1], x[n_2]) = \frac{\partial^2 F(x[n_1], x[n_2])}{\partial x[n_1] \partial x[n_2]}$$

A k-th order multivariate CDF distribution

$$F(x[n_1], x[n_2], \dots, x[n_k]) = \text{Prob}(X[n_1] \leq x[n_1], \dots, X[n_k] \leq x[n_k])$$

and its pdf

$$p(x[n_1], x[n_2], \dots, x[n_k]) = \frac{\partial^k F(x[n_1], \dots, x[n_k])}{\partial x[n_1] \cdots \partial x[n_k]}$$

Mathematically simple, but complicated to evaluate in reality

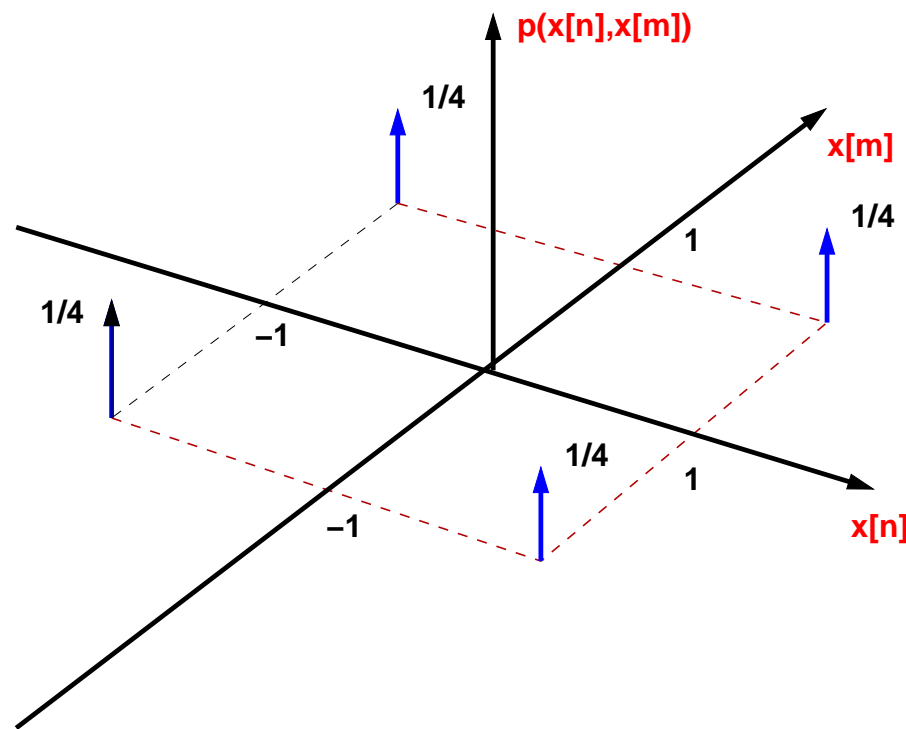
Luckily, real world time series often have “finite memory” (Markov)

Example 1.1. Bivariate pdf

Notice the change in indices (assuming **discrete time** signals)

$$CDF : \quad F(x[n], x[m]) = \text{Prob} \{X[n] \leq x[n], X[m] \leq x[m]\}$$

$$PDF : \quad p(x[n], x[m]) = \frac{\partial^2 F(x[n], x[m])}{\partial x[n] \partial x[m]}$$



Homework: Plot the CDF for this case. (What would happen in \mathbb{C} ?)

Properties of the statistical expectation operator

P1: Linearity

$$E\{ax[n] + by[m]\} = aE\{x[n]\} + bE\{y[m]\}$$

P2: $E\{x[m]y[n]\} \neq E\{x[m]\}E\{y[n]\}$ unless $\{x[m]\}$ and $\{y[n]\}$ are independent random processes

P3: If $y[n] = g(x[n])$ and the pdf of $x[n]$ is $p(x[n])$ then

$$E\{y[n]\} = \int_{-\infty}^{\infty} g(x[n])p(x[n])dx[n]$$

that is, we DO NOT need to know the pdf of $\{y[n]\}$ to find its expected values (when $g(\cdot)$ is a deterministic function).

NB: Think of a saturation-type sensor (microphone)

Example 1.2. Mean for linear systems (use P1 & P2)

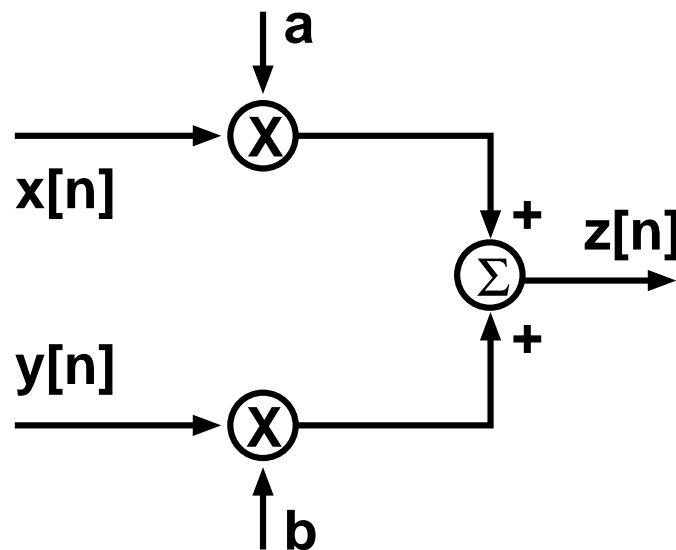
Consider a general linear system given by $z[n] = ax[n] + by[n]$. Find the mean ($E\{x[n]\} = \mu_x$, $E\{y[n]\} = \mu_y$, and $x \perp y$).

Solution:

$$E\{z[n]\} = E\{ax[n] + by[n]\} = aE\{x[n]\} + bE\{y[n]\}$$

that is

$$\mu_z = a\mu_x + b\mu_y$$



This is a consequence of the linearity of the $E\{\cdot\}$ operator.

Example 1.3. Mean for nonlinear systems (use P3)

For a nonlinear system, say the sensor nonlinearity is given by

$$z[n] = x^2[n]$$

using Property P3 of the statistical expectation operator, we have

$$\mu_z = E\{x^2[n]\} = \int_{-\infty}^{\infty} x^2[n]p(x[n])dx[n]$$

This is extremely useful, since most of the real-world signals are observed through sensors, e.g.

microphones, geophones, various probes ...

which are **almost invariably nonlinear** (typically a saturation type nonlinearity)

Dealing with ensembles of random processes

Ensemble \leadsto collection of **all possible realisations** of a **random signal**

The Ensemble Mean

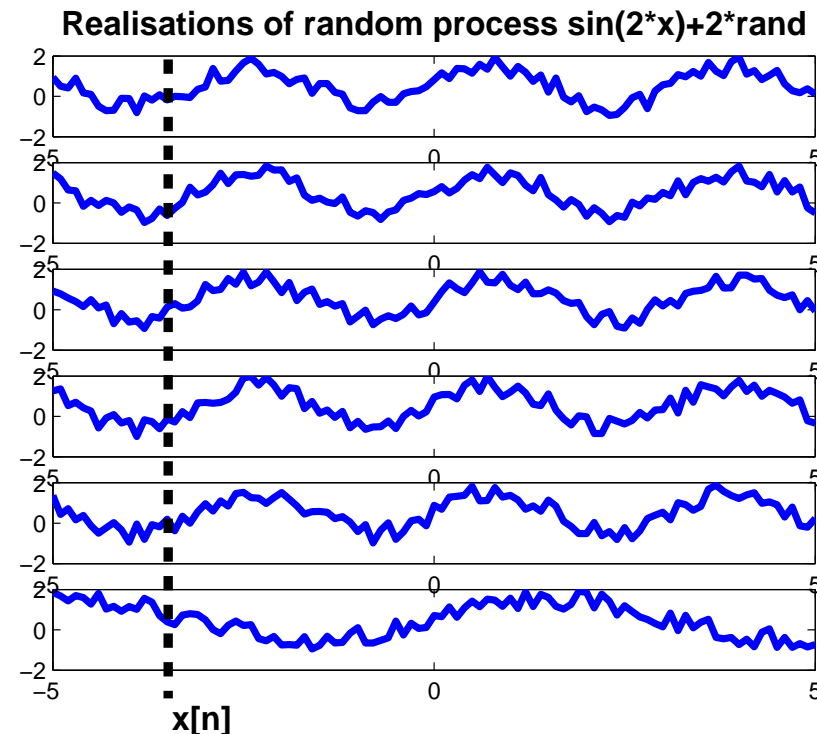
$$\mu_x(n) = \frac{1}{N} \sum_{i=1}^N x_i[n]$$

where $x_i[n]$ \leadsto outcome of i -th experiment at sample n .

For $N \rightarrow \infty$ we have

$$\mu_x(n) = \lim_{N \rightarrow \infty} \frac{1}{N} \sum_{i=1}^N x_i[n]$$

Average both **along** one and **across** all realisations?

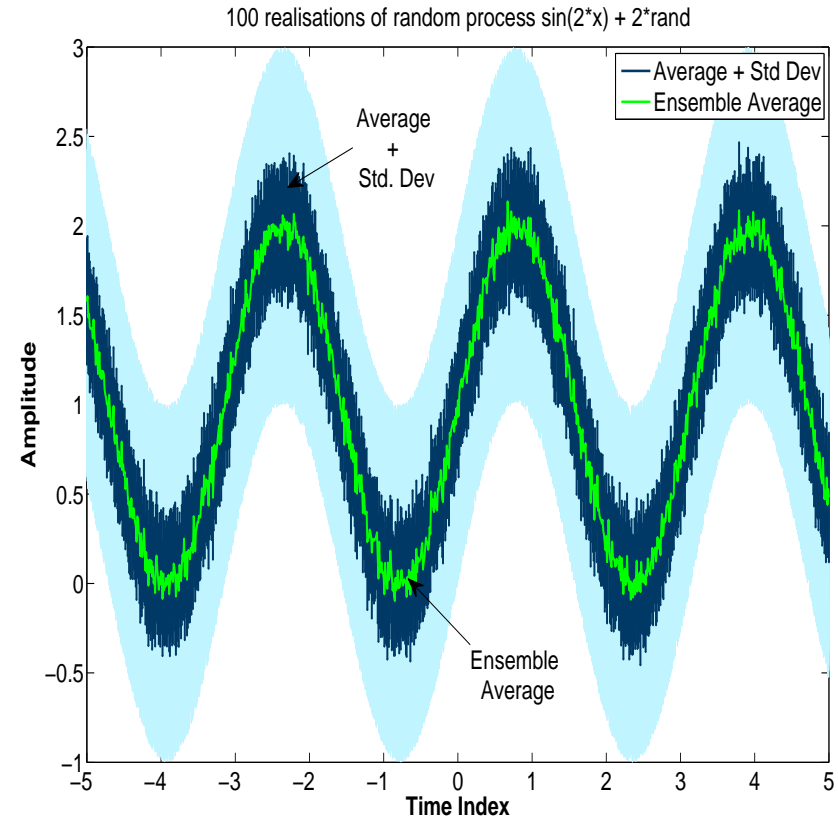
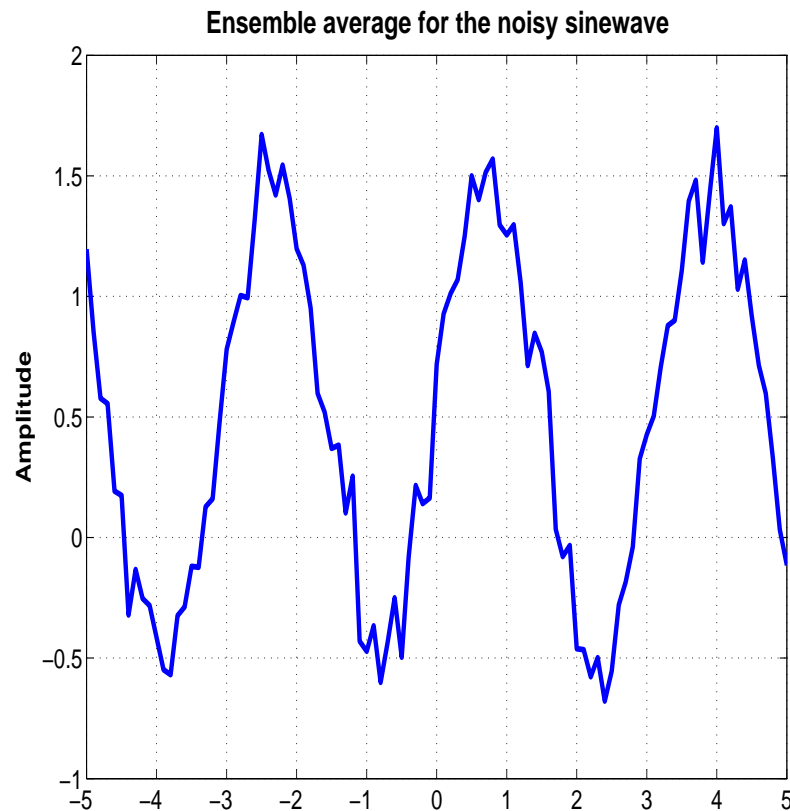


$$\text{Average Statistically} \quad E\{x[n]\} = \mu_x = \int_{-\infty}^{\infty} x[n] p(x[n]) dx[n]$$

Ensemble Average = Ensemble Mean

Our old noisy sine example

stochastic process is a collection of random variables



The pdf at time instant n is different from that at m , in particular:

$$\mu(n) \neq \mu(m) \quad m \neq n$$

Left & Right: Ensemble average

$$\sin(2x) + 2 * rand$$

Left: 6 realisations, **Right:** 100 realisations (and the overlay plot)

Second order statistics: 1) Correlation

- **Correlation (also known as Autocorrelation Function (ACF))**

$$r(m, n) = E\{x[m]x[n]\}, \quad \text{that is}$$

$$r(m, n) = \int_{-\infty}^{\infty} x[m]x[n]p(x[m], x[n])dx[m]dx[n]$$

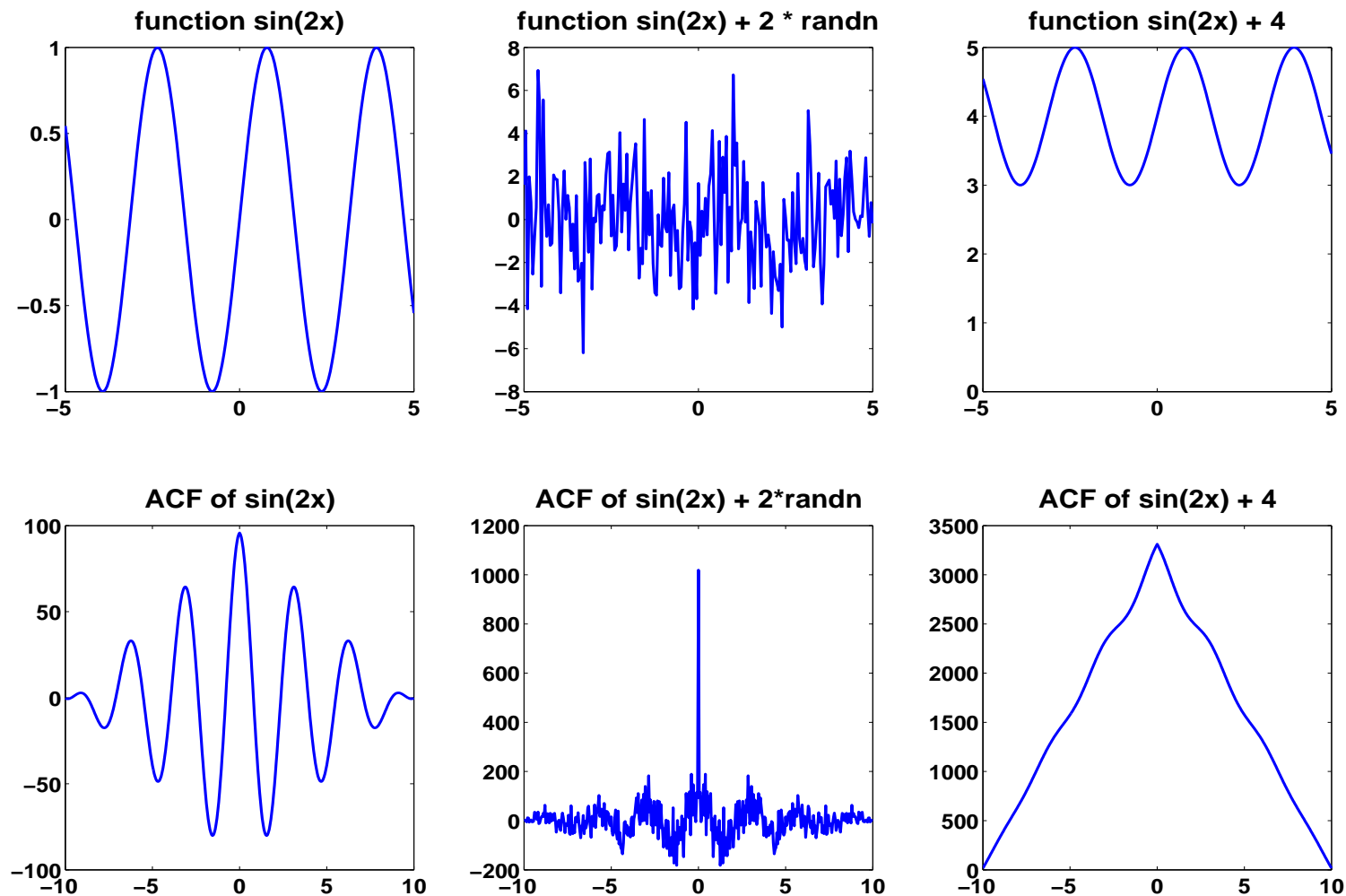
- in practice, for **ergodic signals** we calculate correlations from the **relative frequency perspective**

$$r(m, n) = \lim_{N \rightarrow \infty} \left\{ \frac{1}{N} \sum_{i=1}^N x_i[m]x_i[n] \right\}, \quad (i \text{ denotes the ensemble index})$$

- $r(m, n)$ measures the **degree of similarity** between $x[n]$ and $x[m]$.
- $r(n, n) = E\{x^2[n]\} \Rightarrow$ is the average "power" of a signal
- $r(m, n) = r(n, m) \Rightarrow$ the autocorrelation matrix of all $\{r(m, n)\}$

$$\mathbf{R} = \{r(m, n)\} = E[\mathbf{x}\mathbf{x}^H] \text{ is } \mathbf{symmetric}$$

Example 1.4. Autocorrelation of sinewaves (the need for a covariance function)



⇒ Useful information gets obscured in noise or under a DC offset

Second order statistics: 2) Covariance

- **Covariance** is defined as

$$\begin{aligned}c(m, n) &= E\{(x[m] - \mu(m))(x[n] - \mu(n))\} \\&= E\{x[m]x[n]\} - \mu(m)\mu(n) \\c(n, n) &= \sigma_n^2 = E\{(x[n] - \mu(n))^2\} \quad \text{for } m = n\end{aligned}$$

- Properties:

- $c(m, n) = c(n, m) \Rightarrow$ the covariance matrix for $\mathbf{x} = [x[0], \dots, x[N-1]]^T$ is **symmetric** and given by

$$\mathbf{C} = \{c(m, n)\} = E[\mathbf{x}\mathbf{x}^H], \text{ where } \mathbf{x} = \{x - \mu\}$$

- For zero mean signals, $c(m, n) = r(m, n)$

(see also the Standardisation slide and Example 1.4)

Higher order moments

For a zero-mean stochastic process $\{x[n]\}$:

- Third and fourth order moments

$$\text{Skewness : } R_3(l, m, n) = E\{x[l]x[m]x[n]\}$$

$$\text{Kurtosis : } R_4(l, m, n, p) = E\{x[l]x[m]x[n]x[p]\}$$

- In general, n -th order moment

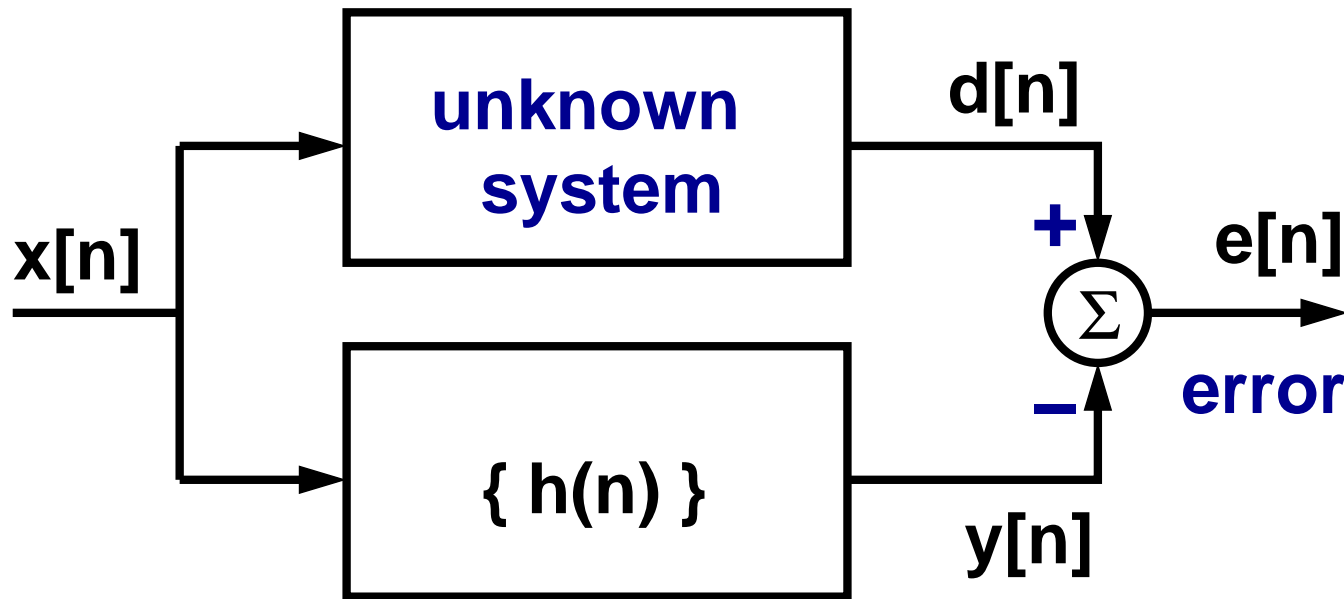
$$R_N(l_1, l_2, \dots, l_n) = E\{x[l_1]x[l_2] \cdots x[l_n]\}$$

Higher order moments can be used to form noise-insensitive statistics (cumulants).

- Important in non-linear signal processing
- Applications: blind source separation

⊗ In many applications the signals are assumed to be, or are reduced to, zero-mean stochastic process.

Example 1.5. Use of statistics in system identification (statistical rather than transfer function based analysis)



Task: Select $\{h(n)\}$ such that $y[n]$ is as similar to $d[n]$ as possible.

Measure of "goodness" is the distribution of the error $\{e[n]\}$.

Ideally, the error should be zero mean, white, and uncorrelated with the output signal

Solution: Minimise error power $E\{e^2[n]\}$ by selecting suitable $\{h(k)\}$

- Cost function: $J = E \left\{ \left(d[n] - \sum_k h(k)x[n-k] \right)^2 \right\}$
- Setting $\nabla_h J = 0$ for $h = h(i)$, gives (you will see more detail later)

$$E\{d[n]x[n-i]\} - \sum_k h(k)E\{x[n-k]x[n-i]\} = 0$$

- The solution $r_{dx}(-i) = \sum_k h(k)r_{xx}(i-k)$ in vector form is

$$\mathbf{h} = \mathbf{R}^{-1}\mathbf{r}_{dx}$$

\Rightarrow The optimal coefficients are **inversely proportional** to the autocorrelation matrix and **directly proportional** to the estimate of the crosscorrelation.

Independence, uncorrelatedness and orthogonality

- Two RV are **independent** if the realisation of one does not affect the distribution of the other, consequently, the joint density is separable:

$$p(x, y) = p(x)p(y)$$

Example: Sunspot numbers on 31 December and Argentinian debt

- Two RVs are **uncorrelated** if their cross-covariance is zero, that is

$$c(x, y) = E[(x - \mu_x)(y - \mu_y)] = E[xy] - E[x]E[y] = 0$$

Example: $x \sim \mathcal{N}(0, 1)$ and $y = x^2$ (impossible to relate through a linear relationship)

- Two RV are **orthogonal** if $r(x, y) = E[xy] = 0$

Example: Two uncorrelated RVs with at least one of them zero-mean

Independence, uncorrelatedness and orthogonality - Properties

- **Independent** RVs are always uncorrelated
- **Uncorrelatedness** can be seen as a 'weaker' form of independence since only the expectation (rather than the density) needs to be separable.
- **Uncorrelatedness** is a measure of **linear** independence. For instance, $x \sim \mathcal{N}(0, 1)$ and $y = x^2$ are clearly **dependent but uncorrelated**, meaning that there is no **linear** relationship between them
- Since $c_{xy} = r_{xy} - m_x m_y$ orthogonal RVs x and y need not be uncorrelated. Furthermore:
 - **uncorrelated** if they are independent and one of them is zero mean
 - **orthogonal** if they are uncorrelated and one of them is zero mean
- For uncorrelated random variables: $\text{var}\{x + y\} = \text{var}\{x\} + \text{var}\{y\}$

Stationarity: Strict and wide sense

- **Strict Sense Stationarity (SSS):** The process $\{x[n]\}$ is SSS if for all k the joint distribution $p(x[n_1], \dots, x[n_k])$ *is invariant under time shifts*, i.e. (all moments considered)

$$p(x[n_1 + n_0], \dots, x[n_k + n_0]) = p(x[n_1], \dots, x[n_k]), \forall n_0$$

As SSS is too strict for practical applications, we consider the more 'relaxed' stationarity condition: .

- **Wide-Sense Stationarity (WSS):** The process $\{x[n]\}$ is WSS if $\forall m, n$:
 - Mean: $E\{x[m]\} = E\{x[m + n]\},$
 - Covariance: $c(m, n) = c(m - n, 0) = c(m - n)$

Note that only the first two moments are considered.

Example of WSS: $x[n] = \sin(2\pi f n + \phi)$, where ϕ is uniformly distributed on $[-\pi, \pi]$

Autocorrelation function $r(m)$ of WSS processes

- i) **Time/shift invariant:** $r(m, n) = r(m - n, 0) = r(m - n)$ (follows from the covariance WSS requirement)
- ii) **Symmetric:** $r(-m) = r(m)$ (follows from the definition)
- iii) $r(0) \geq |r(m)|$ (maximum at $m = 0$)

The signal power = $r(0) \leftrightarrow$ Parseval's relationship

Follows from $E\{(x[n] - \lambda x[n + m])^2\} \geq 0$, i.e.

$$E\{x^2[n]\} - 2\lambda E\{x[n]x[n + m]\} + \lambda^2 E\{x^2[n + m]\} \geq 0 \quad \forall \lambda$$

$$r(0) - 2\lambda r(m) + \lambda^2 r(0) \geq 0 \quad \forall \lambda$$

which is quadratic in λ and required to be positive for all λ , i.e. the equation determinant: $\Delta = r^2(m) - r(0)r(0) \leq 0 \Rightarrow r(0) \geq |r(m)|$.

Properties of ACF – continued

iv) The AC matrix for a stationary $\mathbf{x} = [x[0], \dots, x[L-1]]^T$ is

$$\mathbf{R} = E\{\mathbf{x}\mathbf{x}^T\} = E\left\{ \begin{bmatrix} x_0 \\ x_1 \\ \vdots \\ x_{L-1} \end{bmatrix} [x_0, x_1, \dots, x_{L-1}] \right\} = \begin{bmatrix} r(0) & r(1) & \cdots & r(L-1) \\ r(1) & r(0) & \cdots & \vdots \\ \vdots & \vdots & \ddots & r(1) \\ r(L-1) & r(L-2) & \cdots & r(0) \end{bmatrix}$$

is **symmetric and Toeplitz**.

v) \mathbf{R} is **positive semi-definite**, that is

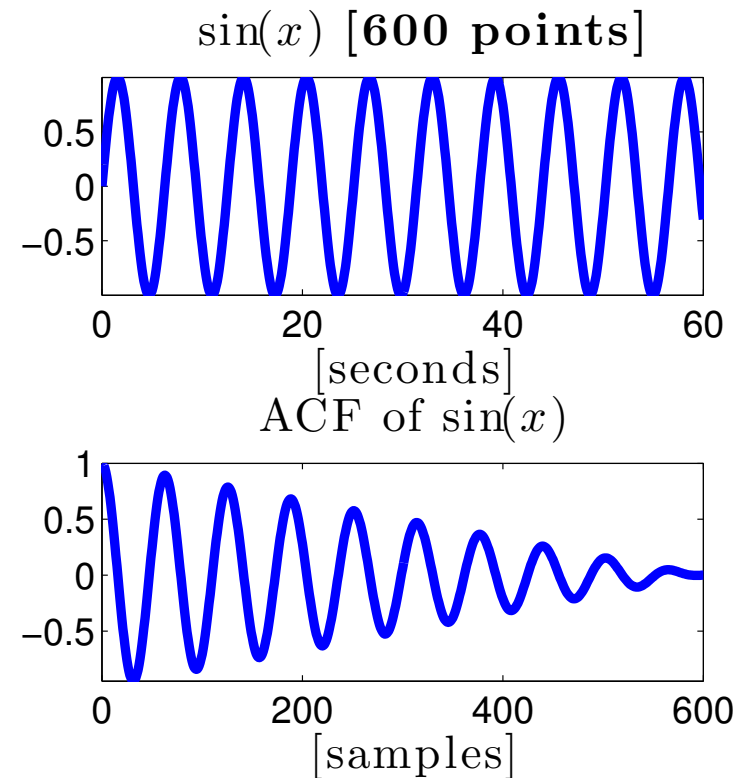
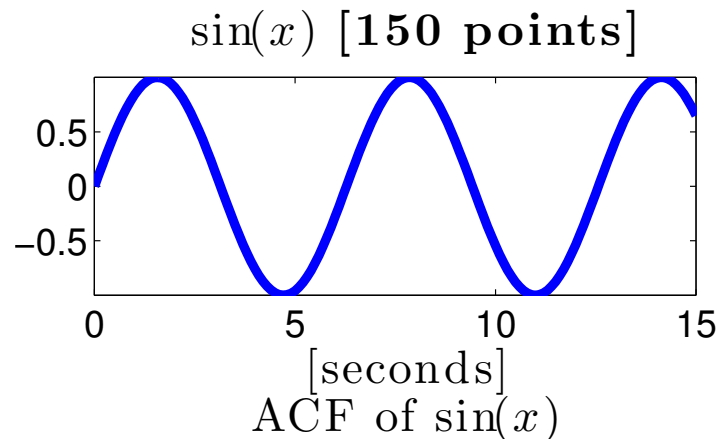
$$\mathbf{a}^T \mathbf{R} \mathbf{a} \geq 0 \quad \forall \mathbf{a} \neq 0$$

which follows from $y = \mathbf{a}^T \mathbf{x}$ and $y^T = \mathbf{x}^T \mathbf{a}$ so that

$$E\{y^2[n]\} = E\{y[n]y^T[n]\} = E\{\mathbf{a}^T \mathbf{x}\mathbf{x}^T \mathbf{a}\} = \mathbf{a}^T E\{\mathbf{x}\mathbf{x}^T\} \mathbf{a} = \mathbf{a}^T \mathbf{R} \mathbf{a} \geq 0$$

Properties of $r(m)$ – contd II

- vi) The autocorrelation function **reflects the basic shape of a signal**, for instance if the signal is periodic, the autocorrelation function will also be periodic



Sinewave and its ACF - Sampling rate=10Hz

Properties of the crosscorrelation

i) $r_{xy}(m) = E\{x[n]y[n+m]\} = r_{yx}(-m)$ (accounts for the lead/trail signal - see also the radar principle in Example 1.6)

ii) If $z[n] = x[n] + y[n]$ then

$$\begin{aligned} r_{zz}(m) &= E\{(x[n] + y[n])(x[n+m] + y[n+m])\} \\ &= r_{xx}(m) + r_{yy}(m) + r_{xy}(m) + r_{yx}(m) \end{aligned}$$

and if $x[n]$ and $y[n]$ are independent or uncorrelated

$$r_{zz}(m) = r_{xx}(m) + r_{yy}(m)$$

(therefore for $m = 0$ we have $\text{var}(z) = \text{var}(x) + \text{var}(y)$)

iii) $r_{xy}^2(m) \leq r_{xx}(0)r_{yy}(0)$ (Same as ACF P(iii) when $x = y$)

Example 1.6. The use of (cross-)correlation

Detection of Tones in Noise:

Consider a noisy tone $x = A \cos(\omega n + \theta)$

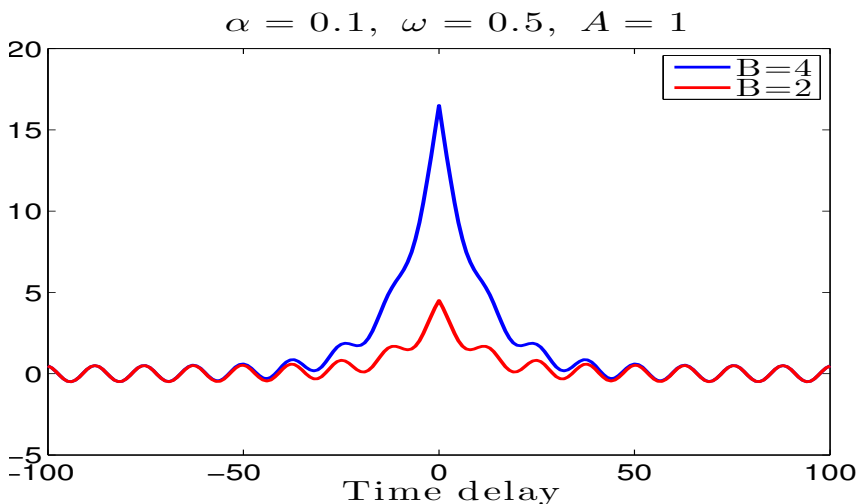
$$y[n] = A \cos(\omega n + \theta) + w[n]$$

$$\begin{aligned} \text{ACF : } R(m) &= E[y[n]y[n+m]] = \\ &= R_x(m) + R_w(m) + R_{xw}(m) + R_{wx}(m) \end{aligned}$$

For $R_w = B^2 \exp(-\alpha|m|)$ & $x \perp w$, then

$$R_y(m) = \frac{1}{2}A^2 \cos(\omega m) + B^2 \exp(-\alpha|m|)$$

- for large m , the ACF \propto the signal
- \exists extract tiny signal from large noise



Principle of Radar:

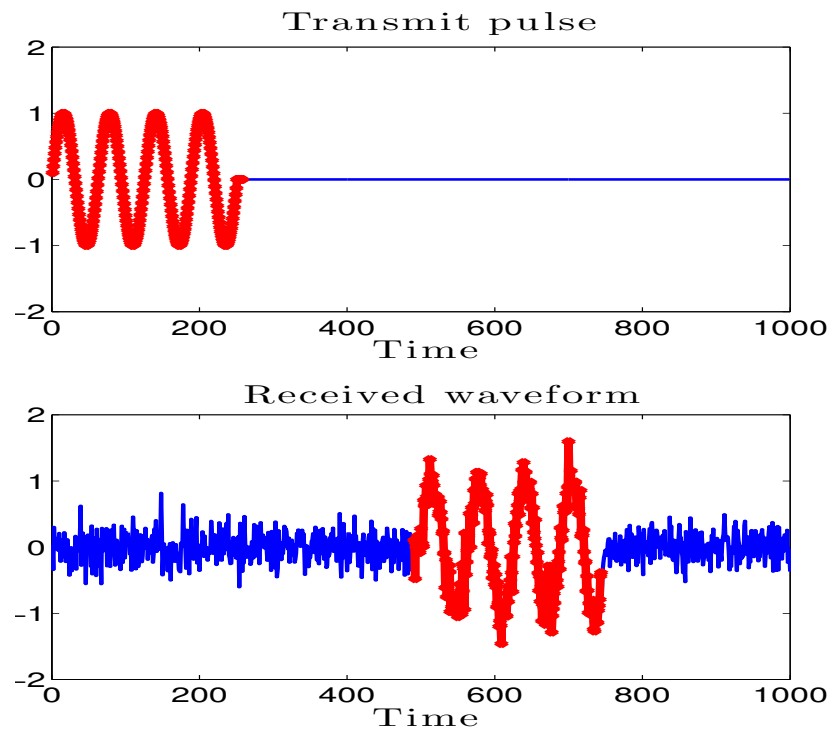
The received signal (see previous slide)

$$y[n] = ax[n - T_0] + w[n], \quad \text{so that}$$

$$\begin{aligned} R_{xy}(\tau) &= E\{x(n)y(n+\tau)\} \\ &= aR_x(\tau - T_0) + R_{xw}(\tau) \end{aligned}$$

Since

$$x \perp w \rightsquigarrow R_{xy}(\tau) = aR_x(\tau - T_0)$$



Example 1.7. Range of a radar

Unbiased estimate of a true radar delay δ_0 that has distribution $\delta \sim \mathcal{N}(\delta_0, \sigma_0^2)$

Q: What is the distribution of the range of the radar, and how should the radar be designed (i.e. what should σ_0 be) so that the range estimate is within 100 of the actual range with a probability of 99%?

A: The range is given by $R = \delta \frac{C}{2}$, therefore, $R \sim \mathcal{N}(\delta_0 \frac{C}{2}, \sigma_0^2 \frac{C^2}{4})$, where $R_0 = \delta_0 \frac{C}{2}$ is the actual true range.

To fulfil the radar design requirement, we need, $\mathbb{P}\{|R - R_0| < 100\} = 0.99$, or equivalently (due to the symmetry of the RV R)

$$\mathbb{P}\left\{\frac{(R - R_0)}{\sigma_0^2 C/2} < \frac{100}{\sigma_0^2 C/2}\right\} = 0.995,$$

and as $\frac{(R - R_0)}{\sigma_0^2 C/2} \sim \mathcal{N}(0, 1)$, we have $P\left(\frac{100}{\sigma_0^2 C/2}; 1, 0\right) = 0.995$. Evaluating this from the expression of the Gaussian CDF in an earlier slide we have

$$\frac{100}{\sigma_0^2 C/2} = 2.58 \Rightarrow \sigma_0 = \sqrt{\frac{200}{2.58 \times 3 \times 10^8}} = 0.51 \text{ milliseconds}$$

NB: By dividing $\mathcal{N}(0, \sigma)$ with σ we standardise pdf to unit variance $\mathcal{N}(0, 1)$.

Power spectral density (PSD)

The **power spectrum** or **power spectral density** $S(f)$ of a process $\{x[n]\}$ is the Fourier transform of its ACF (Wiener–Khinchine Theorem)

$$S(f) = \mathcal{F}\{r_{xx}(m)\} = \sum_{m=-\infty}^{\infty} r_{xx}(m)e^{-j2\pi mf} \quad f \in (-1/2, 1/2], \omega \in (-\pi, \pi]$$

The sampling period T is assumed to be unity, thus f is the *normalised frequency*.

From the inversion formula (Fourier), we can write

$$r_{xx}(m) = \int_{-1/2}^{1/2} S(f)e^{j2\pi mf} df$$

⇒ ACF tells us about the power within the signal (**Average**)

⇒ PSD tell us about the distribution of power across frequencies (**Density**)

PSD properties

i) $S(f)$ is a **positive real** function (**it is a distribution**) $\Rightarrow S(f) = S^*(f)$.

Since $r(-m) = r(m)$ we can write

$$S(f) = \sum_{m=-\infty}^{\infty} r_{xx}(-m)e^{j2\pi mf} = \sum_{m=-\infty}^{\infty} r_{xx}(m)e^{-j2\pi mf}$$

and hence

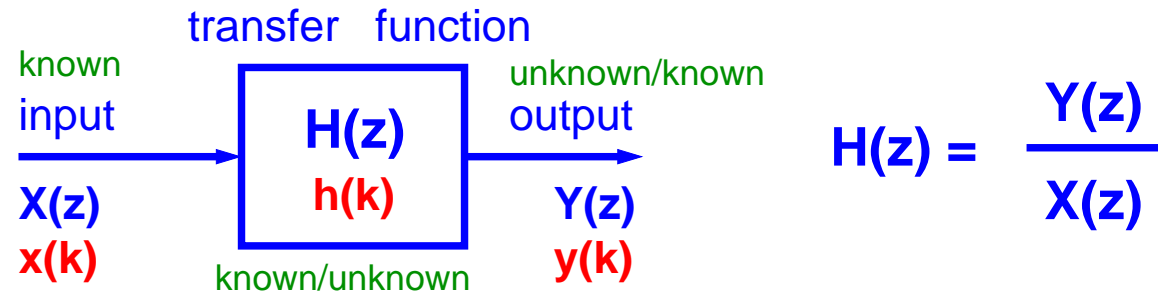
$$S(f) = \sum_{m=-\infty}^{\infty} r_{xx}(m) \cos(2\pi mf) = r_{xx}(0) + 2 \sum_{m=1}^{\infty} r_{xx}(m) \cos(2\pi mf)$$

ii) $S(f)$ is a **symmetric** function, $S(-f) = S(f)$. This follows from the last expression.

$$\text{iii) } r(0) = \int_{-1/2}^{1/2} S(f) df = E\{x^2[n]\} \geq 0.$$

\Rightarrow **the area below the PSD (power spectral density) curve = Signal Power.**

Linear systems



Described by their impulse response $h(n)$ or the transfer function $H(z)$

In the frequency domain (remember that $z = e^{j\theta}$) the transfer function is

$$H(\theta) = \sum_{n=-\infty}^{\infty} h(n)e^{-jn\theta} \quad \{x[n]\} \rightarrow \left| \begin{array}{c} \{h(n)\} \\ H(\theta) \end{array} \right| \rightarrow \{y[n]\}$$

that is

$$y[n] = \sum_{r=-\infty}^{\infty} h(r)x[n-r] = h * x$$

Example 1.8. Linear systems – statistical properties \leadsto mean and variance

i) Mean

$$E\{y[n]\} = E\left\{\sum_{r=-\infty}^{\infty} h(r)x[n-r]\right\} = \sum_{r=-\infty}^{\infty} h(r)E\{x[n-r]\}$$

$$\Rightarrow \mu_y = \mu_x \sum_{r=-\infty}^{\infty} h(r) = \mu_x H(0)$$

[NB: $H(\theta) = \sum_{r=-\infty}^{\infty} h(r)e^{-jr\theta}$. For $\theta = 0$, then $H(0) = \sum_{r=-\infty}^{\infty} h(r)$]

ii) Cross-correlation

$$r_{yx}(m) = E\{y[n]x[n+m]\} = \sum_{r=-\infty}^{\infty} h(r)E\{x[n-r]x[n+m]\}$$

$$= \sum_{r=-\infty}^{\infty} h(r)r_{xx}(m+r) \quad \text{convolution of input ACF and } \{h\}$$

$$\Rightarrow \text{Cross-power spectrum } S_{yx}(f) = \mathcal{F}(r_{yx}) = S_{xx}(f)H(f)$$

Example 1.9. Linear systems – statistical properties \leadsto crosscorrelation (this will be used in AR spectrum)

From $r_{xy}(m) = r_{yx}(-m)$ we have

$r_{xy}(m) = \sum_{r=-\infty}^{\infty} h(r)r_{xx}(m-r)$. Now we write

$$\begin{aligned} r_{yy}(m) &= E\{y[n]y[n+m]\} = \sum_{r=-\infty}^{\infty} h(r)E\{x[n-r]y[n+m]\} \\ &= \sum_{r=-\infty}^{\infty} h(r)r_{xy}(m+r) = \sum_{r=-\infty}^{\infty} h(-r)r_{xy}(m-r) \end{aligned}$$

by taking Fourier transforms we have

$$S_{xy}(f) = S_{xx}(f)H(f)$$

$$S_{yy}(f) = S_{xy}(f)H(-f) = \mathcal{F}(r_{xx})$$

or

$$\mathbf{S}_{yy}(\mathbf{f}) = \mathbf{H}(\mathbf{f})\mathbf{H}(-\mathbf{f})\mathbf{S}_{xx}(\mathbf{f}) = |\mathbf{H}(\mathbf{f})|^2\mathbf{S}_{xx}(\mathbf{f})$$

Output power spectrum = input power spectrum \times squared transfer function

Crosscorrelation and cross-PSD (recap)

- CC of two jointly WSS discrete time signals (**this is not symmetric**)

$$r_{xy}(m) = E\{x[n]y[n+m]\} = r_{yx}(-m)$$

- For $z[n] = x[n] + y[n]$ where $x[n]$ and $y[n]$ are zero mean and independent, we have $r_{xy}(m) = r_{yx}(m) = 0$, therefore

$$\begin{aligned} r_{zz}(m) &= r_{xx}(m) + r_{yy}(m) + r_{xy}(m) + r_{yx}(m) \\ &= r_{xx}(m) + r_{yy}(m) \end{aligned}$$

- Cross Power Spectral Density

$$P_{xy}(f) = \mathcal{F}\{r_{xy}(m)\}$$

Generally a complex quantity and so will contain both the **magnitude** and **phase** information.

Special signals: a) White noise

If the joint pdf is separable

$$p(x[0], x[1], \dots, x[n]) = p(x[0])p(x[1]) \cdots p(x[n]) \quad \forall n$$

where the pdf's $p(x[r])$ are identical $\forall r$, then all the pairs $x[n], x[m]$ are **independent** and $\{x[n]\}$ is said to be an **independent identically distributed (iid) signal**.

Since independent samples $x[n]$ are also uncorrelated, then for a zero-mean signal we have

$$r(n - m) = E\{x[m]x[n]\} = \sigma^2 \delta(n - m)$$

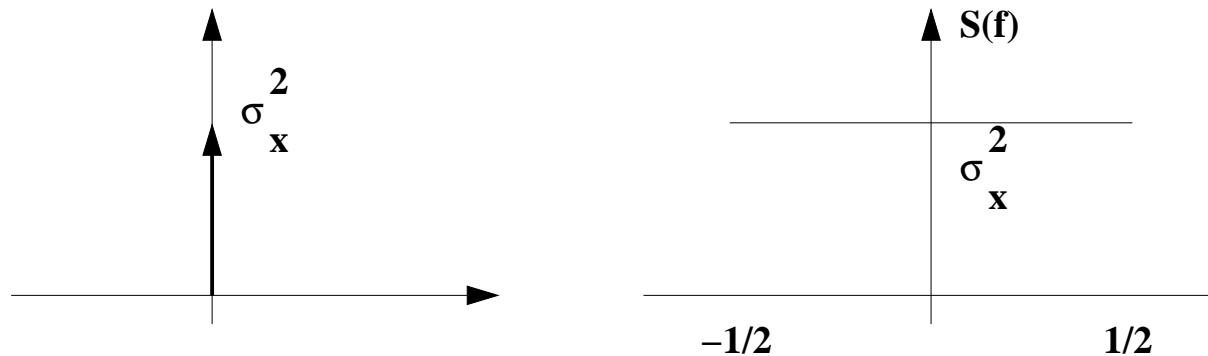
where the variance (signal power) $\sigma^2 = E\{x^2[n]\}$ and

$$\delta(n - m) = \begin{cases} 1 & n = m \\ 0 & \text{elsewhere} \end{cases}$$

where $\delta(n)$ is the Kronecker delta operator

Example 1.10. ACF and power spectrum of white noise

The Fourier transform of WN is constant for all frequencies, hence "white".



- The autocorrelation matrix
$$\mathbf{R} = \sigma_x^2 \mathbf{I} \quad r(m) = \sigma_x^2 \delta(m)$$

Since $E\{x[n]x[n-1]\} = 0$, the variance $r(0) = \sigma_x^2$ is the power of WN.

- The shape of the pdf $p(x[n])$ determines whether the white noise is called Gaussian (WGN), uniform (UWN), Poisson, Laplacian, etc.

From the Wiener–Khinchine Theorem:

$$\text{PSD(White Noise)} = \text{FT(ACF(WN))} = \text{FT}(\delta(t) \text{ function}) = \text{constant}$$

b) First order Markov signals: autoregressive modelling (finite memory in the description of a random signal)

If instead of the iid condition, we have the **first order conditional expectation**, then

$$p(x[n], x[n-1], x[n-2], \dots, x[0]) = p(x[n]|x[n-1])$$

where $p(a|b)$ is defined as the pdf of "a" *conditioned* upon the (possibly correlated) observation "b"

\Rightarrow the signal above is the **first order Markov signal**.

Example: Examine the statistical properties of the signal given by

$$y[n] = ay[n-1] + w[n]$$

where $a = 0.8$ and $w[n] \sim \mathcal{N}(0, 1)$ (see your coursework).

c) Minimum phase signals

Let $\{x[n]\}$ be observed for $n = 0, 1, \dots, N - 1$.

$$X(z) = x[0] + x[1]z^{-1} + \dots + x[N - 1]z^{-(N-1)} =$$
$$A \prod_{i=1}^N (1 - z_i z^{-1}), \quad A(0) = x[0]$$

- $|z_i| \leq 1$, $\forall i$ then $X(z)$ is said to be *minimum phase*
- $|z_i| \geq 1$, $\forall i$, then $X(z)$ is said to be *maximum phase*
- $|z_i| \geq 0$ for some i while for others $|z_i| \leq 1$ then $X(z)$ is said to be of *mixed phase*.

In DSP, the algorithms often rely on the minimum phase property of a signal for stability (of the inverse system) and to be able to have real-time implementation (causality).

d) Gaussian random signals

A signal in which each of the L samples is Gaussian distributed

$$p(x[i]) = \frac{1}{\sqrt{2\pi\sigma_i^2}} e^{-\frac{(x[i]-\mu(i))^2}{2\sigma_i^2}} \quad i = 0, \dots, L-1$$

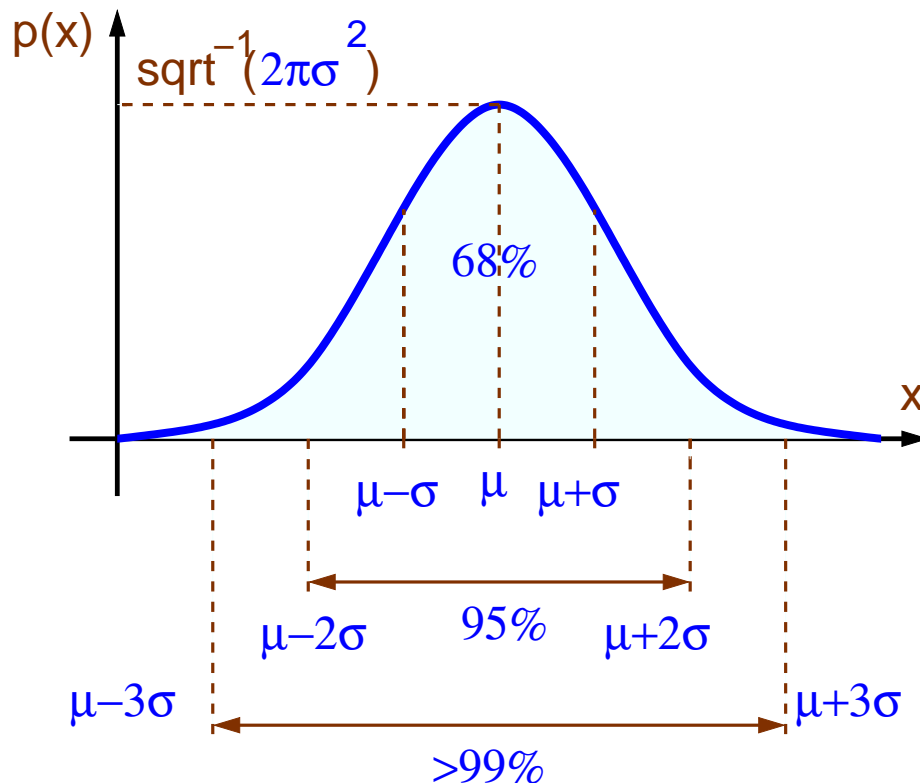
denoted by $\mathcal{N}(\mu(i), \sigma_i^2)$.

The joint pdf of L samples $x[n_0], x[n_1], \dots, x[n_{L-1}]$ is then

$$\begin{aligned} p(\mathbf{x}) &= p(x[n_0], x[n_1], \dots, x[n_{L-1}]) \\ p(\mathbf{x}) &= \frac{1}{[2\pi]^{L/2} \det(\mathbf{C})^{1/2}} e^{-\frac{(\mathbf{x}-\boldsymbol{\mu})^T \mathbf{C}^{-1} (\mathbf{x}-\boldsymbol{\mu})}{2}} = \frac{1}{(2\pi\sigma^2)^{L/2}} e^{-\frac{1}{2\sigma^2} \sum_{n=0}^{L-1} (x[n]-\mu)^2} \end{aligned}$$

where $\mathbf{x} = [x[n_0], x[n_1], \dots, x[n_{L-1}]]$, $\boldsymbol{\mu} = [\mu[n_0], \mu[n_1], \dots, \mu[n_{L-1}]]$ and \mathbf{C} is a covariance matrix with determinant Δ .

e) Properties of a Gaussian distribution



1) If x and y are jointly Gaussian, then for any constants a and b the random variable

$$z = ax + by$$

is Gaussian with mean

$$m_z = am_x + bm_y$$

and variance

$$\sigma_z^2 = a^2\sigma_x^2 + b^2\sigma_y^2 + 2ab\sigma_x\sigma_y\rho_{xy}$$

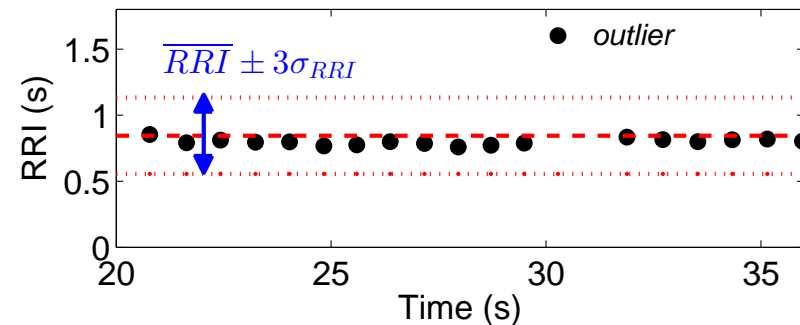
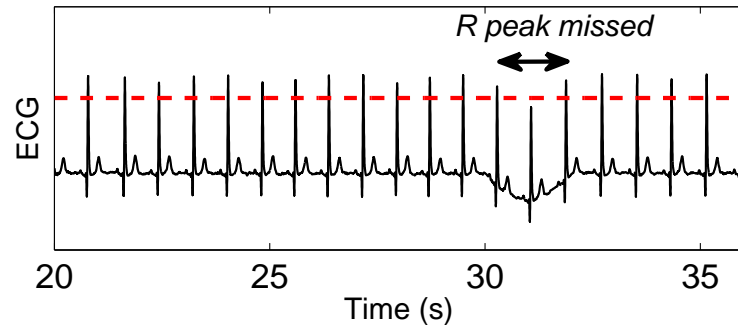
2) If two jointly Gaussian random variables are *uncorrelated* ($\rho_{xy} = 0$) then they are statistically independent,

$$f_{x,y} = f(x)f(y)$$

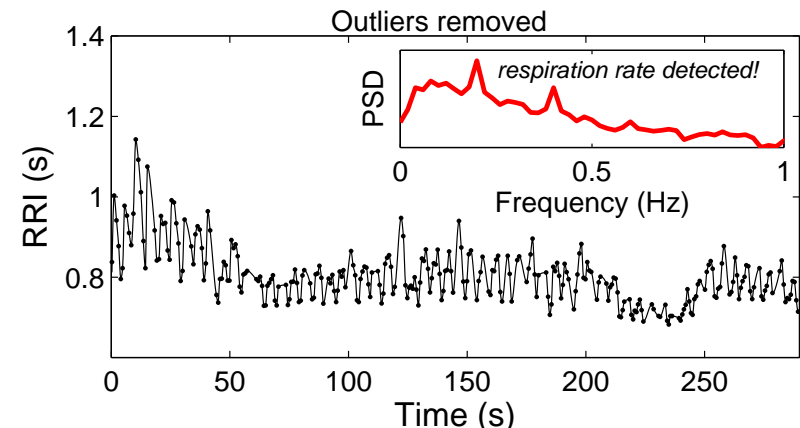
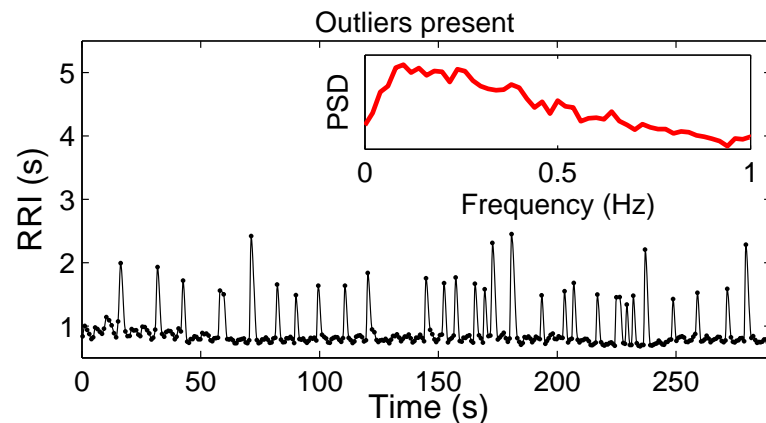
For $\mu = 0$, $\sigma = 1$, the inflection points are ± 1

Example 1.11. Rejecting outliers from cardiac data

- Failed detection of R peaks in ECG [left] causes outliers in R-R interval (RRI, time difference between consecutive R peaks) [right]



- No clear outcome from PSD analysis of outlier-compromised RRI [left], but PSD of RRI with outliers removed reveals respiration rate [right]



f) Conditional mean estimator for Gaussian random variables

3) If x and y are jointly Gaussian random variables then the optimum estimator for y , given by

$$\hat{y} = g(x)$$

that minimizes the mean square error $\xi = E\{[y - g(x)]^2\}$ is a **linear estimator** in the form

$$\hat{y} = ax + b$$

4) If x is Gaussian with zero mean then

$$E\{x^n\} = \begin{cases} 1 \times 3 \times 5 \times \cdots \times (n-1)\sigma_x^n, & n \text{ even} \\ 0, & n \text{ odd} \end{cases}$$

e) Ergodic signals

In practice, we often have only one observation of a signal (real-time)

Then, **statistical averages** are replaced by **time averages**.

This is necessary because

- ensemble averages are generally unknown a priori
- only a single realisation of the random signal is often available

Thus, the ensemble average

$$m_x(n) = \frac{1}{L} \sum_{i=1}^L x_i(n)$$

is therefore replaced by a time average

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

e) Ergodic signals – Example

Consider the random process

$$x(n) = A \cos(n\omega_0)$$

where A is a random variable that is equally likely to assume the value of 1 or 2.

The mean of this process is

$$E\{x(n)\} = E\{A\} \cos(n\omega_0) = 1.5 \cos(n\omega_0)$$

However, for a single realisation of this process, for large N , the sample mean is approximately zero

$$m_x \approx 0, \quad N \gg 1$$

$\Rightarrow x(n)$ is not ergodic and therefore the statistical expectation cannot be computed using time averages on a single realisation.

e) Ergodicity in the mean

DEF: If the sample mean $\hat{m}_x(N)$ of a WSS process converges to m_x , in the mean-square sense, then the process is said to be **ergodic in the mean**, and we write

$$\lim_{N \rightarrow \infty} \hat{m}_x(N) = m_x$$

For the convergence of the sample mean in the mean-square sense

- Asymptotically unbiased

$$\lim_{N \rightarrow \infty} E\{\hat{m}_x(N)\} = m_x$$

Consider the variance of the estimate $\rightarrow 0$ as $N \rightarrow \infty$

$$\lim_{N \rightarrow \infty} \text{Var}\{\hat{m}_x(N)\} = 0 \quad (\text{consistent})$$

e) Ergodicity - Summary

In practice, it is necessary to assume that the single realisation of a discrete time random signal satisfies ergodicity in the mean and autocorrelation.

Mean Ergodic Theorem: Let $x[n]$ be a wide sense stationary (WSS) random process with autocovariance sequence $c_x(k)$, sufficient conditions for $x[n]$ to be ergodic in the mean are that $c_x(k) < \infty$ and

$$\lim_{k \rightarrow \infty} c_x(k) = 0$$

Autocorrelation Ergodic Theorem:mean A necessary and sufficient condition for a WSS Gaussian process with covariance $c_x(k)$ to be autocorrelation ergodic is

$$\lim_{N \rightarrow \infty} \frac{1}{N} \sum_{k=0}^{N-1} c_x(k) = 0$$

Taylor series expansion

Most 'smooth' functions can be expanded into their Taylor Series Expansion (TSE)

$$f(x) = f(x_0) + \frac{f'(x_0)}{1}(x - x_0) + \frac{f''(x_0)}{2!}(x - x_0)^2 + \dots = \sum_{n=1}^{\infty} \frac{f^{(n)}(x_0)}{n!}$$

To show this consider the polynomial

$$f(x) = a_0 + a_1(x - x_0) + a_2(x - x_0)^2 + a_3(x - x_0)^3 + \dots$$

1. To get a_0 \leadsto choose $x = x_0 \Rightarrow a_0 = f(x_0)$
2. To get a_1 \leadsto take derivative of the polynomial above to have

$$\frac{d}{dx}f(x) = a_1 + 2a_2(x - x_0) + 3a_3(x - x_0)^2 + 4a_4(x - x_0)^3 + \dots$$

$$\text{choose } x = x_0 \Rightarrow a_1 = \left. \frac{df(x)}{dx} \right|_{x=x_0} \quad \text{and so on ... } a_k = \frac{1}{k!} \left. \frac{d^k f(x)}{dx^k} \right|_{x=x_0}$$

Power series - contd.

Consider

$$f(x) = \sum_{n=0}^{\infty} a_n x^n \Rightarrow f'(x) = \sum_{n=1}^{\infty} n a_n x^{n-1} \text{ and } \int_0^x f(t) dt = \sum_{n=0}^{\infty} \frac{a_n}{n+1} x^{n+1}$$

1. Exponential function, cosh, sinh, sin, cos, ...

$$e^x = \sum_{n=0}^{\infty} \frac{x^n}{n!} \text{ and } e^{-x} = \sum_{n=0}^{\infty} (-1)^n \frac{x^n}{n!} \Rightarrow \frac{e^x - e^{-x}}{2} = \sum_{n=0}^{\infty} \frac{x^{2n+1}}{(2n+1)!}$$

2. other useful formulas

$$\sum_{n=0}^{\infty} x^n = \frac{1}{1-x} \Rightarrow \sum_{n=1}^{\infty} n x^{n-1} = \frac{1}{(1-x)^2} \text{ and } \frac{1}{1+x^2} = \sum_{n=0}^{\infty} (-1)^n x^{2n}$$

Integrate to obtain $\arctan(x) = \sum_{n=0}^{\infty} (-1)^n \frac{x^{2n+1}}{2n+1}$.

For $x = 1$ we have $\frac{\pi}{4} = 1 = 1/3 + 1/5 - 1/7 + \dots$

Numerical derivatives - examples

- Two-point approximation:

- Forward: $f'(0) = \frac{f(1)-f(0)}{h}$
- Backward: $f'(0) = \frac{f(-1)-f(0)}{h}$

- Three-point approximation:

- $f'(0) = \frac{f(1)-2f(0)+f(-1)}{2h}$
- $f''(0) = \frac{f(1)-2f(0)+f(-1)}{h^2}$

- Five-point approximation (also look up for stencil):

- $f'(0) = \frac{f(-2)-8f(-1)+8f(1)-f(2)}{12h}$
- $f''(0) = \frac{-f(-2)+16f(-1)-30f(0)+16f(1)-f(2)}{12h^2}$

Constrained optimisation using Lagrange multipliers:

Basic principles

Consider a two-dimensional problem:

$$\begin{array}{ll}\text{maximize} & \underbrace{f(x, y)}_{\text{function to max/min}} \\ \text{subject to} & \underbrace{g(x, y) = c}_{\text{constraint}}\end{array}$$

↪ **we look for point(s) where curves f & g touch (but do not cross).**

In those points, the tangent lines for f and g are parallel \Rightarrow so too are the gradients $\nabla_{x,y}f \parallel \lambda \nabla_{x,y}g$, where λ is a scaling constant.

Although the two gradient vectors are parallel they can have different magnitudes

Therefore, we are looking for max or min points (x, y) of $f(x, y)$ for which

$$\nabla_{x,y}f(x, y) = -\lambda \nabla_{x,y}g(x, y) \quad \text{where } \nabla_{x,y}f = \left(\frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}\right) \text{ and } \nabla_{x,y}g = \left(\frac{\partial g}{\partial x}, \frac{\partial g}{\partial y}\right)$$

We can now combine these conditions into one equation as:

$$F(x, y, \lambda) = f(x, y) - \lambda(g(x, y) - c) \quad \text{and solve} \quad \nabla_{x,y,\lambda}F(x, y, \lambda) = \mathbf{0}$$

$$\textbf{Obviously, } \nabla_{\lambda}F(x, y, \lambda) = 0 \quad \Leftrightarrow \quad g(x, y) = c$$

The method of Lagrange multipliers in a nutshell

max/min of a function $f(x, y, z)$ where x, y, z are coupled

Since x, y, z **are not independent** there exists a constraint $g(x, y, z) = c$

Solution: Form a new function

$F(x, y, z, \lambda) = f(x, y, z) - \lambda(g(x, y, z) - c)$ and calculate $F'_x, F'_y, F'_z, F'_\lambda$

Set $F'_x, F'_y, F'_z, F'_\lambda = 0$ and solve for the unknown x, y, z, λ .

Example 1: Economics

Two

factories, A and B make TVs, at a cost $f(x, y) = 6x^2 + 12y^2$ ($x, y = \#TV \in A, B$). Minimise the cost of producing 90 TVs, by finding optimal numbers $\#x$ and $\#y$ at factories A and B.

Solution: The constraint $g: (x+y=90)$, so

$$F(x, y, \lambda) = 6x^2 + 12y^2 - \lambda(x + y - 90)$$

Then: $F'_x = 12x - \lambda$, $F'_y = 24y - \lambda$, $F'_\lambda = -x - y - 90$, and for min / max $\nabla F = 0$

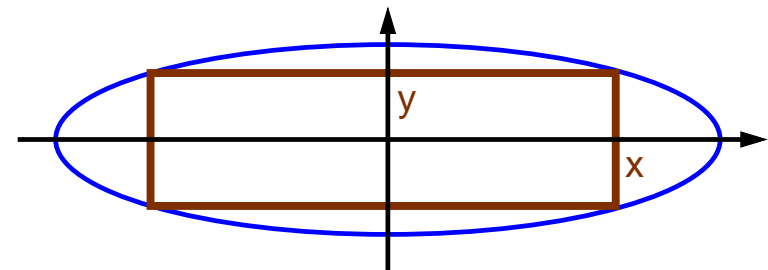
Set to zero $\Rightarrow x = 60, y = 30, \lambda = 720$

Example 2: Geometry

Find the

rectangle of maximal perimeter, inscribed in the ellipse $x^2 + 4y^2 = 4$.

Solution: Constraint ($x^2 + 4y^2 = 4$)



The perimeter $P(x, y) = 4x + 4y$ so

$$F(x, y, \lambda) = 4x + 4y - \lambda(x^2 + 4y^2 - 4)$$

$$\Rightarrow P'_x = \lambda g'_x, P'_y = \lambda g'_y \Rightarrow x = 4y$$

Solve to give: $x = 4/\sqrt{5}, P = 4\sqrt{5}$.

Notes:

○

Notes:

○