Introduction to statistical inference 2

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Chapter 1

Recap from "Introduction to statistical inference 1"

1.1 Random variable

• Random variable is a function from sample space S of an experiment to sample space of the random variable X, which is set of real numbers.

$$X: \mathcal{S} \to \mathcal{X}$$

- The sample space is a set of all possible values random variable can get.
- \mathcal{X} can be
 - an interval of real axis (continuous random variable).

$$\mathcal{X} = [0, 10), \ \mathcal{X} = [0, 10], \ \mathcal{X} = (0, 10)$$

- An uncountable set of integers (discrete random variable)

$$\mathcal{X} = \{0, 1, 2, \ldots\}$$

- A countable set of integers or real numbers (discrete random variable)

$$\mathcal{X} = \{0, 1\}, \ \mathcal{X} = \{0, 0.5, 1\}, \ \mathcal{X} = \{0, 1, \dots, 10\}$$

• Probabilities associated with each value of X are defined by the cumulative distribution function (cdf for short).

$$F_X(x) = P(X \le x)$$
, where $-\infty < x < \infty$

Note: $F_X(x)$ is a step function if X is discrete.

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Note: $F_X(x)$ is a continuous function if X is continuous.

- $F_X(x)$ or F(x) is cdf, if
 - 1) $\lim_{x\to-\infty} F(x) = 0$ and $\lim_{x\to\infty} F(x) = 1$
 - 2) F(x) is non-decreasing
 - 3) F(x) is right-continuous
- Cdf is useful in calculation of any probabilities; for example

$$P(a < X \le b) = F(b) - F(a)$$

Note: Be careful with < and \le when working with discreate random variables.

• The probability density function (pdf for short) is defined for continuous random variable as

$$f_X(x) = F'(x) = \frac{\mathrm{d}F_X(x)}{\mathrm{d}x}, -\infty < x < \infty$$

and

$$\int_{-\infty}^{x} f_X(t) \, \mathrm{d}t = F_X(x)$$

 The probability mass function (pmf for short) is defined for discrete random variables as

$$f_X(x) = P(X = x)$$

$$F_X(x) = \sum_{k=1}^{x} f_X(k)$$

1.2 Transformations of random variable

- Consider a monotonic function $g: \mathcal{X} \to \mathcal{Y}$
- Y = g(X) is also a random variable; function g is called an transformation (muunnos).
- If g(x) is a increasing function of x, then

$$F_Y(y) = F_X(g^{-1}(y))$$

• If g(x) is a decreasing function of x, then

$$F_Y(y) = 1 - F_X(g^{-1}(y))$$

 \bullet The pdf of continuous Y is

$$f_Y(y) = F_Y'(y)$$

1.3 Expected values

$$E(X) = \mu_X = \begin{cases} \int_{-\infty}^{\infty} x f(x) \, \mathrm{d}x & \text{if } X \text{ is continuous} \\ \sum_{x \in \mathcal{X}} x f(x) & \text{if } X \text{ is discrete} \end{cases}$$

$$E(g(X)) = \begin{cases} \int_{-\infty}^{\infty} g(x) f(x) \, \mathrm{d}x & \text{if } X \text{ is continuous} \\ \sum_{x \in \mathcal{X}} g(x) f(x) & \text{if } X \text{ is discrete} \end{cases}$$

1.4 Variance

$$\begin{split} \sigma_X^2 &= Var(X) &= E(X - \mu_X)^2 \\ &= E(X^2 - 2X\mu_X + \mu_X^2) \\ &= E(X^2) - E(2X\mu_X) + E(\mu_X^2) \\ &= E(X^2) - 2\mu_X \underbrace{E(X)}_{\mu_X} + E(\mu_X^2) \\ &= E(X^2) - \mu_X^2 \\ &= E(X^2) - \mu_X^2 \\ sd(X) &= \sqrt{Var(X)} = \sigma_X \end{split}$$

1.5 Bivariate random variables

• For two discrete random variables, the joint pmf is defined as

$$f_{X,Y}(x,y) = P(X = x, Y = y)$$

• For two continuous random variables, we define the joint pdf $f_{X,Y}(x,y)$ as

$$P((X,Y) \in A) = \iint_A f(x,y) dx dy$$

• The expected value for transformation $g(X,Y):\mathbb{R}^2\to\mathbb{R}$ (for example, g(X,Y)=XY or $g(X,Y)=\frac{X}{Y}$) is

$$E(g(X,Y)) = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} g(x,y) f(x,y) \, \mathrm{d}x \, \mathrm{d}y$$

if (X, Y) is continuous, and

$$E(g(X,Y)) = \sum_{x,y \in \mathbb{R}^2} g(x,y) f(x,y)$$

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if (X, Y) is discrete.

• The marginal pmf / pdf for X are

$$f_X(x) = \sum_{y \in \mathbb{R}} f_{X,Y}(x,y)$$
 (pmf)

$$f_X(x) = \int_{-\infty}^{\infty} f_{X,Y}(x,y) \,dy$$
 (pdf)

and correspondingly for Y.

• The conditional pmf / pdf are both defined as

$$f(y \mid x) = \frac{f_{X,Y}(x,y)}{f_X(x)}$$
 (for both discrete and continuous random variables)

and correspondingly for $x \mid y$.

1.6 Independence

• Random variables are said to be independent $(X \perp\!\!\!\perp Y)$ if

$$f_{X,Y}(x,y) = f_X(x)f_Y(y)$$

ullet For independent random variables, conditional distribution $y\mid x$ is

$$f(y \mid x) = f_y(y)$$

regardless of the value of x.

1.7 Covariance

 \bullet Covariance measures the linear association between two random variables X and Y

$$\begin{aligned} cov(X,Y) &= E((X-\mu_X)(Y-\mu_Y)) \\ &= E(XY) - \mu_X \mu_Y \\ cov(X,Y) &= cov(Y,X) \\ corr(X,Y) &= \rho_{XY} = \frac{cov(X,Y)}{\sigma_X \sigma_Y} & \text{where } -1 \leq \rho_{XY} \leq 1 \end{aligned}$$

Note: cov(X, X) = Var(X)

Note: If $X \perp\!\!\!\perp Y$, then cov(X,Y) = 0, but if cov(X,Y) = 0, it does not mean X and Y are necessarily independent.

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1.8 Random vectors

• Random vectors generalize the bivariate random variables to a *n*-variate case.

$$m{X} = egin{bmatrix} X_1 \\ X_2 \\ \vdots \\ X_n \end{bmatrix}$$
 where $X_i, \ i=1,2,\ldots,n$ are scalar random variables

• If x is a continuous random vector, then

$$P(\boldsymbol{X} \in A) = \int \cdots \int_{A} f(\boldsymbol{x}) dx_1 \dots dx_n \quad f : \mathbb{R}^2 \to \mathbb{R}$$

where f(x) is a joint pdf.

• If x is a discrete random vector, then

$$P(X \in A) = \sum ... \sum f(x)$$

where f(x) is a joint pmf.

• Let g(x) be a transformation $g: \mathbb{R}^2 \to \mathbb{R}$. Then, the expected value is

$$E(g(\boldsymbol{X})) = \begin{cases} \int_{-\infty}^{\infty} \dots \int_{-\infty}^{\infty} x f(x) \, \mathrm{d}x_1 \dots \, \mathrm{d}x_n & \text{if } \boldsymbol{X} \text{ is continuous} \\ \sum \dots \sum_{\boldsymbol{X} \in \mathbb{R}^2} g(\boldsymbol{x}) f(\boldsymbol{x}) & \text{if } \boldsymbol{X} \text{ is discrete} \end{cases}$$

• Let us partition the n-variate random vector X as follows:

$$X = \begin{pmatrix} X_1 \\ X_2 \end{pmatrix}$$

where X_1 has length k and X_2 has length n-k.

• The joint pdf of X can be written as

$$f(\boldsymbol{x}) = f(\boldsymbol{x}_1, \boldsymbol{x}_2)$$

• The marginal density of X_1 is

$$f(m{x}_1) = \int\limits_{\mathbb{R}^{n-k}} f(m{x}_1, m{x}_2) \, \mathrm{d}m{x}_2 \quad ext{if } m{X} ext{ is continuous}$$
 $f(m{x}_1) = \sum\limits_{\mathbb{R}^{n-k}} f(m{x}_1, m{x}_2) \qquad \quad ext{if } m{X} ext{ is discrete}$

where $f(x_1)$ is a k-variate pdf/pmf.

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• The conditional pdf/pmf for $X_2 \mid X_1$ is

$$f(\boldsymbol{X}_2 \mid \boldsymbol{X}_1) = \frac{f(\boldsymbol{X}_1, \boldsymbol{X}_2)}{f(\boldsymbol{X}_2)}$$

• Expected value of random vectors is defined as vector

$$E(oldsymbol{x}) = egin{bmatrix} E(X_1) \ E(X_2) \ dots \ E(X_n) \end{bmatrix} = egin{bmatrix} E(X_1) \ E(X_2) \end{bmatrix} = egin{bmatrix} oldsymbol{\mu}_1 \ oldsymbol{\mu}_2 \end{bmatrix}$$

• The variance of a random vector is $n \times n$ symmetric matrix called variance-covariance matrix.

$$Var(\boldsymbol{x})_{n \times n} = \begin{bmatrix} Var(X_1) & cov(X_1, X_2) & cov(X_1, X_3) & \dots & cov(X_1, X_n) \\ cov(X_2, X_1) & Var(X_2) & cov(X_2, X_3) & \dots & cov(X_2, X_n) \\ cov(X_3, X_1) & cov(X_3, X_2) & Var(X_3) & \dots & cov(X_3, X_n) \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ cov(X_n, X_1) & cov(X_n, X_2) & cov(X_n, X_3) & \dots & Var(X_n) \end{bmatrix}$$

$$= \begin{bmatrix} Var(\boldsymbol{x}_1)_{k \times k} & cov(\boldsymbol{x}_1, \boldsymbol{x}_2')_{k \times (n-k)} \\ cov(\boldsymbol{x}_2, \boldsymbol{x}_1')_{(n-k) \times k} & Var(\boldsymbol{x}_2)_{(n-k) \times (n-k)} \end{bmatrix}$$

• The correlation matrix is defined as

$$corr(\boldsymbol{x})_{n \times n} = \begin{bmatrix} 1 & corr(X_1, X_2) & corr(X_1, X_3) & \dots & corr(X_1, X_n) \\ corr(X_2, X_1) & 1 & corr(X_2, X_3) & \dots & corr(X_2, X_n) \\ corr(X_3, X_1) & corr(X_3, X_2) & 1 & \dots & corr(X_3, X_n) \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ corr(X_n, X_1) & corr(X_n, X_2) & corr(X_n, X_3) & \dots & 1 \end{bmatrix}$$

 \bullet If X has a n-variate normal distribution, then, with

$$E(oldsymbol{x}) = oldsymbol{\mu} = egin{bmatrix} oldsymbol{\mu}_1 \ oldsymbol{\mu}_2 \end{bmatrix} \quad ext{and} \quad Var(oldsymbol{x})_{n imes n} = oldsymbol{\Sigma} = egin{bmatrix} oldsymbol{\Sigma}_1 & oldsymbol{\Sigma}_{12} \ oldsymbol{\Sigma}_{21} & oldsymbol{\Sigma}_2 \end{bmatrix}$$

then $X_1 \mid X_2$ has a k-variate normal distribution with

$$E(\boldsymbol{X}_1 \mid \boldsymbol{X}_2) = \boldsymbol{\mu}_1 + \boldsymbol{\Sigma}_{12} \boldsymbol{\Sigma}_2^{-1} (\boldsymbol{X}_2 - \boldsymbol{\mu}_2)$$
$$Var(\boldsymbol{X}_1 \mid \boldsymbol{X}_2) = \boldsymbol{\Sigma}_1 - \boldsymbol{\Sigma}_{12} \boldsymbol{\Sigma}_2^{-1} \boldsymbol{\Sigma}_{21}$$

Note: $Var(\boldsymbol{X}_1 \mid \boldsymbol{X}_2) \leq Var(\boldsymbol{X}_1)$

1.9 Computing using expected values and variances

Let a, b and c be constants, and let X, Y and Z be (scalar) random variables. The following rules hold regardless of the distribution of random variables X, Y and Z.

$$E(c) = c$$

$$E(cX) = cE(X)$$

$$E(X+Y) = E(X) + E(Y)$$

$$E(X+c) = E(X) + c$$

$$E(XY) = E(X)E(Y)$$
 Only if $X \perp \!\!\! \perp Y$.
$$E(g(X)) = g(E(X))$$
 Only in some special cases, like when $g(X)$ is a linear transformation.

$$\begin{split} Var(X+Y) &= Var(X) + Var(Y) + 2cov(X,Y) \\ Var(X-Y) &= Var(X) + Var(Y) - 2cov(X,Y) \\ Var(aX) &= a^2 \cdot Var(X) \\ cov(X,Y+Z) &= cov(X,Y) + cov(X,Z) \\ cov(aX,bY) &= ab \cdot cov(XY) \\ Var(X+a) &= Var(X) \\ cov(X+a,Y+b) &= cov(X,Y) \\ E(X) &= E_Y \left[E_{X|Y}(X\mid Y) \right] \\ Var(X) &= E_Y \left[Var_{X|Y}(X\mid Y) \right] + Var_Y \left[Var_{X|Y}(X\mid Y) \right] \end{split}$$

Let a and b be fixed vectors and X and Y random vectors so that the dimensions in the equations match.

$$\begin{split} E(\boldsymbol{a}'\boldsymbol{X}) &= \boldsymbol{a}'E(\boldsymbol{X}) \\ Var(\boldsymbol{a}'\boldsymbol{X}) &= \boldsymbol{a}'Var(\boldsymbol{X})\boldsymbol{a} \\ cov(\boldsymbol{a}'\boldsymbol{X},\boldsymbol{b}'\boldsymbol{Y}) &= \boldsymbol{a}'cov(\boldsymbol{X},\boldsymbol{Y})\boldsymbol{b} \end{split}$$
 Compare to $Var(aX) = a^2 \cdot Var(X) = a \cdot Var(X) \cdot a$

Note: These equations need to be remembered by hearth!

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Chapter 2

Random samples

Definition 2.1. Random variables $X_1, ..., X_n$ are called random sample of size n from population f(x), if $X_1, ..., X_n$ are mutually independent random variables and the marginal pdf/pmf if each X_i is the same function f(x). Alternatively, $X_1, ..., X_n$ are called independent and identically distributed random variables (i.i.d.) with pdf/pmf f(x).

Note: Sample X_1, \ldots, X_n can also be denoted by X, where $X = \begin{bmatrix} X_1 & \ldots & X_n \end{bmatrix}^T$

• If follows from the mutual independence, of X_1, \ldots, X_n that the joint pdf or pmf of X is

$$f(x_1, \dots, x_2) = f(x_1)f(x_2)\dots f(x_n) = \prod_{i=1}^n f(x_i)$$

• All univariate marginal distributions $f(x_i)$ are the same by definition 2.1.

Example 2.1. Let X_1, \ldots, X_n be a random sample from $Exponential(\beta)$ population. X_i specifies the time until failure for n identical cellphones. The exponential pdf is

$$f(x_i) = \frac{1}{\beta} e^{-x_i/\beta}$$

The joint pdf pf the sample is

$$f(x_1,x_2,\ldots,x_n\mid\beta)=\prod_{i=1}^n\frac{1}{\beta}e^{-x_i/\beta}=\frac{1}{\beta^n}e^{\frac{1}{\beta}\sum x_i}$$
 Recall: $a^ba^c=a^{b+c}$

What is the probability that all n cellphones last more than 2 years?

$$P(X_1 > 2, \dots, X_n > 2) = \int_2^{\infty} \dots \int_2^{\infty} f(x_1, \dots, x_n) \, \mathrm{d}x_1 \dots \, \mathrm{d}x_n$$

$$= \int_2^{\infty} \dots \int_2^{\infty} \prod_{i=1}^{1} \frac{1}{\beta} e^{-x_i/\beta} \, \mathrm{d}x_1 \dots \, \mathrm{d}x_n$$

$$= \int_2^{\infty} \dots \int_2^{\infty} \prod_{i=2}^{1} \frac{1}{\beta} e^{-x_i/\beta} \, \int_2^{\infty} \frac{1}{\beta} e^{-x_1/\beta} \, \mathrm{d}x_1 \, \mathrm{d}x_2 \dots \, \mathrm{d}x_n$$
Integral over the exponential pdf
$$\int_2^{\infty} f(x_1) \, \mathrm{d}x_1 = 1 - F(2)$$

$$= 1 - (1 - e^{-\frac{1}{\beta}^2}) = e^{\frac{-2}{\beta}}$$

$$= e^{\frac{-2}{\beta}} \int_2^{\infty} \dots \int_2^{\infty} \prod_{i=2}^{1} \frac{1}{\beta} e^{-x_i/\beta} \, \mathrm{d}x_2 \dots \, \mathrm{d}x_n = \dots$$
Identical to original integral except that there are only n-1 terms to be integrated
$$= e^{-2/\beta} \cdot e^{-2/\beta} \int_2^{\infty} \dots \int_2^{\infty} \prod_{i=3}^{1} \frac{1}{\beta} e^{-x_i/\beta} \, \mathrm{d}x_3 \dots \, \mathrm{d}x_n$$

$$= (e^{-2/\beta})^n = e^{-2n/\beta}$$

A much more simpler solution: notice that $P(X_i > 2) = 1 - F(2) = e^{-1/\beta}$. Because X_i 's are independent, then also events $(X_i > 2)$ are independent event, and so

$$P(\text{All } X_i > 2) = \prod_{i=1}^n P(X_i > 2) = (e^{-2/\beta})^n = e^{-2n/\beta}$$

Illustration See Exponential Sample.R for the case where n=2 and $\beta=3$. Notice that R uses parametrization $\lambda=\frac{1}{\beta}$ for the exponential distribution, so in R, exponential pdf is $f(x\mid \lambda)=\lambda e^{-\lambda x}$.

Note: Definition 2.1 assumes that X_1, \ldots, X_n are independent. In practice, we often do analysis on dependent data. For that use, we could define term "dependent random sample". Mathematical treatment of dependent sample requires more specific description of dependence structure, e.g. through spatial or temporal autocorrelation models, or explicit models for grouped data.

Dependent data are modeled by assuming that random vectors X_1, \ldots, X_n are vectors of appropriate length, and there are n independent realizations of them in the data. Quite often n=1 and each replicate to X_1 , which is a rather long vector.

Example 2.2. Let $X = \begin{bmatrix} X(u_1) & X(u_2) & X(u_3) \end{bmatrix}^T$ include random variables X at locations u_1 , u_2 and u_3 . Assume that X_1 is normally distributed and the correlation $\rho_{ij} = corr(X(u_i), X(u_j))$ depends only on the spatial distance $s_{ij} = \|u_i - u_j\|$ between locations u_i and u_j .

The marginal means and variances of $X(u_i)$ are $E(X(u_i)) = \mu$ for all i and $Var(X(u_i)) = \sigma^2$ for all i. Consider case where $u_1 = (1,0), u_2 = (0,0), u_3 = (0,2)$ and $\rho_{ij} = e^{-s_{ij}/2}$. Correlations and covariances related to different random variables can be seen on table 2.1.

Table 2.1: Correlations between the random variable pairs of example 2.2

Pair	s_{ij}	$ ho_{ij}$	$cov(X(u_i), X(u_j))$
1,2	1	0.61	$0.61\sigma^2$
1,3	$\sqrt{5}$	0.33	$0.33\sigma^2$
2,3	2	0.37	$0.37\sigma^2$

Let

$$\mu = \begin{bmatrix} \mu \\ \mu \\ \mu \end{bmatrix}$$
 and $\Sigma = \sigma^2 \begin{bmatrix} 1 & 0.61 & 0.33 \\ 0.61 & 1 & 0.37 \\ 0.33 & 0.37 & 1 \end{bmatrix}$

The joint pdf of X is 3 variate normal distribution with mean μ and variance Σ

$$f(x_1 \mid \boldsymbol{\mu}, \boldsymbol{\Sigma}) = \frac{1}{\sqrt{(2\pi)^3}} |\boldsymbol{\Sigma}|^{-1/2} e^{\frac{1}{2}(\boldsymbol{x}_1 - \boldsymbol{\mu})' \boldsymbol{\Sigma}^{-1}(\boldsymbol{x}_1 - \boldsymbol{\mu})}$$

Note: In random sample, we assume that X_1, \ldots, X_n are identically distributed. In the case of random vectors, whis is specified by saying that the marginal distributions of the elements of X_i are identical.

Note: We can also have independent replicates of random vectors, e.g. if we have grouped data where the groups are independent replicates from the process that generates the groups, and the observations within the groups are dependent. This is related to mixed-effects models.

Note: Definition 2.1 specifies sampling from an infinite population (or population model): the sampling procedure does not change the population.

The population may also be finite, like numbers in hat. In that case, the sampling can be made with replacement: whenever a number has been drawn, the value is recorded but the number is put back to the hat. This procedure, which is useful e.g. in Bootstrapping, fulfils the conditions of definition 2.1.

If we do the sampling without replacement (which often makes much sense), then the conditions of definition 2.1 are not fulfilled: each draw changes the population by removing one unit from the finite population. This leads to so called design-based inference, which is covered in the literature of sampling theory.

The difference to definition 2.1 is important especially if the sample size n is large compared to the size of population, but may be unimportant if N >> n. In this course, we are considering only sampling according to definition 2.1.

Random samples can be summarized to a well defined summary called statistic.

Definition 2.2. Let $X = X_1, \dots, X_n$ be a random sample of size n from population f(x) and let T(X) be a real-valued or vector-valued function, whose domain includes

the sample space of X. The random variable Y = T(X) is called a statistic (statisti-ikka, tunnusluku). The probability distribution of Y is called the sampling distribution of Y (otantajakauma).

Examples of statistics

Sample minimum min(X)

Sample maximum max(X)

Sample median median(X)

Also other like sample mean, variance, standard deviation, quartiles etc.

Example 2.3. Let X be a random sample of size 5 from Exponential(2) population. R script statistics.R illustrates the distributions of min(X), max(X), median(X) and mean(X) through simulation. Script repeates the sampling from Exponential distribution M=10000 times and illustrates the distribution of the above mentioned sample statistic when n=5.

Definition 2.3. The sample mean is the statistic defined by

$$\bar{X} = \frac{1}{n} \sum_{i=1}^{n} X_i$$

Definition 2.4. The sample variance is the statistic defined by

$$S^{2} = \frac{1}{n-1} \sum_{i=1}^{n} (X_{i} - \bar{X})^{2}$$

- The sample standard deviation is the statistic defined by $S = \sqrt{S^2}$.
- The observed values of these random variables are denoted by \bar{X} , s^2 and s.

Theorem 2.1. Let X_1, \ldots, X_n be any numbers. Then

a)
$$min_a \sum_{i=1}^n (x_i - a)^2 = \sum_{i=1}^n (x_i - \bar{x})^2$$

b) $(n-1)s^2 = \sum_{i=1}^n (x_i - \bar{x})^2 = \underbrace{\sum_{i=1}^n x_i^2 - n\bar{x}^2}_{\text{Compare to}}$
 $Var(\mathbf{X}) = E(\mathbf{X}^2) - (E(\mathbf{X}))^2$

Lemma 2.1. Let X_1, \ldots, X_n be a random sample form a population and the g(X) be a function such that $E(g(X_i))$ and $Var(g(X_i))$ exist. Then

$$E\left[\sum_{i=1}^{n} g(X_i)\right] = n \cdot E(g(X_i))$$

and

$$Var(\sum_{i=1}^{n} g(X_i)) = n \cdot Var(g(X_i))$$

Proof

$$E(\sum g(X_i)) = \sum E(g(X_i)) = *$$

Because X_i 's are identically distributed

$$* = nE(g(X_i))$$

$$Var(\sum g(X_i)) = E\left[\sum g(X_i) - E(\sum g(X_i))\right]^2$$
$$= E\left\{\sum \left[\underbrace{g(X_i) - E(g(X_i))}_{t_i}\right]^2 = *\right\}$$

With t_i , we can get following

$$\begin{split} E(\sum t_i \sum t_i) \\ &= E(\sum t_i^2 + \sum_{i \neq j}^{\sum} t_i t_j) \\ &= \sum E(t_i^2) + \sum_{i \neq j}^{\sum} E(t_i t_j) \end{split}$$

which we can expand back to original equation

$$\begin{split} * &= \underbrace{\sum E(\left[g(X_i) - E(g(X_i))\right]^2)}_{\sum E(t_i^2)} + \underbrace{\sum \sum_{i \neq j} E\left[(g(X_i) - E(g(X_i)))(g(X_j) - E(g(X_j)))\right]}_{\sum_{i \neq j} E(t_i t_j)} \\ &= \sum Var(g(X_i)) + \sum_{i \neq j} cov(g(X_i), g(X_j)) \\ &= \underbrace{n \cdot Var(X_1)}_{\text{Because } X_i \text{'s}} + n(n-1) \cdot \underbrace{0}_{\substack{\text{Because } X_i \text{'s} \text{ are independent}}} = n \cdot Var\left[g(X_1)\right] \end{split}$$

Note: The proof of $E(\sum g(X_i))$ did not utilize the assumption of independence. Therefore it is valid also for dependent samples. The poof of $Var(\sum g(X_i))$ used the assumption of independence. Therefore it is not valid for dependent samples.

• For dependent samples, we get

$$Var(\sum g(X_i)) = \sum_{i=1}^{n} \sum_{j=1}^{n} cov \left[g(X_i), g(X_j) \right]$$

• Proof left as an exercise (recall, that $Var(X_i) = cov(X_i, X_i)$).

$$Var(\boldsymbol{x}) = \begin{bmatrix} Var(X_1) & cov(X_1, X_2) & \dots & cov(X_1, X_n) \\ cov(X_2, X_1) & Var(X_2) & \dots & cov(X_2, X_n) \\ \vdots & \vdots & \ddots & \vdots \\ cov(X_n, X_1) & cov(X_n, X_2) & \dots & Var(X_n) \end{bmatrix}$$

$$Var(\sum X_i) = \sum \sum cov(X_i, X_j)$$

• That is, the sum of all elements of matrix Var(X).

Note: This extends the rule

$$Var(X_1 + X_2) = Var(X_1) + Var(X_2) + 2 \cdot cov(X_1, X_2)$$

to general sum $\sum X_i$.

Theorem 2.2. Let X_1, \ldots, X_n be a random sample form a population with mean μ and variance $\sigma^2 < \infty$. Then

a)
$$E(\bar{X})=\mu$$

b) $Var(\bar{X})=\frac{\sigma^2}{n}$ $E(\bar{X}-\mu)^2$
c) $E(S^2)=\sigma^2$

• Proof for a and b are familiar from ISI1 course. Proof of c applier the theorem 2.1 and is left as an exercise.

Note: Because $E(\bar{X}) = \mu$, the statistic \bar{X} is said to be an unbiased estimator (harhaton estimatori) of population mean μ . Also S^2 is an unbiased estimator of population variance σ^2 .

Example 2.4. Illustrate these results through simulation. Intuitively, mean of an observed value of a statistic over a large number of samples should be close to expected value of the statistic in question. Therefore, (a), (b) and (c) of theorem 2.2 can be demonstrated by simulating M samples of a fixed size n and exploring how the means of the sample values of \bar{x} , $(\bar{x} - \mu)^2$ and s^2 behave as $M \to \infty$.

R-script demonstratebias . R implements this by assuming that the population has the Uniform(0,10) distribution and n=10.

2.1 Distribution of sum

Theorem 2.3. (Convolution formula)

If X and Y are independent, continuous random variables with pdf's $f_X(x)$ and $f_Y(y)$, then the pdf of the sum Z = x + Y is

$$f_Z(z) = \int_{-\infty}^{\infty} f_X(w) f_Y(z - w) dw$$

Proof: See Casella Berger, p 215-216.

Illustration: See example 1.20 of notes.pdf

Note: To compute the complete distribution of a sum $Z = \sum_{i=1}^{n}$, where X_i 's do not need to be identically distributed, but they are independent, we need to apply theorem 2.3 iteratively. E.g. to find the distribution of $X_1 + X_2 + X_3$, you may first find the distribution of $Z = X_1 + X_2$ using theorem 2.3 and thereafter use theorem 2.3 again to find the pdf of $Z + X_3$. This is not trivial in very general case, but there are easier ways to do this, e.g. when X_i 's are identically distributed and independent (i.i.d).

However, recall that in any case, the moments are easy

$$E(\sum X_i) = \sum E(X_i)$$

$$Var(\sum X_i) = \sum Var(X_i) \quad \text{(If X_i's are independent)}$$

2.2 Sampling from Normally distributed population

Theorem 2.4. Let $X_{n\times 1}$ be random sample of size n from a $N(\mu, \sigma^2)$ distribution and let $\bar{X} = \frac{1}{n} \sum_{i=1}^n X_i$ and $S^2 = \frac{1}{n-1} \sum_{i=1}^n (X_i - \bar{X})^2$. Then

- a) \bar{X} and S^2 are independent random variables.
- b) \bar{X} has $N(\mu, \frac{\sigma^2}{n})$
- c) $(n-1)S^2/\sigma^2$ has chi-squared distribution with parameter n-1. Parameter n-1 is commonly called the "degrees of freedom".

Note: The chi-squared distribution with p degrees of freedom has the pdf

$$f_X(x) = \frac{1}{\Gamma(p/2)} X^{p/2-1} e^{-x/2}$$
Gamma-function

• Gamma -function: $\Gamma(z) = \int_0^\infty x^{z-1} e^{-x} dx$

Lemma 2.2. We use notation χ_p^2 for a chi-squared random variables with p degrees of freedom

• If Z is a N(0,1) random variable, then $Z^2 \sim \chi_1^2$, that is, the square of a standard normal random variable is a chi-squared random variable.

• If X_1, \ldots, X_n are independent and $X_i \sim \chi^2_{p_i}$ then $\sum X_i \sim \chi^2_{\sum p_i}$. That is, independent chi-squared random variables add to a chi-squared random variable, and the degrees of freedom also add.

Example 2.5. Illustrate the result of theorem 2.4 using R-script. In file normalchi.R, we implement the following:

Repeat M times

- 1. Generate random variables $X_i \sim N(0, 1)$, where $n = 1, \dots, 9$
- 2. Compute $Z = \sum x_i^2$ for each sample
- 3. Plot the histogram of the M obtained values of Z and compare to the distribution χ_n^2 .

Note: If X_n is a random sample from $N(\mu, \sigma^2)$ population, then the random variable

$$Z = \frac{\bar{X} - \mu}{\sigma / \sqrt{n}} \sim N(0, 1)$$

where $\mu = E(\bar{X})$ and $\sigma/\sqrt{n} = Var(\bar{X})$.

- This transformation could be used to make inference on μ using the observed \bar{x} , if σ^2 is known.
- However, that is not the case, and therefore we use $\frac{\bar{X} \mu}{\sigma/\sqrt{n}}$ which can also be written as

$$\frac{(\bar{X} - \mu)/(\sigma/\sqrt{n})}{\sqrt{S^2/\sigma^2}}$$

where $(\bar{X}-\mu)/(\sigma/\sqrt{n})\sim N(0,1)$ and $\sqrt{S^2/\sigma^2}\sim \chi^2_{n-1}.$

Theorem 2.5. Let X be a random sample from $N(\mu, \sigma^2)$. The random variable

$$T = \frac{\bar{X} - \mu}{S/\sqrt{n}}$$

has Student's t-distribution with parameter (n-1) "degrees of freedom".

In general, a random variable T has Student's t-distribution with p degrees of freedom $(T \sim t_p)$ if T has the pdf

$$f_T(t) = \frac{\Gamma(\frac{p+1}{2})}{\Gamma(\frac{p}{2})} \frac{1}{(p\pi)^{1/2}} \frac{1}{(1+t^2/p)^{(p+1)/2}}$$

Proof: See Casella Berger, p.223-224.

Illustration: File normalchi.R

• We use samples of sizes $3, 4, \dots, 100$ and M = 10000.

- Plot empirical histogram of $\frac{\bar{X}-\mu}{s/\sqrt{n}}$
- Compare to t_{n-1} distribution and to $N(0, \sigma^2)$.
- $\mu = 5$ and $\sigma^2 = 3^2$.
- Because $S^2 \to \sigma^2$ as n increases, the difference between N(0,1) and t_{n-1} becomes meaningless as $n \to \infty$.

Note: t_p has only p moments. Especially $E(T_p) = 0$ if p > 1.

- A third important derived distribution under sampling from a normal population is the Snedecor's/Fisher's F-distribution.
- It is the theoretical distribution of the ratio of variances.

Example 2.6. Let X be a random sample from $N(\mu_X, \sigma_X^2)$ and let Y be a random sample from $N(\mu_Y, \sigma_Y^2)$. If we were interested in comparing the variability in these two populations, a quality of interest would be σ_X^2/σ_Y^2 . Information about σ_X^2/σ_Y^2 is contained in s_X^2/s_Y^2 ; ratio of the sample variances.

The F-distribution allows this comparison by giving te distribution for random variable

$$\frac{s_X^2/s_Y^2}{\sigma_X^2/\sigma_Y^2} = \frac{s_X^2/\sigma_X^2}{s_Y^2/\sigma_Y^2}$$

where ratios s_X^2/σ_X^2 and s_Y^2/σ_Y^2 are independent, scaled χ^2 -distributed random variables.

Definition 2.5. Let $X_{n\times 1}$ be a random sample from a $N(\mu_X, \sigma_X^2)$ population and let $Y_{m\times 1}$ be a random sample from an independent $N(\mu_Y, \sigma_X^2)$ population. The random variable

$$F = \frac{s_X^2/\sigma_X^2}{s_Y^2/\sigma_Y^2}$$

has the F-distribution with parameters (numerator and denominator degrees of freedom) n-1 and m-1. Equivalently, random variable F has the F-distribution with p and q degrees of freedom ($F \sim F_{p,q}$) if F has the pdf

$$f_F(x) = \frac{\Gamma(\frac{p+q}{2})}{\Gamma(\frac{p}{2})\Gamma(\frac{q}{2})} \left(\frac{p}{q}\right)^{p/2} \frac{x^{p/2-1}}{\left[1 + \frac{p}{q}x\right]^{(p+q)/2}}$$

Note:

- a) If $X \sim F_{p,q}$, then $1/X \sim F_{q,p}$
- b) If $X \sim t_q$, then $X^2 \sim \chi_1^2$ (recall that if $X \sim N(0,1)$, then $X^2 \sim \chi_1^2$)

2.3 Convergence

• What happens if $n \to \infty$.

Theorem 2.6. Markov inequality

Let X be a random variable such that $P(X \ge 0) = 1$ (i.e. X gets only positive values). Then, for every number t > 0

$$P(x > t) \le \frac{E(X)}{t}$$

Proof: Consider only case where *X* is discrete random variable.

$$\begin{split} E(X) &= \sum_X x f(x) \\ &= \sum_{x < t} x f(x) + \sum_{x \ge t} x f(x) \end{split}$$

Because $X \ge 0$, both terms are positive

$$E(X) \geq \sum_{X \geq t} x f(x) \geq \sum_{X \geq t} t f(x) = t \underbrace{\sum_{X \geq t} f(x)}_{P(X \geq t)}$$
$$= t \cdot P(X \geq t)$$

$$E(X) \ge t \cdot P(X \ge t) || : t \ (>0)$$

$$\frac{E(X)}{t} \ge P(X \ge t)$$

Example 2.7. Let X be a non-negative random variable with E(X) = 1

$$P(X \ge 100) \le \frac{E(X)}{100} = 0.01$$

Theorem 2.7. Chebyshev's inequality

Let X be a random variable such that Var(X) exists. Then for every number t>0

$$P(|X - E(X)| \ge t) \le \frac{\sigma^2}{t^2}$$

Proof: Let $Y = (X - E(X))^2$. Now $P(Y \ge 0) = 1$ and E(Y) = Var(X)

$$P(|X - E(X)| \ge t) = P(Y \ge t^2) \le \frac{E(Y)}{t^2}$$

$$P(|X - E(X)| \ge t) \le \frac{Var(X)}{t^2}$$

Example 2.8. If $Var(X) = \sigma^2$ and we select $t = 3\sigma$

$$P(|X - E(X)| \ge 3\sigma) \le \frac{\sigma^2}{(3\sigma)^2} = \frac{1}{9}$$

Definition 2.6. A sequence of random variables X_1, X_2, \ldots converges in probability (konvergoi todennäköisyysmielessä) to a random variable X if for every $\epsilon > 0$

$$\lim_{n \to \infty} P\left(|X_n - X| \ge \epsilon\right) = 0$$

or equivalently

$$\lim_{n \to \infty} P\left(|X_n - X| < \epsilon\right) = 1$$

A commonly used notation for this kind of convergence is $X_n \xrightarrow{P} X$. The sequence X_1, \ldots, X_n is not usually an i.i.d. sample but e.g. a statistic based on a sample of size n. Also, X is often a fixed constant, as it is in the following theorem.

Theorem 2.8. (The weak law of large numbers (Heikko suurten lukujen laki), WLLN) Let X_1, X_2, \ldots be i.i.d. random variables with expected value $E(X_i) = \mu$ and variance $Var(X_i) = \sigma^2 < \infty$. Define $\bar{X}_n = \frac{1}{n} \sum_{i=1}^n X_i$. Then, for every $\epsilon > 0$

$$\lim_{n \to \infty} P(|\bar{X}_n - \mu| < \epsilon) = 1$$

meaning sample mean convergences in probability to population mean $(\bar{X}_n \xrightarrow{P} \mu)$. We will have the sample mans of an i.i.d. sample arbitrarily close to μ if just n is high enough.

Proof: Use Chebyshev's inequality for the complement event:

$$P(|\bar{X}_n - \mu| \ge \epsilon) \le \frac{\overbrace{Var(\bar{X}_n)}^{-\sigma^2/n}}{\epsilon^2} = \frac{\sigma^2}{n\epsilon^2}$$

$$P(|\bar{X}_n - \mu| < \epsilon) = 1 - P(|\bar{X}_n - \mu| \ge \epsilon) \ge 1 - \frac{\sigma^2}{n\epsilon^2}$$
 where $\lim_{n \to \infty} \frac{\sigma^2}{n\epsilon^2} = 0$ and $\lim_{n \to \infty} 1 - \frac{\sigma^2}{n\epsilon^2} = 1$.

Note The property that the same sample quantity approaches a fixed constant as $n \to \infty$ is called consistency (konsistenssi).

Note WLLN was already used in R-script demonstrateBias.R

Note WLLN also justifies the use of a histogram from a large number of replicates as an approximation of the true density as $n \to \infty$. This is because every class of the histogram gives $P(c_1 \le x < c_1) = F(c_2) - F(c_1)$. Weather x belongs to a class $[c_1, c_2[$ is a Bernoulli(p) distributed random variable with

$$p = E(Y) = F(c_1) - F(c_2)$$

Note Another way of specifying the law of large numbers is the strong law of strong numbers (SLLN).

$$P(\lim_{n\to\infty}\bar{X}_n=\mu)=1.$$

 \bar{X} converges to μ with probability 1, meaning almost sure convergence.

Definition 2.7. A sequence of random variables converges in distribution to random variable X if

$$\lim_{n \to \infty} F_{X_n}(x) = F_X(x)$$

at all points of x where $F_X(x)$ is a continuous cdf. Notation used for convergence in distribution is $X_n \stackrel{d}{\to} X$.

Example 2.9. (Maximum of uniforms)

If X_1, X_2, \ldots are i.i.d. Uniform(0,1) distributed random variables and let $X_{(n)} = \max_{1 \le i \le n} X_i$. Let us explore if and to where $X_{(n)}$ converges in distribution.

When $n \to \infty$, $X_{(n)} \to 1$ and as $X_{(n)} < 1$, we have for any $\epsilon > 0$

$$P(|X_{(n)} - 1| \ge \epsilon) = P(1 - X_{(n)} \ge \epsilon) = P(X_{(n)} \le 1 - \epsilon)$$

Because sample is independent

$$\begin{split} P(X_{(n)} < 1 - \epsilon) &= P(\operatorname{All} X_i \leq 1 - \epsilon) \\ &= \underbrace{\left[\underbrace{P(X \leq 1 - \epsilon)}_{F_X(1 - \epsilon)}\right]^n}_{n} \quad X \sim Unif(0, 1) \\ &= (1 - \epsilon)^n \qquad \lim_{n \to \infty} (1 - \epsilon)^n = 0 \\ &\to X_{(n)} \xrightarrow{P} 1 \end{split}$$

Let us take $\epsilon = \frac{t}{n}$ to rewrite $P(X_{(n)} \le 1 - \epsilon)$ as

$$P(X_{(n)} \le 1 - \frac{t}{n}) = (1 - \frac{t}{n})^n$$

for which $\lim_{n\to\infty} (1-\frac{t}{n})^n = e^{-t}$.

$$\begin{split} X_{(n)} & \leq 1 - \frac{t}{n} \\ X_{(n)} - 1 & \leq -\frac{t}{n} \\ 1 - X_{(n)} & \leq \frac{t}{n} \\ n(1 - X_{(n)}) & \leq t \\ P(n(1 - X_{(n)}) & \leq t) = e^{-t} \\ P(n(1 - X_{(n)}) & \geq t) = \underbrace{1 - e^{-t}}_{\text{Exponential(1) cdf}} \end{split}$$

Illustration: See R-script MaxOfUniforms.R

How fast does the convergence occur, i.e. how large does the n need to be to have $n(1-X_{(n)}) \sim Exponential(1)$?

We generate M=10000 samples of each of the sample sizes n=2,3,4,5,10,20. For each sample, find $X_{(n)}$ and compute $y=n(1-X_{(n)})$. Plot the histogram if y and compare to Exponential(1) pdf.

The approximation looks quite nice already, when $n \ge 10$. But notice the behaviour in the tails. Especially $n(1 - X_{(n)})$ is bounded to interval [0, n].

Theorem 2.9. (The central limit theorem (CTL) (Keskeinen raja-arvolause))

Let X_1, X_2, \ldots be a sequence of i.i.d. random variables with finite $E(X) = \mu$ and finite variance $Var(X) = \sigma^2 > 0$. Let $\bar{X}_n = \sum_{i=1}^n X_i$. Let $G_{Z_n}(z)$ denote the cdf of

$$Z_n = \frac{\sqrt{n}(\bar{X}_n - \mu)}{\sigma} = \frac{\bar{X}_n - \overbrace{\mu}^{E(\bar{X}_n)}}{\underbrace{\sigma/\sqrt{n}}_{sd(\bar{X}_n)}}$$

Then $\lim_{n \to \infty} G_{Z_n} z = \underbrace{\int_{-\infty}^z \underbrace{\frac{1}{\sqrt{2\pi}} e^{-y^2/2} \, \mathrm{d}y}_{N(0,1) \, \mathrm{pdf}}}_{N(0,1) \, \mathrm{cdf}}$ That is: Z_n has limited standard normal

distribution, meaning $Z_n \xrightarrow{d} N(0,1)$. Equivalently, $\bar{X} \xrightarrow{d} N(\mu, \sigma^2/n)$. Notes:

- We assume only finite μ and σ^2 and end up to normlity.
- The rate of convergence is affected by the original distribution of X_i
 - The closer the distribution of X is the Normal curve, the faster the convergence is.
 - The rate needs to be evaluated case by case.
- Provides an all purpose approximation og the distribution of sums of i.i.d. random variables.

Proof: See Casella Berger pages 237-238.

Illustration: See R-script uniffmean.R. Script takes M=10000 samples using each of the following sample sizes n=1,2,3,4,5,6 from a Uniform(0,10) population, and demonstrates the distribution of \bar{X} using a histogram and compares it to $N(\mu,\sigma^2/n)$ distribution. Expected value and variance of Uniform(0,10) population are

$$\mu = \frac{b-a}{2} = \frac{10}{2} = 5$$

$$\sigma^2 = \frac{1}{12}(b-a)^2 = \frac{(10-0)^2}{12} = \frac{100}{12} = \frac{25}{3}$$

The visual evaluation of histogram gives an impression that the approximation is good already when n > 4. However, we are usually interested in the behaviour in the tails,

which needs more careful evaluation, e.g. compute the 0.95th quartile of the Normal distribution. Check which proportion of the simulated samples had the mean above this quartile. The proportion should be close to 0.05 if the approximation is good.

Theorem 2.10. (Slutsky) If $X_n \xrightarrow{d} X$ and if $Y_n \xrightarrow{P} a$, where a is a constant, then

a)
$$Y_n X_n \xrightarrow{d} aX$$

b)
$$X_n + Y_n \xrightarrow{d} X + a$$

Note Proof for this is not shown.

Example 2.10. (Normal approximation with estimated variance)

The central limit theorem said that under mild conditions

$$\frac{\sqrt{n}(\bar{X}_n - \mu)}{\sigma} \xrightarrow{d} N(0, 1)$$

However, σ^2 is often unknown, but can be unbiasedly and consistently estimated by $s_n^2 = \frac{1}{n-1} \sum (X_i - \bar{X}_n)^2$. It can also be shown that (see Casella Berger example 5.5.3 and exercise 3.32)

$$\frac{\sigma}{s_n} \xrightarrow{P} 1$$

Note

$$\frac{\sqrt{n}(\bar{X}_n - \mu)}{s_n} = \underbrace{\frac{\sigma^2}{s_n}}_{\frac{\sigma}{s_n} \xrightarrow{P} 1} \underbrace{\frac{\sqrt{n}(\bar{X}_n - \mu)}{\sigma^2}}_{\frac{\sqrt{n}(\bar{X}_n - \mu)}{\sigma} \xrightarrow{d} N(0, 1)}$$

So even tough the variance is estimated, $\frac{\bar{X}-\mu}{s/\sqrt{n}}$ converges to a normal distribution. However, intuitively, the convergence should be more slow than when σ is known.

Chapter 3

Estimation

In estimation, we want to use sample X_1, \ldots, X_n to make inference on an unknown population parameter θ . It means that we want to find a statistic $T(\mathbf{X})$ that somewhat optimally captures the information of the sample on the parameter of interest, θ .

Most commonly. θ is the population mean, but it can also be related to population variance, covariance, median, maximum ect. Very often the population parameters, especially the mean, is not a scalar number, but it's function of some known characteristics of sampled units i, and therefore θ includes parameters of that function.

Example 3.1. Simple linear regression

Consider case of simple linear regression, where X is a fixed known predictor, and Y is a random variable we want to predict with values of X. The model can be of following for

$$Y_i = \mu(X_i) + e_i$$

Where $\mu(X_i)$ is a fixed mean of Y

$$\mu_i = \mu(X_i) = \beta_0 + \beta_1 X_i$$

and e_i is the random error related to value of Y_i

$$e_i \sim N(0, \sigma^2)$$

This would mean that distribution of Y_i would be

$$Y_i \sim N(\mu(x_i), \sigma^2)$$

In this case, we are interested in estimating the the regression coefficients

$$oldsymbol{ heta} = egin{bmatrix} eta_0 \ eta_1 \end{bmatrix}$$

Example 3.2. Lets consider a random variable

$$Y_i = \begin{cases} 1 & \text{if tree } i \text{ is dead} \\ 0 & \text{if tree } i \text{ is alive} \end{cases}$$

Parameter of interest might be the probability of a tree being dead.

$$Y_i \sim Bernoulli(p), \ \theta = p$$

If we also know the age of the tree, and it is justified to assume that older trees are more commonly dead than younger trees, then

$$Y_i \sim Bernoulli(p_i)$$

$$p_i = \beta_0 + \beta_1 Age_i, \rightarrow \boldsymbol{\theta} = \begin{bmatrix} \beta_0 \\ \beta_1 \end{bmatrix}$$

Because $\beta_0 + \beta_1 \text{Age}_i$ is not restricted to [0, 1], a better model uses some function for $p(\text{Age}_i)$. Most commonly we use

$$log\left(\frac{p_i}{1-p_i}\right) = \beta_0 + \beta_1 Age_i$$

Which is a logit transformation $[0,1] \to \mathbb{R}$. This provides us with logistic regression. **Note:** See also example 1.2 and 1.3 in notes.pdf

3.1 Two principles

A sufficient statistic (tyhjentävä statistiikka) for a population parameter θ is such statistic $T(\boldsymbol{X})$, that in certain sense it captures all the information about θ contained in the sample.

3.2 The sufficient principle

If T(X) is a sufficient statistic for θ , then any inference about θ should depend on X only through the value of T(X).

Definition 3.1. A statistic T(X) is sufficient statistic for θ if the conditional distribution of the sample given T(X)[X|T(X)] does not depend on θ .

Example 3.3. Let X_1, \ldots, X_n be i.i.d. from $N(\mu, \sigma^2)$ population, where σ^2 is known. Now the sample mean \bar{X} is a sufficient statistic for μ .

That is, if we have the complete data and you know only the sample mean, we both still have all information on μ the sample includes. However, if a third person knows

only $min(\boldsymbol{X})$, $max(\boldsymbol{X})$ and $median(\boldsymbol{X})$, he still does not have much information on μ as you have when you know \bar{X} .

If also σ^2 is unknown, then $T(X) = \begin{bmatrix} \bar{X} \\ s^2 \end{bmatrix}$ is sufficient for $\theta = \begin{bmatrix} \mu \\ \sigma^2 \end{bmatrix}$ For more formal discussion on the topic, see Casella Berger section 6.2.

3.3 The likelihood function

Definition 3.2. Let $f(x, \theta)$ denote the joint pdf or pmf of the sample $X = X_1, \dots, X_n$. Then Given that X = x has been observed, the function of θ defined by

$$L(\boldsymbol{\theta}|\boldsymbol{x}) = f(\boldsymbol{x}|\boldsymbol{\theta})$$

is called the likelihood function (uskottavuusfunktio).

Note We do not assume X_1, \ldots, X_n to be independent.

 \bullet If X is a discrete random vector, then

$$L(\boldsymbol{\theta}|\boldsymbol{X}) = P_{\boldsymbol{\theta}}(\boldsymbol{X} = \boldsymbol{x})$$

• If we compare the likelihood at two parameter points $\theta = \theta_1$ and $\theta = \theta_2$, and find out that

$$P_{\theta_1}(\mathbf{X} = \mathbf{x}) = L(\theta_1|\mathbf{x}) > L(\theta_2|\mathbf{x}) = P_{\theta_2}(\mathbf{X} = \mathbf{x})$$

then the sample we observed is more likely to have occurred when $\theta = \theta_1$ than when $\theta = \theta_2$.

- $\rightarrow \theta_1$ is more plausible (uskottava) value for θ than θ_2 .
- If X is a continuous, real valued random variable, then for a small value ϵ

$$P_{\theta}(X - \epsilon < X < X + \epsilon) \approx 2\epsilon f(X, \theta) = 2\epsilon L(\theta|X)$$

therefore

$$LR = \frac{P_{\theta_1}(X - \epsilon < X < X + \epsilon)}{P_{\theta_2}(X - \epsilon < X < X + \epsilon)} \approx \frac{L(\theta_1|X)}{L(\theta_2|X)}$$

 \rightarrow If the LR > 1, then the sample was more likely to occur when $\theta = \theta_1$, and if LR < 1, then the sample was more likely to occur when $\theta = \theta_2$.

Example 3.4. Let X have negative binomial distribution. The Negative binomial distribution specifies the pmf for the number of successes before a fixed number of a failures (r) in a series of Bernoulli trials with success probability p. The pmf

$$P(X = x | r, p) = {r + x - 1 \choose x} p^r (1 - p)^x$$
, where $x = 0, 1, ...$ and $0 \le p \le 1$

Consider an experiment where r=3 and it was observed that X=2. The likelihood as a function of $\theta=p$

$$L(p|2) = P_p(X=2) = {4 \choose 2} p^3 (1-p)^2$$

 \rightarrow The likelihood is a 5th order polynomial with respect to p.

In general, if X = x has been observed,

$$L(p|x) = P_p(X = x) = {3 + x - 1 \choose 2} p^3 (1 - p)^x$$

in which case the likelihood is a polynomial of order x + 3.

Example 3.5. Let $X_{3\times 1}$ have 3-variate normal distribution so that all components have a common, unknown mean of μ and a known, common variance $\sigma^2=3.69$. In addition, $cov(X_1,X_2)=2.25$ and $cov(X_1,X_3)=cov(X_2,X_3)=0$.

This kind of situation could stem from a grouped structure of the data: X_1 and X_2 might belong to the same group (e.g. school class) and X_3 may originate from another group. That is

$$m{X} \sim N(m{\mu}, m{\Sigma})$$
, where $\mu = egin{bmatrix} \mu \\ \mu \\ \mu \end{bmatrix}$ and $m{\Sigma} = egin{bmatrix} 3.69 & 2.25 & 0 \\ 2.25 & 3.69 & 0 \\ 0 & 0 & 3.69 \end{bmatrix}$

The likelihood of this model for observation $\boldsymbol{x} = \begin{bmatrix} 6 \\ 5 \\ 4 \end{bmatrix}$ has been illustrated in R-script MVNlikelihood.R.

Note The likelihood function is not a pdf as a function of θ . Therefore we say, that $_1$ is more plausible than θ_2 (not more probable). Note also, that θ is thought to be fixed, therefore it makes no sense to specify a pdf for it. (In Bayesian statistics, θ is thought to be random.)

3.4 The likelihood principle

If x and y are two observed sample points such that $L(\theta|x)$ is proportional to $L(\theta|y)$, that is there exists a constant C(x, y) such that

$$L(\theta|\mathbf{x}) = C(\mathbf{x}, \mathbf{y})L(\theta|\mathbf{y})$$
 for all θ

then the conclusions drawn from x and y should be identical. Note C(x, y) may be different for different (x, y) pairs, but does not dependent on θ .

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3.5 Point estimation

Definition 3.3. A point estimator is any function $W(X_1, \dots, X_n)$. That is, any statistic is a point estimator.

 We now first consider the methods to find estimators and there after consider criteria to evaluate, whether the estimators are good or bad.

Note Estimator is a function of the sample X_1, \ldots, X_n and estimate is the numerical value the estimator got after the sample was actually taken.

- → Estimator is a random variable, and estimate is its realized value in the sample.
- Often, there is a natural, intuitive candidate fro a point estimator, e.g. \bar{X} is a intuitive choice for a point estimator of μ .

3.6 Estimating the population mean using least squares

Based on theorem 2.1, the ordinary least squares estimator of parameter μ using observed sample values y_1, \ldots, y_n is found by

$$min_{\theta} \sum_{i=1}^{n} (y_i - \theta)^2$$

The solution is $\theta = \bar{y}$ The approach also generalized to cases where $E(Y_i)$ is a function of a fixed variable x_i . Consider a model $Y_i = \beta_0 + \beta_1 x_i + \epsilon_i$, where Y_i is a random variable and x_i is another fixed observable variable. We can β_0 such that $E(\epsilon_i) = 0$ and

$$E(Y_i) = E(\underbrace{\beta_0 + \beta_1 x_i}_{\text{fixed}} + \underbrace{\epsilon_i}_{E(\epsilon_i) = 0}) = \beta_0 + \beta_1 x_i = E(Y_i | x_i) = \mu(x_i, \beta)$$

Parameters β are "second order" parameters. They quantify the (assumed) linear relationship between x_i and $E(Y_i)$. Using the idea of theorem 2.1, β can be estimated using

$$min_{\beta_0,\beta_1} \sum_{i=1}^{n} (y_i - (\beta_0 + \beta_1 x_i))^2$$

The solutions are

$$\widehat{\beta_1} = \frac{\sum (x_i - \bar{x})(y_i - \bar{y})}{\sum (y_i - \bar{y})^2}$$
 and $\widehat{\beta_0} = \bar{y} - \beta_1 \bar{x}$

• A more general formulation results by defining

$$m{X} = egin{bmatrix} 1 & X_1 \ dots & dots \ 1 & X_n \end{bmatrix}, \; m{Y} = egin{bmatrix} Y_1 \ dots \ Y_n \end{bmatrix}, \; m{eta} = egin{bmatrix} eta_0 \ eta_1 \end{bmatrix}, \; m{\epsilon} = egin{bmatrix} \epsilon_1 \ dots \ \epsilon_n \end{bmatrix},$$

and then the model is $oldsymbol{Y} = oldsymbol{X}'oldsymbol{eta} + oldsymbol{\epsilon}$

• The least squares estimator of β is found by minimizing

$$\epsilon' \epsilon = (y - X'\beta)(bmy - X'\beta)$$

- ullet The solution is $\widehat{oldsymbol{eta}}=(oldsymbol{X}'oldsymbol{X})^{-1}oldsymbol{X}'oldsymbol{y}$
- This is a general solution for the multiple regression case as well.