Introduktion til sandsynlighedsteori og statistik

January 7, 2018

1 Gennerelt

Bog

Instruktor:

Navn: Jeanett Pelck Petersen

Mail: 201303926@post.au.dk eller jeanett_92@hotmail.dk

Kontor: F1

2 Basic Concepts

- Probabillity theory is allows us to describe random phenomena in the world around us
- When we say something is random we say that our knowlegde of the outcome is limited

2.1 Sums

Geometric series formular for any $a, x \in \mathbb{R}$:

$$a + ax + ax^{2} + ax^{3} + \dots + ax^{n-1} = \sum_{k=0}^{n-1} ax^{k} = a \frac{1 - x^{n}}{1 - x}$$
 (1)

For |x| < 1

$$a + ax + ax^{2} + ax^{3} + \dots = \sum_{k=0}^{\infty} ax^{k} = a \frac{1}{1-x}$$
 (2)

2.2 Sets

$$A = \{x \mid x \text{ satisfies some property}\} \text{ or } A = \{x : x \text{ satisfies some property}\}$$
 (3)

- "|" and ":" are pronounced "such that"
- Two sets are equal if they have the same elements
- The universal set is the set of all things
 - E.g. if we roled a die the universal set would be: $S = \{1, 2, 3, 4, 5, 6\}$
 - Universal set is often denoted with S
- Vendiagrams is a digram showing the relationship between one or more sets and the universal set

2.2.1 Set operations

- A union of sets is written: $\bigcup_{i=1}^{n} A_i$
- The **complement** of a set *A* is denoted by A^c or \hat{A}
- Two sets are mutually exclusive or disjoined if they don't share any elements
- A collection of nonempty sets is a partition af a set A if they are disjoined and their union is A
- The **Multiplication principle** says that number of elements in $A \times B$ is $|A| \times |B|$

2.2.2 Cardinality of a set

- The **cardinary** of a set is basically the size of the set and is therefore denoted |A|
- If *A* is a finite set, the **cardinary** is simply the number of elements in the set
- Inclusion-exclusion principle: $|A \cup B| = |A| + |B| |A \cap B|$
- There are two types of infinte sets called **countable** (e.g. \mathbb{N} , \mathbb{Q} and \mathbb{Z}) and **uncountable** (e.g. \mathbb{R})
- Sets are **countable** if the following is true
 - 1. If it is a finite set or
 - 2. It can be put down one to one with the natural numbers which is called infinite countable

A set is called **uncountable** if it is not **countable**

- Any set on a real line [a,b],(a,b),[a,b),[a,b],(a,b), or (a,b) where a < b is uncountable
- Any subset of \mathbb{N} , \mathbb{Q} and \mathbb{Z} is uncountable
- Any subset of a countable set is countable.
- Any superset of an uncountable set is uncountable.
- If A_1, A_2, \cdots is a list of countable sets, then the set $\bigcup_i A_i$ is also countable.
- If A and B is countable then $A \times B$ is also countable

2.3 Functions

$$f: A \to B$$
 (4)

- The input of a function is called the domain
- The output of a function is called the codomain
- The range of the function is all the possible values of f(x)

2.4 Random experiments

- Outcome is the result of a random experiment
- The set of all possible outcomes is called the **sample space**
- When repeatting a random experiment several times each one of them is called a trial
- Our goal with random experiments is to assign a probability to certain events
- An **event** is a collection of certain outcomes

• A **probabillity** measure P(A) is assigned to an event A between 0 and 1 which is how likely the event is

Axioms of probability:

- Axiom 1: For any event, $P(A) \ge 0$
- Axiom 2: Probability of the sample space *S* is P(S) = 1
- Axiom 3: If $A_1, A_2, A_3 \cdots$ are disjoint events, then $P(A_1 \cup A_2 \cup A_3 \cdots) = P(A_1) + P(A_2), P(A_3) \cdots$

Rules from axioms of probability:

- 1. For any event $A,P(A^c) = 1 P(A)$
- 2. The probability of the empty set is zero, i.e., $P(\emptyset) = 0$
- 3. For any event A, $P(A) \le 1$
- 4. $P(A B) = P(A) P(A \cap B)$
- 5. $P(A \cup B) = P(A) + P(B) P(A \cap B)$
- 6. If $A \subset B$ then $P(A) \leq P(B)$

Notation: $-P(A \cap B) = P(A \text{ and } B) = P(A, B) - P(A \cup B) = P(A \text{ or } B)$

- To find the probality of a event there are usually two steps
 - 1. Use the specific information that we have about the random experiment
 - 2. Use the probability axioms
- A probabillity P on S is **discret** if S is countable
- If $A \subset S$ is an event, then A is also countable, and by the third axiom of probability we can write

$$P(A) = P(\bigcup_{s_j \in A} \{s_j\}) = \sum_{s_j \in A} P(s_j)$$

In a finite sample space *S*, where are outcomes are equally possible, the probability of any event *A* can be found by

$$P(A) = \frac{|A|}{|S|}$$

2.5 Conditional Probability

If *A* and *B* are two events in a sample space S, then the **conditional probability of** *A* **given** *B* is defined as

$$P(A|B) = \frac{P(A \cap B)}{P(B)}, \text{ when } P(B) > 0$$
(5)

- A conditional probability is itself a probability measure, so it satisfies the probability axioms
- When B is a subset of A the probability of A is 1

For tree events A, B and C P(C) > 0 we have:

- $P(A^c|C) = 1 P(A|C)$
- $P(\emptyset|C) = 0$
- $P(A|C) \le 1$
- $P(A B|C) = P(A|C) P(A \cap B|C)$
- $P(A \cup B|C) = P(A|C) + P(B|C) P(A \cap B|C)$
- $P(A|C) \leq P(B|C)$

Inclusion-exclusion principle:

- $P(A \cup B) = P(A) + P(B) P(A \cap B)$
- $P(A \cup B \cup C) = P(A) + P(B) + P(C) P(A \cap B) P(A \cap C) P(B \cap C) + P(A \cap B \cap C)$

Chain rule for conditional probability:

- 1. $P(A \cap B) = P(A)P(B|A) = P(B)P(A|B)$
- 2. $P(A_1 \cap A_2 \cap \cdots \cap A_n) = P(A_1)P(A_2|A_1)P(A_3|A_2,A_1)\cdots P(A_n|A_{n-1}A_{n-2}\cdots \cap A_n)$

2.5.1 Independence

Definition of independence:

$$P(A \cap B) = P(A)P(B) \land P(A|B) = P(A) \tag{6}$$

- If A, B events are disjoined where P(A), P(B) > 0, then A and B is not independent
- If A and B are independent then
 - 1. A and B^c are independent
 - 2. A^c and B are independent
 - 3. A^c and B^c are independent
- If A_1, A_2, \dots, A_n are independent then:

$$P(A_1 \cup A_2 \cup \dots \cup A_n) = 1 - (1 - P(A_1))(1 - P(A_2)) \cdots (1 - P(A_n)). \tag{7}$$

2.5.2 Total probability

Law of total probability - If B_1 , B_2 , B_3 , \cdots is a partition of the sample space S, then for any event A we have

$$P(A) = \sum_{i} P(A \cap B_i) = \sum_{i} P(A|B_i)P(B_i). \tag{8}$$

2.5.3 Bayers rule

• For any two events *A* and *B*, where $P(A) \neq 0$, we have:

$$P(B|A) = \frac{P(A|B)P(B)}{P(A)} \tag{9}$$

• If B_1, B_2, B_3, \cdots form the partition of the sample space S, and A is any event with $P(A) \neq 0$, we have

$$P(B_j|A) = \frac{P(A|B_j)P(B_j)}{\sum_{i} P(A|B_i)P(B_i)}$$
(10)

2.5.4 Conditional independence

Two events A and B are **conditionally independed** if given an event C where P(C) > 0

$$P(A \cap B|C) = P(A|C)P(B|C) \tag{11}$$

3 Combinators

- Multiplication Principle:
 - If we preform r experiments such that the kth experiment has n_k possible outcomes, for $k = 1, 2, \ldots, r$. The there are a total of $n_1 \cdot n_2 \cdot \cdots \cdot n_r$ possible outcomes for the sequence of r experiments
- Sampling from a set means choosing an element from that set.
- **Sampling with replacement** means that we put eachs objects back after each draw, in **sampling without replacement** we don't
- If order matters it is called **ordered sampling** otherwise, it is called **unordered**
- If we have *n* elements there is *n*! ways of ordering them

Sampling type	Posiblities
ordered sampling with replacement	n^k
ordered sampling without replacement	$P_k^n = \frac{n!}{(n-k)!}$
unordered sampling without replacement	$\binom{n}{k} = \frac{n!}{k!(n-k)!}$ $\binom{n+k-1}{k}$
unordered sampling with replacement	$\binom{n+k-1}{k}$
3	

3.1 Unordered sampling without replacement

- **K-permutation** of the elements in a set is choosing k elements, ordered with no repetion
- $\binom{n}{k}$ is called the **binomial coefficient**
- The total way to divide n distinct objects into two groups A and B such that group A consists of k objects and group B consists of n-k objects is $\binom{n}{k}$
- Binomial theorem states that for an integer n > 0

$$(a+b)^n = \sum_{k=0}^n \binom{n}{k} a^k b^{n-k}.$$
 (12)

- Binomial coefficient rules:
 - 1. We have $\sum_{k=0}^{n} {n \choose k} = 2^{n}$
 - 2. For $0 \le k < n$, we have $\binom{n+1}{k+1} = \binom{n}{k+1} + \binom{n}{k}$
 - 3. We have $\binom{m+n}{k} = \sum_{i=0}^k \binom{m}{i} \binom{n}{k-i}$ (Vandermonde's identity).
- A **Bernoulli Trial** is a random experiment that has two possible outcomes that we can label "success" or "failure" such as
 - Tossing a coin were you can define heads or tails as "failure" or "success"
 - You take a test, where the outcomes are pass or fail
- Success is denoted by *p*
- The probability of failure is denoted by q = p 1
- An experiment where *n* independent Bernoulli trials is performed and we count the number of success, it is called a **binomial** experiment
- **Binomial formula:** For n independent Bernoulli trials where each trial has success probability p, the probability of k successes is given by

$$P(k) = \binom{n}{k} p^k (1-p)^{n-k}$$

- Multinomial formula:
 - If $n = n1 + n2 + ... + n_r$, where all $ni \ge 0$ are integers, then the number of ways to divide n distinct objects to r distinct groups of sizes $n1, n2, ..., n_r$ is given by

$$\binom{n}{n_1, n_2, \dots, n_r} = \frac{n!}{n_1! n_2! \dots n_r!}.$$

• Multinomial theoroem:

$$(x_1 + x_2 + \dots + x_r)^n = \sum_{\substack{n_1 + n_2 + \dots + n_r = n \\ n_1, n_2, \dots, n_r}} \binom{n}{n_1, n_2, \dots, n_r} x_1^{n_1} x_2^{n_2} \dots x_r^{n_r}$$
(13)

Discrete Random Variables

4.1 Random variables

- A random variable is a real-valued function that assigns a numerical value to each possible outcome of the random experiment
- Random variables are shown by capital letters such as X, Y, Z where as number values are shown by small letters such as x, y, z
- A random variable X is a function from the sample space to real numbers

$$X: S \to \mathbb{R} \tag{14}$$

- The range of a random variable X is shown by Range(X) or R_X is the set of possible values of X
- *X* is a discrete random variable if its range is countable

4.2 Probability Mass Function (PMF)

• If *X* is a discrete random variable then we can list the elements in R_x :

$$R_x = \{x_1, x_2, x_3, \dots\} \tag{15}$$

• The event $A = \{X = x_k\}$ is defined as the set of outcomes s in the sample space S where there corresponding value of X is equal to x_k

$$A = \{ s \in S \mid X(s) = x_k \} \tag{16}$$

• The probabilities of events $\{X = x_k\}$ are formally shown by the **probability mass function (pmf)** of *X*

Definition 3.1

Let X be a discrete random variable with range $R_x = \{x_1, x_2, x_3, ...\}$. This function is called the probablity masse function (PMF) of *X*:

$$P_X(x_k) = P(X = x_k), \text{ for } k = 1, 2, 3, ...,$$
 (17)

- For discrete random variables the PMF is also called the probability distribution
- Properties of PMF:
 - 1. $0 \le P_X(x) \le 1 \text{ for all } x$

 - 2. $\sum_{x \in R_X} P_X(x) = 1$ 3. For any set $A \subset R_X$, $P(X \in A) = \sum_{x \in A} P_X(x)$
- If it is desired that the PMF to take numbers not in R_X we define $P_X(x)$ to be

$$P_X(x) = \begin{cases} P(X = x) & \text{if } x \text{ is in } R_X \\ 0 & \text{otherwise} \end{cases}$$

4.3 Independent Random Variables

Definition 3.2

Consider two discrete random variables *X* and *Y*, *X* and *Y* is independent if

$$P(X = x, Y = y) = P(X = x)P(Y = y), \quad \text{for all } x, y.$$
 (18)

• If two random variables are independent you can write

$$P(X \in A, Y \in B) = P(X \in A)P(Y \in B),$$
 for all sets A and B . (19)

• The following is also true is two random variables are independent

$$P(Y = y \mid X = x) = P(Y = y), \text{ for all } x, y.$$
 (20)

Definition 3.3

Consider n discrete random variables $X_1, X_2, X_3, ..., X_n$. We say that $X_1, X_2, X_3, ..., X_n$ are independent if

$$P\left(X_{1}=x_{1},X_{2}=x_{2},...,X_{n}=x_{n}\right)=P(X_{1}=x_{1})P(X_{2}=x_{2})...P(X_{n}=x_{n}), \text{ for all } x_{1},x_{2},...,x_{n}.$$
(21)

4.4 Special Distributions

4.4.1 Bernoulli Distribution

Definition 3.4

A random variable X is said to be a Bernoulli random variable with parameter p, shown as $X \sim Bernoulli(p)$, if its PMF is given by where 0

$$P_X(x) = \begin{cases} p & \text{for } x = 1\\ 1 - p & \text{for } x = 0\\ 0 & \text{otherwise} \end{cases}$$

- The bernoulli random variable is also called the **indicator** random variable.
- The indicator random variable *I*^A is defined for an event *A* as

$$I_A = \begin{cases} 1 & \text{if the event } A \text{ occurs} \\ 0 & \text{otherwise} \end{cases}$$

• The indicator random variable for an event A has the Bernoulli distribution with parameter p = P(A), so we can write

$$I_A \sim Bernoulli(P(A)).$$

4.4.2 Geometric Distribution

- A geometric distribution is can usually be thought of as repeating independent Bernoulli trials until the first success
- In a geometric distribution we usually use q as 1 p
- Some define the geometric district distribution as number of failures before success here it is defined as the numbers of experiments that lead to success

Definition 3.5

• A random variable X is said to be a geometric random variable with parameter p, shown as $X \sim Geometric(p)$ if its PMF is given by where 0

$$P_X(k) = \begin{cases} p(1-p)^{k-1} & \text{for } k = 1, 2, 3, \dots \\ 0 & \text{otherwise} \end{cases}$$

4.4.3 Binomial Distribution

• A random experiment involving a binomial distribution can be a coin that has P(H) = p. The coin is tossed n times and we define X to be the total number of heads observed.

Definition 3.6

• A random variable X is said to be a binomial random variable with parameters n and p, shown as $X \sim Binomial(n, p)$, if its PMF is given by where 0

$$P_X(k) = \begin{cases} \binom{n}{k} p^k (1-p)^{n-k} & \text{for } k = 0, 1, 2, \dots, n \\ 0 & \text{otherwise} \end{cases}$$

Lemma 3.1

• If $X_1, X_2, ..., X_n$ are independent Bernoulli(p) random variables, then the random variable X defined by $X = X_1 + X_2 + ... + X_n$ has a Binomial(n, p) distribution.

4.4.4 Negative Binomial (Pascal) Distribution

- The negative binomial distribution is a generalization of the geometric distribution.
- It relates to the random experiment of repeated independent trials until observing *m* successes.
- Pascal(1, p) = Geometric(p)

Definition 3.7

• A random variable X is said to be a Pascal random variable with parameters m and p, shown as $X \sim Pascal(m, p)$, if its PMF is given by where 0

$$P_X(k) = \begin{cases} \binom{k-1}{m-1} p^m (1-p)^{k-m} & \text{for } k = m, m+1, m+2, m+3, \dots \\ 0 & \text{otherwise} \end{cases}$$

4.4.5 Hypergeometric Distribution

Hypergeometric Distribution can be thought of as chosing k marbles at random out of a bag
with b blue maples and r red maples and then the random variable X is the number of blue

marbles

Definition 3.8

• A random variable X is said to be a Hypergeometric random variable with parameters b,r and k, shown as $X \sim Hypergeometric(b,r,k)$, if its range is $R_X = \{\max(0,k-r), \max(0,k-r) + 1, \max(0,k-r) + 2, ..., \min(k,b)\}$, and its PMF is given by

$$P_X(x) = \begin{cases} \frac{\binom{b}{x}\binom{r}{k-x}}{\binom{b+r}{k}} & \text{for } x \in R_X \\ 0 & \text{otherwise} \end{cases}$$

4.4.6 Poisson Distribution

 The poisson distribution is often used counting occurences of certain events in a inteval of time and space

Definition 3.9

• A random variable X is said to be a Poisson random variable with parameter λ , shown as $X \sim Poisson(\lambda)$, if its range is $R_X = \{0, 1, 2, 3, ...\}$, and its PMF is given by

$$P_X(k) = \begin{cases} \frac{e^{-\lambda}\lambda^k}{k!} & \text{for } k \in R_X \\ 0 & \text{otherwise} \end{cases}$$

4.5 Cumulative Distribution Function

Definition 3.10

The cumulative distribution function (CDF) of random variable *X* is defined as

$$F_X(x) = P(X \le x)$$
, for all $x \in \mathbb{R}$ (22)

• For all $a \leq b$, we have

$$P(a < X \le b) = F_X(b) - F_X(a)$$
(23)

4.6 Expectation

- The expected value is defined as the weighted value in the range
- Expected value = mean = average
- Different notations for expected value of *X*: $EX = E[X] = E(X) = \mu_X$

Definition 3.11

• Let X be a discrete random variable with range $R_X = \{x_1, x_2, x_3, ...\}$ (finite or countably infinite). The expected value of X, denoted by EX is defined as

$$EX = \sum_{x_k \in R_X} x_k P(X = x_k) = \sum_{x_k \in R_X} x_k P_X(x_k)$$
 (24)

Theorem 3.2

• Expectation is linear

- We have
 - * E[aX + b] = aEX + b for $a, b \in \mathbb{R}$
 - * $E[X_1 + X_2 + \cdots + X_n] = EX_1 + EX_2 + \cdots + EX_n$, for any set of random variables X_1, X_2, \cdots, X_n

4.7 Functions of Random Variables

- If *X* is a random variable and Y = g(X), then *Y* itself is a random variable.
 - Thus, we can talk about its PMF, CDF, and expected value.
 - The range of *Y* can be written as

$$R_Y = \{g(x) | x \in R_X\} \tag{25}$$

• If the PMF of X is already known the PMF of Y can be found by

$$P_{Y}(y) = P(Y = y)$$

$$= P(g(X) = y)$$

$$= \sum_{x:g(x)=y} P_{X}(x)$$
(26)

• Law of the unconscious statistician (LOTUS) for discrete random variables:

$$E[g(X)] = \sum_{x_k \in R_X} g(x_k) P_X(x_k)$$

4.8 Variance

• The variance of a random variable X, with mean $EX = \mu_X$, is defined as

$$Var(X) = E[(X - \mu_X)^2].$$

• The **standard deviation** of a random variable *X* is defined as

$$SD(X) = \sigma_X = \sqrt{Var(X)}$$

• Computational formula for the variance:

$$Var(X) = E[X^2] - [EX]^2$$

• **Theorem 3.3:** For a random variable *X* and real numbers *a* and *b*,

$$Var(aX + b) = a^2 Var(X)$$

• **Theorem 3.4:** If X_1, X_2, \dots, X_n are independent random variables and $X = X_1 + X_2 + \dots + X_n$, then

$$Var(X) = Var(X_1) + Var(X_2) + \cdots + Var(X_n)$$

Continuous Random Variables

• A random variable X with CDF $F_X(x)$ is said to be continuous if $F_X(x)$ is a continuous function for all $x \in \mathbb{R}$

Definition 4.2

• Consider a continuous random variable X with an absolutely continuous CDF $F_X(x)$. The function $f_X(x)$ is defined by:

$$f_X(x) = \frac{dF_X(x)}{dx} = F_X'(x)$$
, if $F_X(x)$ is differentiable at x

is called the probability density function (PDF) of X

• Since the PDF is the derivative of the CDF we can find the CDF in the following way

$$F_X(x) = \int_{-\infty}^x f_X(u) du.$$

• Consider a continuous random variable X with a PDF $f_X(x)$

- 1. $f_{-}X(x) \ge 0$ for all $x \in \mathbb{R}$ 2. $\int_{-\infty}^{\infty} f_X(u) du = 1$
- 3. $P(a < X \le b) = F_X(b) F_X(a) = \int_a^b f_X(u) du$ 4. For at set $A: P(X \in A) = \int_A f_X(u) du$

• The range of a continuous random variable is defined as the set of real numbers x for which the PDF is larger than zero

$$R_X = \{x | f_X(x) > 0\}$$

5.1 **Expected value and Variance**

• The expected value of a continuous random variable is defined as

$$EX = \int_{-\infty}^{\infty} x f_X(x) dx$$

• Law of the unconscious statistician (LOTUS) for continuous random variables:

$$E[g(X)] = \int_{-\infty}^{\infty} g(x) f_X(x) dx$$

12

• For a continuous random variable we can write variance as

$$Var(X) = E[(X - \mu_X)^2] = \int_{-\infty}^{\infty} (x - \mu_X)^2 f_X(x) dx$$
 (27)

$$= EX^{2} - (EX)^{2} = \int_{-\infty}^{\infty} x^{2} f_{X}(x) dx - \mu_{X}^{2}$$
 (28)

5.2 Functions

Theorem 4.1 - If X is a continuous random variable and $g : \mathbb{R} \to \mathbb{R}$ is a strictly monotonic differentiable function. Let Y = g(x). Then the PDF of Y is given by

$$f_Y(y) = \begin{cases} f_x(g^{-1}(x)) \cdot |g^{-1'}(y)| \\ 0 \end{cases}$$
 if $g(x) = y$ does not have a solution

Theorem 4.2 - If we have a random variable X with domain R_X and let Y = g(X). If we can partition R_X into a finite number of intervals where g(x) is strictly monotone and differentiable on each partition. Then the PDF for Y is given by

$$f_Y(y) = \sum_{i=1}^n \frac{f_X(x_i)}{|g'(x_i)|} = \sum_{i=1}^n f_X(x_i) \cdot \left| \frac{dx_i}{dy} \right|$$

where x_1, x_2, \ldots, x_n are real solutions to g(x) = y

5.3 Special Distributions

5.3.1 Uniform Distribution

• A continuous random variable X is said to have a uniform distribution over the interval [a, b], shown as $X \sim Uniform(a, b)$, if its PDF is given by

$$f_X(x) = \begin{cases} \frac{1}{b-a} & a < x < b \\ 0 & x < a \text{ or } x > b \end{cases}$$

• Its mean is given by

$$EX = \frac{a+b}{2}$$

• Its variance is given by

$$Var(X) = \frac{(b-a)^2}{12}$$

5.3.2 Exponential Distribution

• A continuous variable X is said to have an exponential distribution with parameter $\lambda > 0$, shown as $X \sim Exponential(\lambda)$ if its PDF is given by

$$f_X(x) = \begin{cases} \lambda e^{-\lambda x} & x > 0 \\ 0 & \text{otherwise} \end{cases}$$

- If $X \sim Exponential(\lambda)$ then $EX = \frac{1}{\lambda}$ and $Var(X) = \frac{1}{\lambda^2}$
- If X is exponential with a parameter $\lambda > 0$, then X is a **memoryless** random variable

$$P(X > x + a \mid X > a) = P(X > x), \text{ for } a, x \ge 0.$$

Normal (Gaussian) Distribution

• A continuous random variable Z is said to be a standard normal (standard Gaussian) random variable, shown as $Z \sim N(0,1)$ is its PDF is given by

$$f_Z(z) = \frac{1}{\sqrt{2\pi}} \exp\left\{-\frac{z^2}{2}\right\}, \quad \text{for all } z \in \mathbb{R}.$$

- If $Z \sim N(0,1)$, then EZ = 0 and Var(Z) = 1
- The CDF of the normal distribution is denoted Φ function:

$$\Phi(x) = P(Z \le x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{x} \exp\left\{-\frac{u^2}{2}\right\} du.$$

- Some properties of the Φ function
 - 1. $\lim_{x \to \infty} \Phi(x) = 1, \lim_{x \to -\infty} \Phi(x) = 0$

 - 2. $\Phi(0) = \frac{1}{2}$ 3. $\Phi(-x) = 1 \Phi(x)$ for all $x \in \mathbb{R}$
- If Z is a standard normal random variable and $X = \sigma Z + \mu$, then X is a random normal variable with mean μ and variance σ^2

$$X \sim N(\mu, \sigma)$$

• If *X* is a normal random variable with mean μ and variance σ^2 i.e. $X \sim N(\mu, \sigma^2)$ then

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

$$F_X(x) = P(X \le x) = \Phi\left(\frac{x-\mu}{\sigma}\right)$$
,

$$P(a < X \le b) = \Phi\left(\frac{b-\mu}{\sigma}\right) - \Phi\left(\frac{a-\mu}{\sigma}\right).$$

• $X \sim N(\mu_X, \sigma_X^2)$ and Y = aX + b where $a, b \in \mathbb{R}$ then $Y \sim N(\mu_Y, \sigma_Y^2)$ where

$$\mu_Y = a\mu_X + b$$
, $\sigma_Y^2 = a^2\sigma_X^2$.

5.3.4 Gamma Distribution

• A continuous random variable is said to have a gamma distribution with parameters a > 0and $\lambda > 0$, shown as $X \sim Gamma(\alpha, \lambda)$ if its PDF is given by

$$f_X(x) = \begin{cases} \frac{\lambda^{\alpha} x^{\alpha - 1} e^{-\lambda x}}{\Gamma(\alpha)} & x > 0 \\ 0 & \text{otherwise} \end{cases}$$

• A gamma distribution with parameter one is the an exponential distribution $Gamma(1, \lambda) =$ Exponential(λ)

• If $X \sim Gamma(\alpha, \lambda)$ then

$$EX = \frac{\alpha}{\lambda}, \quad Var(X) = \frac{\alpha}{\lambda^2}.$$

• For any positive real number a:

1.
$$\Gamma(\alpha) = \int_0^\infty x^{\alpha-1} e^{-x} dx$$

2.
$$\int_0^\infty x^{\alpha-1} e^{-\lambda x} dx = \frac{\Gamma(\alpha)}{\lambda^{\alpha}}, \text{ for } \lambda > 0;$$
3.
$$\Gamma(\alpha+1) = \alpha \Gamma(\alpha);$$

3.
$$\Gamma(\alpha + 1) = \alpha \Gamma(\alpha)$$
;

4.
$$\Gamma(n) = (n-1)!$$
, for $n = 1, 2, 3, \dots$;

5.
$$\Gamma(\frac{1}{2}) = \sqrt{\pi}$$

5.4 Mixed random variables

 There is some random variables which is neither continuous or discrete they are called Mixed random variables

- They are a mixture of both

 The CDF of a mixed random variable can be written as the sum of a continuous and a staircase function

$$F_Y(y) = C(y) + D(y)$$

• If differentiate the continuous part of the CDF we get c(y) which is not a valid CDF because it does not summarize to one

$$c(y) = \frac{dC(y)}{dy}$$
, wherever $C(y)$ is differentiable.

• Let $\{y_1, y_2, y_3, ...\}$ be a set of jump points of D(y) for which $P(X = y_k) > 0$ We then have

$$\int_{-\infty}^{\infty} c(y)dy + \sum_{y_k} P(Y = y_k) = 1.$$

The expected value can be obtained as

$$EY = \int_{-\infty}^{\infty} yc(y)dy + \sum_{y_k} y_k P(Y = y_k)$$

5.4.1 The delta function

Definition 4.3

We define the delta function $\delta(x)$ as a function with the following properties:

1.
$$\delta(x) = \begin{cases} \infty & x = 0 \\ 0 & \text{otherwise} \end{cases}$$

2. $\delta(x) = \frac{d}{dx}u(x)$ where u(x) is a unit step

3.
$$\int_{-\epsilon}^{\epsilon} \delta(x) dx = 1$$
 for any $\epsilon > 0$

4. For any $\epsilon > 0$ and any function g(x) that is continuous over $(x_0 - \epsilon, x_0 + \epsilon)$ we have

$$\int_{-\infty}^{\infty} g(x)\delta(x-x_0)dx = \int_{x_0-\epsilon}^{x_0+\epsilon} g(x)\delta(x-x_0)dx = g(x_0)$$
 (29)

• For a discrete random variable X with range $RX = \{x_1, x_2, x_3, \dots\}$ and PMF $PX(x_k)$, we define the (generalized) probability density function (PDF) as

$$f_X(x) = \sum_{x_k \in R_X} P_X(x_k) \delta(x - x_k).$$

• The (generalized) PDF of a mixed random variable can be written in the form

$$f_X(x) = \sum_k a_k \delta(x - x_k) + g(x),$$

• where ak = P(X = xk), and $g(x) \ge 0$ does not contain any delta functions. Furthermore, we have

$$\int_{-\infty}^{\infty} f_X(x) dx = \sum_{k} a_k + \int_{-\infty}^{\infty} g(x) dx = 1$$

6 Joint Distributions

6.1 For two discrete variables

6.1.1 Joint PMF

• The **joint probability mass function** of two discrete random variable *X* and *Y* is defined as

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

• The joint range for *X* and *Y* can be defined as

$$R_{XY} = \{(x, y) \mid P_{XY}(x, y) > 0\}$$

• For two discrete random variables we have

$$\sum_{(x_i, y_j) \in R_{XY}} P_{XY}(x_i, y_j) = 1$$

• To find $P((X,Y) \in A \text{ for any set } A \subset \mathbb{R} \text{ we have}$

$$P((X,Y) \in A) = \sum_{(x_i,y_j) \in (A \cap R_{XY})} P_{XY}(x_i,y_j)$$

• The marginal PMFs of *X* and *Y*

$$P_X(x) = \sum_{y_j \in R_Y} P_{XY}(x, y_j), \quad \text{for any } x \in R_X$$

$$P_Y(y) = \sum_{x_i \in R_X} P_{XY}(x_i, y), \quad \text{for any } y \in R_Y$$
(30)

6.1.2 Joint CDF

• The **joint cumulative distribution function** of two random variables *X* and *Y* is defined as:

$$F_{XY}(x,y) = P(X \le x, Y \le y).$$

• Marginal CDFs of *X* and *Y*:

$$F_X(x) = F_{XY}(x, \infty) = \lim_{y \to \infty} F_{XY}(x, y), \quad \text{for any } x,$$

$$F_Y(y) = F_{XY}(\infty, y) = \lim_{x \to \infty} F_{XY}(x, y), \quad \text{for any } y$$
 (31)

For a joint CDF the following must be true

$$F_{XY}(\infty, \infty) = 1,$$

 $F_{XY}(-\infty, y) = 0,$ for any y ,
 $F_{XY}(x, -\infty) = 0,$ for any x .

• **Lemma 5.1:** For two random variables X and Y, and real numbers $x1 \le x2$, $y1 \le y2$, we have

$$P(x_1 < X \le x_2, y_1 < Y \le y_2) = F_{XY}(x_2, y_2) - F_{XY}(x_1, y_2) - F_{XY}(x_2, y_1) + F_{XY}(x_1, y_1).$$

6.1.3 Conditioning and Independence

• For a discrete variable *X* and a event *A*, the **conditional PMF** of *X* given *A* is defined as

$$\begin{aligned} P_{X|A}(x_i) &= P(X = x_i | A) \\ &= \frac{P(X = x_i \text{ and } A)}{P(A)}, \text{ for any } x_i \in R_X. \end{aligned}$$

• The **conditional CDF** of *X* given *A* is defined as

$$F_{X|A}(x) = P(X \le x|A).$$

• For discrete variables *X* and *Y* the **conditional PMF** of *X* given *Y* and vice versa is defined as

$$P_{X|Y}(x_i|y_j) = \frac{P_{XY}(x_i, y_j)}{P_Y(y_j)},$$

 $P_{Y|X}(y_j|x_i) = \frac{P_{XY}(x_i, y_j)}{P_X(x_i)}$

for any $x_i \in R_X$ and $y_i \in R_Y$.

X

• Two discrete random variables *X* and *Y* are independent if

$$P_{XY}(x,y) = P_X(x)P_Y(y)$$
, for all x,y .

Or

$$F_{XY}(x,y) = F_X(x)F_Y(y)$$
, for all x,y .

• Conditional expectation of X

$$E[X|A] = \sum_{x_i \in R_X} x_i P_{X|A}(x_i),$$

$$E[X|Y = y_j] = \sum_{x_i \in R_X} x_i P_{X|Y}(x_i|y_j)$$

• Law of total probability:

$$P(X \in A) = \sum_{y_j \in R_Y} P(X \in A | Y = y_j) P_Y(y_j)$$
, for any set A .

- Law of total expectation
 - 1. If B_1, B_2, B_3, \ldots is a partition of the sample space S

$$EX = \sum_{i} E[X|B_i]P(B_i)$$

2. For random variable *X* and discrete variable *Y*

$$EX = \sum_{y_j \in R_Y} E[X|Y = y_j] P_Y(y_j)$$

6.1.4 Functions of Two Random Variables

• For two random variables X and Y and Z = g(X, Y) where $g : \mathbb{R}^2 \to \mathbb{R}$

$$P_Z(z) = P(g(X,Y) = z)$$

$$= \sum_{(x_i,y_i) \in A_z} P_{XY}(x_i,y_j), \text{ where } A_z = \{(x_i,y_j) \in R_{XY} : g(x_i,y_j) = z\}.$$

• Law of the unconscious statistician (LOTUS) for two discrete random variables:

$$E[g(X,Y)] = \sum_{(x_i,y_i) \in R_{XY}} g(x_i,y_j) P_{XY}(x_i,y_j)$$

6.1.5 Conditional Expectation

• Let *X* and *Y* be two random variables and *g* and *h* be two functions then the following is true

$$E[g(X)h(Y)|X] = g(X)E[h(Y)|X]$$

• Law of Iterated Expectations

$$E[X] = E[E[X|Y]]$$

- If *X* and *Y* are independent variables then
 - 1. E[X|Y] = EX
 - 2. E[g(X)|Y] = E[g(X)]
 - 3. E[XY] = EXEY
 - 4. E[g(X)h(Y)] = E[g(X)]E[h(Y)]
- Law of Total Variance

$$Var(X) = E[Var(X|Y)] + Var(E[X|Y])$$

6.2 For two continuous random variables

6.2.1 Joint PDF

• Two random variable X and Y are **jointly continuous** if there exists a nonnegative function such that we have $f_{XY} : \mathbb{R}^2 \to \mathbb{R}$, such that for any set $A \in \mathbb{R}^2$ we have:

$$P((X,Y) \in A) = \iint_A f_{XY}(x,y) dxdy$$

- The function $f_{XY}(x, y)$ is called the **joint probability density function (PDF)** of *X* and *Y*
- The range of (X, Y) in a joint PDF is

$$R_{XY} = \{(x,y) \mid f_{X,Y}(x,y) > 0\}.$$

• The following must be true for a joint PDF

$$\int_{-\infty}^{\infty} \int_{-\infty}^{\infty} f_{XY}(x, y) dx dy = 1$$

• Marginal PDFs

$$f_X(x) = \int_{-\infty}^{\infty} f_{XY}(x, y) dy$$
, for all x , $f_Y(y) = \int_{-\infty}^{\infty} f_{XY}(x, y) dx$, for all y .

6.2.2 Joint CDF

• The **joint cumulative function** of two random variables *X* and *Y* is defined as:

$$F_{XY}(x,y) = P(X \le x, Y \le y).$$

- The joint CDF satisfies the following properties
 - 1. $F_X(x) = F_{XY}(x, \infty)$ for any x (marginal CDF of X)
 - 2. $F_Y(y) = F_{XY}(\infty, y)$ for any y (marginal CDF of Y)
 - 3. $F_{XY}(\infty,\infty)=1$

 - 4. $F_{XY}(-\infty, y) = F_{XY}(x, -\infty) = 0$ 5. $P(x_1 < X \le x_2, y_1 < Y \le y_2) =$ $F_{XY}(x_2,y_2) - F_{XY}(x_1,y_2) - F_{XY}(x_2,y_1) +$
 - 6. If *X* and *Y* are independent then $F_{XY}(x,y) = F_X(x)F_Y(y)$

Conditioning and independence

• If X is a continuous random variable, and A is an event that a < X < b (where possibly $a = -\infty$ or $b = \infty$, then

$$F_{X|A}(x) = \begin{cases} 1 & x > b \\ \frac{F_X(x) - F_X(a)}{F_X(b) - F_X(a)} & a \le x < b \\ 0 & x < a \end{cases}$$

$$f_{X|A}(x) = \begin{cases} \frac{f_X(x)}{P(A)} & a \le x < b \\ 0 & \text{otherwise} \end{cases}$$

• In general for a random variable *X* and an event *A* we have the following

$$E[X|A] = \int_{-\infty}^{\infty} x f_{X|A}(x) dx,$$

$$E[g(X)|A] = \int_{-\infty}^{\infty} g(x) f_{X|A}(x) dx,$$

$$Var(X|A) = E[X^{2}|A] - (E[X|A])^{2}$$

- For two jointly continuous random variables *X* and *Y*, we have:
 - 1. The conditional PDF of *X* given Y = y:

$$f_{X|Y}(x|y) = \frac{f_{XY}(x,y)}{f_{Y}(y)}$$

2. The conditional probability that $X \in A$ given Y = y

$$P(X \in A|Y = y) = \int_{A} f_{X|Y}(x|y)dx$$

3. The conditional CDF of X given Y = y

$$F_{X|Y}(x|y) = P(X \le x|Y = y) = \int_{-\infty}^{x} f_{X|Y}(x|y) dx$$

- For two jointly continuous random variables *X* and *Y*, we have:
 - 1. Expected value of X given Y = y

$$E[g(X)|Y = y] = \int_{-\infty}^{\infty} g(x) f_{X|Y}(x|y) dx$$

2. Conditional LOTUS

$$E[g(X)|Y = y] = \int_{-\infty}^{\infty} g(x) f_{X|Y}(x|y) dx$$

3. Conditional variance of *X* given Y = y

$$Var(X|Y = y) = E[X^{2}|Y = y] - (E[X|Y = y])^{2}$$

• Two continuous random variables *X* and *Y* are independent if and only if

$$f_{XY}(x,y) = f_X(x)f_Y(y)$$
, for all x,y

$$F_{XY}(x,y) = F_X(x)F_Y(y)$$
, for all x,y

• If two continuous random variables *X* and *Y* are independent then the following is true

$$E[XY] = EXEY,$$

$$E[g(X)h(Y)] = E[g(X)]E[h(Y)].$$

• Law of Total Probability

$$P(A) = \int_{-\infty}^{\infty} P(A|X=x) f_X(x) \ dx$$

• Law of Total Expectation

$$E[Y] = \int_{-\infty}^{\infty} E[Y|X = x] f_X(x) \ dx = E[E[X|Y]]$$

• Law of Total Variance

$$Var(Y) = E[Var(Y|X)] + Var(E[Y|X])$$

6.2.4 Functions of two random variables

LOTUS for two continuous random variables

$$E[g(X,Y)] = \int_{-\infty}^{\infty} \int_{-\infty}^{\infty} g(x,y) f_{XY}(x,y) \ dxdy$$

• **Theorem 5.1** Let X and Y be two jointly continuous random variables. Let $(Z, W) = g(X, Y) = (g_1(X, Y), g_2(X, Y))$, where $g : \mathbb{R}^2 \to \mathbb{R}^2$ is a continuous one-to-one (invertible) function with continuous partial derivatives. Let $h = g^{-1}$, i.e., $(X, Y) = h(Z, W) = (h_1(Z, W), h_2(Z, W))$. Then Z and W are jointly continuous and their joint PDF, $f_{ZW}(z, w)$, for $(z, w) \in R_{ZW}$ is given by

$$f_{ZW}(z, w) = f_{XY}(h_1(z, w), h_2(z, w))|J|,$$

where *J* is the Jacobian of *h* defined by

$$J = \det \begin{bmatrix} \frac{\partial h_1}{\partial z} & \frac{\partial h_1}{\partial w} \\ \frac{\partial h_2}{\partial z} & \frac{\partial h_2}{\partial w} \end{bmatrix} = \frac{\partial h_1}{\partial z} \cdot \frac{\partial h_2}{\partial w} - \frac{\partial h_2}{\partial z} \frac{\partial h_1}{\partial w}.$$

• Let *X* and *Y* be two jointy continous random variables and Z = X + Y, then

$$f_Z(z) = \int_{-\infty}^{\infty} f_{XY}(w, z - w) dw = \int_{-\infty}^{\infty} f_{XY}(z - w, w) dw.$$

If X and Y are also independed then

$$f_Z(z) = f_X(z) * f_Y(z)$$

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

• Therem 5.2: If $X \sim N(\mu_X, \sigma_X^2)$ and $Y \sim N(\mu_Y, \sigma_Y^2)$ are independed then

$$X + Y \sim N\left(\mu_X + \mu_Y, \sigma_X^2 + \sigma_Y^2\right).$$

6.3 Covariance and Correlation

• The covariance between *X* and *Y* is defined as

$$Cov(X,Y) = E[(X - EX)(Y - EY)] = E[XY] - (EX)(EY)$$

The covariance has the following properties

- 1. Cov(X, X) = Var(X)
- 2. If *X* and *Y* are independed then Cov(X, Y) = 0
- 3. Cov(X, Y) = Cov(Y, X)
- 4. Cov(aX, Y) = aCov(X, Y)
- 5. Cov(X + c, Y) = Cov(X, Y)
- 6. Cov(X + Y, Z) = Cov(X, Z) + Cov(Y, Z)
- 7. More generally

$$\operatorname{Cov}\left(\sum_{i=1}^{m} a_i X_i, \sum_{j=1}^{n} b_j Y_j\right) = \sum_{i=1}^{m} \sum_{j=1}^{n} a_i b_j \operatorname{Cov}(X_i, Y_j).$$

- If Cov(X, Y) = 0 X and Y are not necessarily independed
- Generally for $a, b \in \mathbb{R}$

$$Var(aX + bY) = a^{2}Var(X) + b^{2}Var(Y) + 2abCov(X, Y)$$

- The **correlation coefficient** of two random variables *X* and *Y* is defined as the covariance of the standardized versions of *X* and *Y*
 - The sandardized versions of *X* and *Y* is defined as

$$U = \frac{X - EX}{\sigma_X}, \quad V = \frac{Y - EY}{\sigma_Y}$$

• The **correlation coefficient** denoted by ρ_{XY} or $\rho(X,Y)$ is obtained by normalizing the covariance

$$\rho_{XY} = \rho(X, Y) = \frac{\text{Cov}(X, Y)}{\sqrt{\text{Var}(X) \text{Var}(Y)}} = \frac{\text{Cov}(X, Y)}{\sigma_X \sigma_Y}$$

- Properties of the correlation coefficient
 - 1. $-1 \le \rho(X, Y) \le 1$
 - 2. If $\rho(X,Y) = 1$ then Y = aX + b, where a > 0
 - 3. If $\rho(X,Y) = -1$ then Y = aX + b, where a < 0
 - 4. $\rho(aX + b, cY + d) = \rho(X, Y)$ for a, c > 0
- **Definition 5.2:** Consider two random variable *X* and *Y*:
 - If $\rho(X,Y) = 0$ we say that X and Y are **uncorrelated**
 - If $\rho(X,Y) > 0$ we say that X and Y are **positivly** correlated
 - If $\rho(X,Y)$ < 0 we say that X and Y are **negativly** correlated
- In general the following is true about two random variables X and Y

$$Var(X + Y) = Var(X) + Var(Y) + 2Cov(X, Y)$$

• If *X* and *Y* are uncorrelated then

$$Var(X + Y) = Var(X) + Var(Y)$$

• More generally if $X_1, X_2, ..., X_n$ are pairwise uncorrelated i.e. $\rho(X_i, X_i) = 0$ for $i \neq j$, then

$$Var(X_1 + X_2 + ... + X_n) = Var(X_1) + Var(X_2) + ... + Var(X_n).$$

• If *X* and *Y* are independed then they are uncorrelated

7 Multiple random variables

- All the concepts used for two random variables can be extended to more variables
- Let $X_1, X_2, ..., X_n$ be n discrete variables then the joint PMF of $X_1, X_2, ..., X_n$ is defined as

$$P_{X_1,X_2,...,X_n}(x_1,x_2,...,x_n) = P(X_1 = x_1,X_2 = x_2,...,X_n = x_n).$$

• For n jointly continuous random variables $X_1, X_2, ..., X_n$ the joint PDF is defined to be the function $f_{X_1X_2...X_n}(x_1, x_2, ..., x_n)$ such that the probability of any set $A \subset \mathbb{R}^n$ is given by the integral of the PDF over the set A. In particular, for a set $A \subset \mathbb{R}^n$, we can write

$$P\bigg((X_1,X_2,\cdots,X_n)\in A\bigg)=\int\cdots\int_A\cdots\int f_{X_1X_2\cdots X_n}(x_1,x_2,\cdots,x_n)dx_1dx_2\cdots dx_n.$$

• The marginal PDF of X_i can be obtained by integrating all other X_i 's

$$f_{X_1}(x_1) = \int_{-\infty}^{\infty} \cdots \int_{-\infty}^{\infty} f_{X_1 X_2 \dots X_n}(x_1, x_2, \dots, x_n) dx_2 \cdots dx_n$$

• The join CDF of *n* random variables $X_1, X_2, ..., X_n$ is defined as

$$F_{X_1,X_2,...,X_n}(x_1,x_2,...,x_n) = P(X_1 \le x_1,X_2 \le x_2,...,X_n \le x_n)$$

7.1 Independence

• Random variables $X_1, X_2, ..., X_n$ are independent if for all $(x_1, x_2, ..., x_n) \in \mathbb{R}^n$

$$F_{X_1,X_2,...,X_n}(x_1,x_2,...,x_n) = F_{X_1}(x_1)F_{X_2}(x_2)\cdots F_{X_n}(x_n).$$

• If $X_1, X_2, ..., X_n$ are discrete, then they are independent if for all $(x_1, x_2, ..., x_n) \in \mathbb{R}^n$, we have

$$P_{X_1,X_2,...,X_n}(x_1,x_2,...,x_n) = P_{X_1}(x_1)P_{X_2}(x_2)\cdots P_{X_n}(x_n).$$

• If $X_1, X_2, ..., X_n$ are continuous, then they are independent if for all $(x_1, x_2, ..., x_n) \in \mathbb{R}^n$, we have p

$$f_{X_1,X_2,...,X_n}(x_1,x_2,...,x_n) = f_{X_1}(x_1)f_{X_2}(x_2)\cdots f_{X_n}(x_n)$$

• If random variables X_1, X_2, \dots, X_n are independent, then we have

$$E[X_1X_2\cdots X_n]=E[X_1]E[X_2]\cdots E[X_n].$$

Definition 6.1

 Random variables are said to be independent and identically distributed (i.i.d.) if they are independent and have the same marginal distributions:

$$F_{X_1}(x) = F_{X_2}(x) = ... = F_{X_n}(x)$$
, for all $x \in \mathbb{R}$

Then the following must also be true

$$E[X_1X_2\cdots X_n]=E[X_1]^n$$

8 Statistical Inference 1: Classical methods

• **Statistical Inference** is a collection of methods that deal with drawing conclusions from data that are prone to random variation

8.1 Random sampling

- The collection of random variables $X_1, X_2, X_3, \dots, X_n$ is said to be a random sample of size n if they are independent and identically distributed i.e.
 - 1. $X_1, X_2, X_3, \dots, X_n$ are independent random variables and
 - 2. they have the same distribution i.e.

$$F_{X_1}(x) = F_{X_2}(x) = ... = F_{X_n}(x),$$
 for all $x \in \mathbb{R}$.

• The **point estimator** is a function of the random sample $\hat{\Theta} = h(X_1, X_2, \dots, X_n)$ that is used to estimate an unknown quantity

- Properties of random sampling
 - 1. the X_i 's are independent

2.
$$F_{X_1}(x) = F_{X_2}(x) = \dots = F_{X_n}(x) = F_X(x)$$

3.
$$EX_i = EX = \mu < \infty$$

4.
$$0 < \operatorname{Var}(X_i) = \operatorname{Var}(X) = \sigma^2 < \infty$$

• The **sample mean** is defined as

$$\overline{X} = \frac{X_1 + X_2 + \dots + X_n}{n}.$$

- Properties of the sample mean
 - 1. $\overline{EX} = \mu$
 - 2. $Var(\overline{X}) = \frac{\sigma^2}{n}$
 - 3. Weak Law of Large Numbers (WLLN):

$$\lim_{n\to\infty} P(|\overline{X}-\mu| \ge \epsilon) = 0.$$

4. Central limit Theorem: The random variable

$$Z_n = \frac{\overline{X} - \mu}{\sigma / \sqrt{n}} = \frac{X_1 + X_2 + \dots + X_n - n\mu}{\sqrt{n}\sigma}$$

converges in distribution to the standard normal random variable as n goes to infinity that is

$$\lim_{n\to\infty} P(Z_n \le x) = \Phi(x), \quad \text{for all } x \in \mathbb{R}$$

where $\Phi(x)$ is the standard normal CDF

• If we let $X_1, X_2, ... X_n$ be a random sample from a continuous distribution function with CDF $F_X(x)$ then if we order X_i 's from the smallest to the largest and denote the resulting sequence of random variables as

$$X_{(1)}, X_{(2)}, \cdots, X_{(n)}.$$

- $X_{(1)}, X_{(2)}, \cdots, X_{(n)}$ is called the **order statistics** of the random sample $X_1, X_2, \ldots X_n$
- Theorem 8.1: Let $X_1, X_2, ... X_n$ be a random sample from a continuous distribution with CDF $F_X(x)$ and PDF $f_X(x)$. Let $X_{(1)}, X_{(2)}, ..., X_{(n)}$ be the order statistics of the random sample $X_1, X_2, ... X_n$. Then the CDF and PDF of $X_{(i)}$ are given by

$$f_{X_{(i)}}(x) = \frac{n!}{(i-1)!(n-i)!} f_X(x) \left[F_X(x) \right]^{i-1} \left[1 - F_X(x) \right]^{n-i}, \tag{32}$$

$$F_{X_{(i)}}(x) = \sum_{k=i}^{n} {n \choose k} \left[F_X(x) \right]^k \left[1 - F_X(x) \right]^{n-k}. \tag{33}$$

Also the join PDF of $X_{(1)}, X_{(2)}, \dots, X_{(n)}$ is given by

$$f_{X_{(1)},\dots,X_{(n)}}(x_1,x_2,\dots,x_n) = \begin{cases} n! f_X(x_1) f_X(x_2) \dots f_X(x_n) & \text{for } x_1 \le x_2 \le x_2 \dots \le x_n \\ 0 & \text{otherwise} \end{cases}$$

8.2 Point Estimation

• If θ is the unknown parameter to be estimated, where θ is a fixed (non-random) quantity, we estimate θ by defining a point estimator $\hat{\Theta}$ that is a function of the random sample i.e.

$$\hat{\Theta} = h(X_1, X_2, \cdots, X_n).$$

• If $\theta = EX$ we may choose the $\hat{\Theta}$ to be the sample mean

$$\hat{\Theta} = \overline{X} = \frac{X_1 + X_2 + \dots + X_n}{n}$$

8.2.1 Evaluating estimators

• Let $\hat{\Theta} = h(X_1, X_2, \dots, X_n)$ be a point estimator for θ . The **bias** of the point estimator $\hat{\Theta}$ is defined by

$$B(\hat{\Theta}) = E[\hat{\Theta}] - \theta.$$

- In general we want to have a bias close to 0
- Let $\hat{\Theta} = h(X_1, X_2, \dots, X_n)$ be a point estimator for a parameter θ . We say that $\hat{\Theta}$ is an **unbiased** of estimator of θ if

$$B(\hat{\Theta}) = 0$$
, for all possible values of θ .

• The **mean squared error** (MSE) of a point estimator $\hat{\theta}$ is defined as

$$MSE(\hat{\Theta}) = E[(\hat{\Theta} - \theta)^2].$$

- The MSE is a measure of the distance between $\hat{\Theta}$ and θ
- A smaller MSE is generally indicative a better estimator
- The MSE can generally be written as

$$MSE(\hat{\Theta}) = Var(\hat{\Theta}) + B(\hat{\Theta})^2.$$

- We say that an estimator is **consistent** if as the sample size n get larger $\hat{\Theta}$ converges to the real value of θ
- Let $\hat{\Theta}_{-1}$, $\hat{\Theta}_{-2}$, ... $\hat{\Theta}_{-n}$..., be a sequence of point estimators for θ . We say that $\hat{\Theta}_n$ is a **consistent** estimator of θ if

$$\lim_{n\to\infty} P(|\hat{\Theta}_n - \theta| \ge \epsilon) = 0, \text{ for all } \epsilon > 0.$$

Theorem 8.2

• Let $\hat{\Theta}_{-1}$, $\hat{\Theta}_{-2}$, ... be a sequence of point estimators of θ if

$$\lim_{n\to\infty} MSE(\hat{\Theta}_n) = 0,$$

then $\hat{\Theta}_n$ is a consistent estimator of θ

8.2.2 For mean and variance

• Let $X_1, X_2, X_3..., X_n$ be a random sample with mean $EX_i = \mu < \infty$ and variance $0 < Var(X_i) = \sigma^2 < \infty$. Then the **sample variance** of this random sample is defined as

$$S^{2} = \frac{1}{n-1} \sum_{k=1}^{n} (X_{k} - \overline{X})^{2} = \frac{1}{n-1} \left(\sum_{k=1}^{n} X_{k}^{2} - n \overline{X}^{2} \right).$$

• The sample variance is an unbiased estimator of σ^2 . The **sample standard deviation** is defined as

$$S = \sqrt{S^2}$$

and is commonly used as an estimator for σ . Nevertheless, S is a biased estimator for σ

8.2.3 Maximum likelihood estimation

- Let $X_1, X_2, ..., X_n$ be a random sample from a distribution with a parameter θ . Suppose that we have observed $X_1 = x_1, X_2 = x_2, ..., X_n = x_n$
 - 1. If X_i 's are discrete, then the **likelihood function** is defined as

$$L(x_1,x_2,\cdots,x_n;\theta)=P_{X_1X_2\cdots X_n}(x_1,x_2,\cdots,x_n;\theta).$$

2. If X_i 's are jointly continuous, then the likelihood function is defined as

$$L(x_1,x_2,\cdots,x_n;\theta)=f_{X_1X_2\cdots X_n}(x_1,x_2,\cdots,x_n;\theta).$$

• In some problems it is easier to work with the log likelihood function given by

$$\ln L(x_1, x_2, \cdots, x_n; \theta).$$

• Let $X_1, X_2, ..., X_n$ be a random sample from a distribution with a parameter θ . Given that we have observed $X_1 = x_1, X_2 = x_2, ..., X_n = x_n$ a maximum likelihood estimate of θ , shown by $\hat{\theta}_{ML}$ is the value of θ that maximizes the likelihood function

$$L(x_1, x_2, \cdots, x_n; \theta)$$

- A maximum estimator (MLE) of a parameter θ shown by $\hat{\Theta}_{ML}$ is a random variable
 - $\hat{\Theta}_{ML}(X_1, X_2, \dots, X_n)$ whose value when X_1, X_2, \dots, X_n is given by $\hat{\theta}_{ML}$

- Asymptotic properties of MLEs: Let $X_1, X_2, ..., X_n$ be a random sample from a distribution with a parameter θ . Let $\hat{\Theta}_{ML}$ denote the maximum likelihood estimator (MLE) of θ . Then under some mild regularity conditions:
 - 1. $\hat{\Theta}_{ML}$ is asymptotically consistent i.e.

$$\lim_{n\to\infty} P(|\hat{\Theta}_{ML} - \theta| > \epsilon) = 0$$

2. $\hat{\Theta}_{ML}$ is asymptotically unbiased i.e.

$$\lim_{n\to\infty} E[\hat{\Theta}_{ML}] = \theta.$$

3. As *n* becomes large $\hat{\Theta}_{ML}$ is approximately a random variable. More precisely the random variable

$$\frac{\hat{\Theta}_{ML} - \theta}{\sqrt{\operatorname{Var}(\hat{\Theta}_{ML})}}$$

converges in distribution to N(0,1)

8.3 Interval estimation

- Let $X_1, X_2, ... X_n$ be a random sample from a distribution with parameter θ that is to be estimated, the goal is to find two estimators for θ
 - 1. the low estimator $\hat{\Theta}_l = \hat{\Theta}_l(X_1, X_2, \dots, X_n)$
 - 2. the high estimator $\hat{\Theta}_h = \hat{\Theta}_h(X_1, X_2, \cdots, X_n)$
- The interval estimator is given by the interval $[\hat{\Theta}_l, \hat{\Theta}_h]$
- Let $X_1, X_2, ... X_n$ be a random sample from a distribution with parameter θ that is to be estimated. An **interval estimator** with **confidence level** 1α consists of two estimators $\hat{\Theta}_l(X_1, X_2, ..., X_n)$ and $\hat{\Theta}_h(X_1, X_2, ..., X_n)$ such that

$$P\Big(\hat{\Theta}_l \leq \theta \text{ and } \hat{\Theta}_h \geq \theta\Big) \geq 1 - \alpha,$$

for every possible value of θ . Equivalently, we say that $[\hat{\Theta}_l, \hat{\Theta}_h]$ is a $(1 - \alpha)100\%$ **confidence interval** for θ

8.3.1 Find interval estimators

Pivotal Quantity

Let $X_1, X_2, ..., X_n$ be an random sample from a distribution with parameter θ that is to be estimated. The random variable Q is said to be a pivot or a pivotal quantity, if it has the following properties

1. It is a function of the observed data X_1, X_2, \dots, X_n and the unknown parameter θ , but it does not depend on any other unknown parameters:

$$Q=Q(X_1,X_2,\cdots,X_n,\theta).$$

2. The probability distribution of Q does not depend on θ or any other unknown parameters

Interval estimators:

- Assumptions: A random sample $X_1, X_2, ..., X_n$ is given from a distribution with known variance $Var(X_i) = \sigma^2 < \infty$; n is large
- Parameter to be Estimated: $\theta = EX_i$
- Confidence Interval: $\left[\overline{X} z_{\frac{\alpha}{2}} \frac{\sigma}{\sqrt{n}}, \overline{X} + z_{\frac{\alpha}{2}} \frac{\sigma}{\sqrt{n}}\right]$ is approximately a $(1 \alpha)100\%$ confidence interval for θ
- Assumptions: A random sample $X_1, X_2, ..., X_n$ is given from a $Bernoulli(\theta)$; n is large
- Parameter to be Estimated: θ
- Confidence Interval: $\left[\overline{X} \frac{z_{\frac{\alpha}{2}}}{2\sqrt{n}}, \overline{X} + \frac{z_{\frac{\alpha}{2}}}{2\sqrt{n}}\right]$ is approximately $(1 \alpha)100\%$ confidence interval for θ . This is a conservative confidence interval as it is obtained using an upper bound for σ
- Assumptions: A random sample X_1, X_2, \ldots, X_n is given from a *Bernoulli*(θ); n is large
- Parameter to be Estimated: θ
- Confidence Interval: $\left[\overline{X} z_{\frac{\alpha}{2}} \sqrt{\frac{\overline{X}(1-\overline{X})}{n}}, \overline{X} + z_{\frac{\alpha}{2}} \sqrt{\frac{\overline{X}(1-\overline{X})}{n}}\right]$ is approximately a $(1-\alpha)100\%$ confidence interval for θ
- Assumptions: A random sample $X_1, X_2, ..., X_n$ is given from a distribution with unknown variance $Var(X_i) = \sigma^2 < \infty$; n is large
- Parameter to be Estimated: $\theta = EX_i$
- Confidence Interval: If *S* is the sample standard deviation

$$S = \sqrt{\frac{1}{n-1} \sum_{k=1}^{n} (X_k - \overline{X})^2} = \sqrt{\frac{1}{n-1} \left(\sum_{k=1}^{n} X_k^2 - n \overline{X}^2 \right)}$$

then the interval

$$\left[\overline{X}-z_{rac{lpha}{2}}rac{S}{\sqrt{n}},\overline{X}+z_{rac{lpha}{2}}rac{S}{\sqrt{n}}
ight]$$

is approximately a $(1-\alpha)100\%$ confidence interval for θ

8.4 Confidence intervals for normal samples

8.4.1 The Chi-Squared Distribution

• **Definition 8.1:** If $Z_1, Z_2, ..., Z_n$ are independent standard normal random variables, the random variable Y defined as

$$Y = Z_1^2 + Z_2^2 + \dots + Z_n^2$$

is said to have a chi-squared distribution with n degrees of freedom shown by

$$Y \sim \chi^2(n)$$

- Properties of the chi-squared distribution:
 - The chi-squared distribution in a special case of the gamma distribution. More specifically

$$Y \sim Gamma\left(\frac{n}{2}, \frac{1}{2}\right).$$

Thus

$$f_Y(y) = \frac{1}{2^{\frac{n}{2}}\Gamma(\frac{n}{2})}y^{\frac{n}{2}-1}e^{-\frac{y}{2}}, \text{ for } y > 0.$$

- 2. EY = n, Var(Y) = 2n
- 3. For any $P \in [0,1]$ and $n \in \mathbb{N}$, we define $\chi^2_{p,n}$ as the real value for which

$$P(Y > \chi_{p,n}^2) = p,$$

where $U \sim \chi^2(n)$

• **Theorem 8.3:** Let $X_1, X_2, ..., X_n$ be i.i.d. $N(\mu, \sigma)$ random variables. Also let S^2 be the standard variance for this random sample. Then the random variable Y is defined as

$$Y = \frac{(n-1)S^2}{\sigma^2} = \frac{1}{\sigma^2} \sum_{i=1}^{n} (X_i - \overline{X})^2$$

has a chi-squared distribution with n-1 degrees of freedom, i.e., $Y \sim \chi^2(n-1)$. Moreover, \hat{X} and S^2 are independent random variables.

8.4.2 The t Distribution

• **Definition 8.2:** Let $Z \sim N(0,1)$, and $Y \sim \chi^2(n)$, where $n \in \mathbb{N}$. Also assume that Z and Y are independent. The random variable T defined as

$$T = \frac{Z}{\sqrt{Y/n}}$$

is said to have a *t*-distribution with *n* degrees of freedom shown by

$$T \sim T(n)$$

• Properties:

- 1. The *t*-distribution has a bell-shaped PDF centered at 0, but its PDF is more spread out than the normal PDF
- 2. ET = 0, for n > 0. But ET, is undefined for n = 1.
- 3. $Var(T) = \frac{n}{n-2}$ for n < 2. But variance is undefined for n = 1, 2
- 4. As *n* becomes large, the *t* density approaches the standard normal PDF. More formally, we can write

$$T(n) \xrightarrow{d} N(0,1).$$

5. For any $p \in [0,1]$ and $n \in$, we define $t_{p,n}$ as the real value for which

$$P(T > t_{p,n}) = p.$$

• Since the *t*-distribution has a symmetric PDF, we have

$$t_{1-p,n}=-t_{p,n}.$$

• **Theorem 8.4:** Let $X_1, X_2, ..., X_n$ be i.i.d. $N(\mu, \sigma)$ random variables. Also let S^2 be the standard variance for this random sample. Then let the random variable T defined as

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

has a *t*-distribution with n-1 degrees of freedom, i.e. $T \sim T(n-1)$

8.4.3 More Interval estimators

- Assumptions: A random sample X_1, X_2, \dots, X_n is given from a $N(\mu, \sigma^2)$ distribution where $Var(X_i) = \sigma^2$ is known
- Parameter to be estimated: $\mu = EX_i$
- Confidence interval $\left[\overline{X} z_{\frac{\alpha}{2}} \frac{\sigma}{\sqrt{n}}, \overline{X} + z_{\frac{\alpha}{2}} \frac{\sigma}{\sqrt{n}}\right]$ is a $(1 \alpha)100\%$ confidence interval for mu
- Assumptions: A random sample $X_1, X_2, ..., X_n$ is given from a $N(\mu, \sigma^2)$ distribution where $Var(X_i) = \sigma^2$ are unknown
- Parameter to be estimated: $\mu = EX_i$
- Confidence interval $\left[\overline{X} t_{\frac{\alpha}{2},n-1} \frac{S}{\sqrt{n}}, \overline{X} + t_{\frac{\alpha}{2},n-1} \frac{S}{\sqrt{n}}\right]$ is a $(1-\alpha)100\%$ confidence interval for mu
- Assumptions: A random sample $X_1, X_2, ..., X_n$ is given from a $N(\mu, \sigma^2)$ distribution where $\mu = EX_i$ and $Var(X_i) = \sigma^2$ are unknown
- Parameter to be estimated: $Var(X_i) = \sigma^2$
- Confidence interval $\left[\frac{(n-1)S^2}{\chi^2_{\frac{\alpha}{2},n-1}},\frac{(n-1)S^2}{\chi^2_{1-\frac{\alpha}{2},n-1}}\right]$ is a $(1-\alpha)100\%$ confidence interval for mu

8.5 Hypothesis testing

- Let *S* be the set of possible values for θ , we can partition *S* into two disjoint sets S_0 and S_1 . Then let H_0 be the hypothesis that $\theta \in S_0$ and let H_1 be the hypothesis that $\theta \in S_1$
 - − H_0 (the **null** hypothesis): $\theta \in S_0$
 - **–** H_1 (the **alternative** hypothesis): θ ∈ S_1
- **Definition 8.3:** Let $X_1, ..., X_n$ be a random sample of interest. A **statistic** is a real valued function of data. For example the sample mean defined as

$$W(X_1, X_2, \cdots, X_n) = \frac{X_1 + X_2 + \dots + X_n}{n},$$

is a statistic. A test statistic is a statistic based on which we build our test

- To decide whether to choose H_0 or H_1 we
 - choose a test statistic $W(X_1, X_2, \dots, X_n)$
 - define a set $A \subset \mathbb{R}$ as the possible values of W for which we would accept H_0 called the **acceptance region**
 - the set $R = \mathbb{R} A$ is said to be the **rejection region**
- **Type 1 error** is defined as the event that we reject H_0 when H_0 is true.
 - The probability of a type 1 error is

$$P(\text{type I error} \mid \theta) = P(\text{Reject } H_0 \mid \theta)$$
(34)

$$= P(W \in R \mid \theta), \quad \text{for } \theta \in S_0. \tag{35}$$

If the probability of a type 1 error satisfies

$$P(\text{type I error}) \leq \alpha$$
, for all $\theta \in S_0$,

then we say that the test has **significance level** α or simply the test is a α test

- **Type 2 error** is defined as the event that we accept H_0 when H_0 is false.
 - The probability is a function of θ and is shown by β :

$$\beta(\theta) = P(\text{Accept } H_0 \mid \theta), \quad \text{for } \theta \in S_1.$$

8.5.1 Hypothesis test for mean

• To decide between the following hypothesis

$$H_0: \mu = \mu_0$$

$$H_1: \mu \neq \mu_0$$

the null hypothesis is a simple hypothesis and the alternative is a two-sided, this hypothesis test is called a two sided hypothesis test

Table 8.2: Two-sided hypothesis testing for the mean: $H_0: \mu = \mu_0, H_1: \mu \neq \mu_0$.

Case	Test Statistic	Acceptance Region
$X_i \sim N(\mu, \sigma^2), \sigma { m known}$	$W = \frac{\overline{X} - \mu_0}{\sigma / \sqrt{n}}$	$ W \le z_{\frac{\alpha}{2}}$
n large, X_i non-normal	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$ W \le z_{\frac{\alpha}{2}}$
$X_i \sim N(\mu, \sigma^2)$, σ unknown	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$ W \le t_{\frac{a}{2}, n-1}$

• If the hypothesis is defined as follows

$$\mu \leq \mu_0 \vee \mu \geq mu_0$$

$$\mu > \mu_0 \lor \mu < mu_0$$

the null hypothesis and the alternative are one-sided this is called a one-sided hypothesis test:

Table 8.3: One-sided hypothesis testing for the mean: $H_0: \mu \leq \mu_0$, $H_1: \mu > \mu_0$.

Case	Test Statistic	Acceptance Region
$X_i \sim N(\mu, \sigma^2), \sigma { m known}$	$W = \frac{\overline{X} - \mu_0}{\sigma / \sqrt{n}}$	$W \leq z_{\alpha}$
n large, X_i non-normal	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$W \le z_{\alpha}$
$X_i \sim N(\mu, \sigma^2), \sigma { m unknown}$	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$W \le t_{\alpha,n-1}$

Table 8.4. One-sided hypothesis testing for the mean: $H_0: \mu \geq \mu_0, H_1: \mu < \mu$	8.4: One-sided hypothesis testing for the mean: H	$I_0: \mu >$	$\mu_0, H_1: \mu < \mu$
---	---	--------------	-------------------------

Case	Test Statistic	Acceptance Region
$X_i \sim N(\mu, \sigma^2)$, σ known	$W = \frac{\overline{X} - \mu_0}{\sigma / \sqrt{n}}$	$W \ge -z_{\alpha}$
n large, X_i non-normal	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$W \ge -z_{\alpha}$
$X_i \sim N(\mu, \sigma^2)$, σ unknown	$W = \frac{\overline{X} - \mu_0}{S / \sqrt{n}}$	$W \ge -t_{\alpha,n-1}$

8.5.2 P-values

- **P-value** is the lowest significance level α that results in rejecting the null hypothesis
- Consider a hypothesis test for choosing between H₀ and H₁. Let W be the test statistic, and w₁ be the observed value of W
 - 1. Assume H_0 is true
 - 2. The P-value is P(type I error) then the test threshold c is chosen to be $c = w_1$

8.5.3 Likelihood ratio test

• Let $X_1, ..., X_n$ be a random sample from a distribution with parameter θ . Suppose we have observed $X_1 = x_1, ..., X_n = x_n$. To decide between two simple hypothesis

$$H_0: \theta = \theta_0$$

$$H_1: \theta = \theta_1$$

we define

$$\lambda(x_1, x_2, \cdots, x_n) = \frac{L(x_1, x_2, \cdots, x_n; \theta_0)}{L(x_1, x_2, \cdots, x_n; \theta_1)}.$$

To perform a **likelihood ratio test (LRT)**, we choose a constant c. We reject H_0 if $\lambda < c$ and accept it if $\lambda > c$. The value of c can be chosen based on the desired a

• Let $X_1, ..., X_n$ be a random sample from a distribution with parameter θ . Suppose we have observed $X_1 = x_1, ..., X_n = x_n$. Define

$$\lambda(x_1,x_2,\cdots,x_n)=\frac{\sup\{L(x_1,x_2,\cdots,x_n;\theta):\theta\in S_0\}}{\sup\{L(x_1,x_2,\cdots,x_n;\theta):\theta\in S\}}.$$

To perform a **likelihood ration test (LRT)**, we choose a constant c in [0,1]. We reject H_0 if $\lambda < c$ and accept it if $\lambda > c$. The value of c can be chosen based on the desired a

9 Discrete-Time Markov Chains

• Consider the random process $\{X_n, n=1,2,\cdots\}$, where $R_{X_i}=S\subset\{0,1,2,\cdots\}$. We say this process is a Markov chain if

$$P(X_{m+1} = j | X_m = i, X_{m-1} = i_{m-1}, \dots, X_0 = i_0) = P(X_{m+1} = j | X_m = i),$$

for all $m, j, i, i_0, i_1, \dots, i_{m-1}$. If the number of states is finite. e.g. $S = \{0, 1, 2, \dots, r\}$ we call it a **finite** Markov chain

- If $X_n = j$, we say that the process is in state j
- The numbers $P(X_{m+1} = j | X_m = i)$ are called the **transition probabilities**
 - It is assumed that they do not depend on time. That is $P(X_{m+1}=j|X_m=i \text{ do not depend on } m$
- The following is defined

$$p_{ij} = P(X_{m+1} = j | X_m = i).$$

we have in particular

$$p_{ij} = P(X_1 = j | X_0 = i) (36)$$

$$= P(X_2 = j | X_1 = i) (37)$$

$$= P(X_3 = j | X_2 = i) = \cdots.$$
 (38)

if the process is in state i, it will a transition to state j with probability p_{ij} .

9.1 State Transition Matrix and Diagram

- The transitions probabilities is often listed in a matrix
 - is called the state transition matrix or transition probability matrix and is usually shown by P
 - Assuming the states are $1, 2, \dots, r$, then the state transition matrix is shown by

$$P = \begin{bmatrix} p_{11} & p_{12} & \dots & p_{1r} \\ p_{21} & p_{22} & \dots & p_{2r} \\ \vdots & \vdots & \vdots & \vdots \\ p_{r1} & p_{r2} & \dots & p_{rr} \end{bmatrix}.$$

for $p_{ij} \ge 0$, and for all i we have

$$\sum_{k=1}^{r} p_{ik} = \sum_{k=1}^{r} P(X_{m+1} = k | X_m = i) = 1.$$

- A Markov chain is usually shown by a state transition diagram
 - If there is no arrow from state *i* to state *j*, then $p_{ij} = 0$

9.2 Probability distributions

• Given a Markov chain $\{X_n, n = 0, 1, 2, ...\}$ where $X_n \in S = \{1, 2, \cdots r\}$. The following is defined

$$\pi^{(n)} = \begin{bmatrix} P(X_n = 1) & P(X_n = 2) & \cdots & P(X_n = r) \end{bmatrix}$$
,

• The following is then true

$$\pi^{(n+1)} = \pi^{(n)} P$$
, for $n = 0, 1, 2, \cdots$;
 $\pi^{(n)} = \pi^{(0)} P^n$, for $n = 0, 1, 2, \cdots$.

• Given a Markov chain $\{X_n, n = 0, 1, 2, ...\}$ the *n*-step probabilities $p_{ij}^{(n)}$ is defined as

$$p_{ij}^{(n)} = P(X_n = j | X_0 = i), \text{ for } n = 0, 1, 2, \cdots,$$

and the n-step transition matrix, $P^{(n)}$, as

$$P^{(n)} = \begin{bmatrix} p_{11}^{(n)} & p_{12}^{(n)} & \dots & p_{1r}^{(n)} \\ p_{21}^{(n)} & p_{22}^{(n)} & \dots & p_{2r}^{(n)} \\ \vdots & \vdots & \vdots & \vdots \\ p_{r1}^{(n)} & p_{r2}^{(n)} & \dots & p_{rr}^{(n)} \end{bmatrix}.$$

• The Chapman-Kolmogorov equation can be written as

$$p_{ij}^{(m+n)} = P(X_{m+n} = j | X_0 = i)$$
$$= \sum_{k \in S} p_{ik}^{(m)} p_{kj}^{(n)}.$$

• The *n* step matrix is given by

$$P^{(n)} = P^n$$
, for $n = 1, 2, 3, \cdots$.

9.3 Classification of States

- A state j is **accessible** from state i, written as $i \to j$, if $p_{ij}^{(n)} > 0$ for some n. It is assumed that every state is accessible from itself since $p_{ii}^{(0)} = 1$
- Two state *i* and *j* are said to **communicate**, written as *i* ↔ *j* if they are accessible from each other, in other words

$$i \leftrightarrow j$$
 means $i \rightarrow j$ and $j \rightarrow i$.

- Communication is an equivalence relation, which means
 - every state communicates with itself $i \leftrightarrow i$;
 - if $i \leftrightarrow j$ then $j \leftrightarrow i$;

- if $i \leftrightarrow j$ and $j \leftrightarrow k$ then $i \leftrightarrow k$
- Markov chains can be partitioned into communicating classes such that only member of the same class communicate with each other
- A Markov chain is said to be irreducible if all the states communicate with each other
- For any state *i* we define

$$f_{ii} = P(X_n = i, \text{ for some } n \ge 1 | X_0 = i).$$

State *i* is **recurrent** if $f_{ii} = 1$, and is **transient** if $f_{ii} < 1$

- If two states are in the same class they are either both recurrent or transient
 - Therefore we can say that a class is transient if all the states in it are transient and the same is the case for recurrent
- Consider a discrete-time Markov chain. Let V be the total number of visits to state i
 - 1. If i is a recurrent state, then

$$P(V = \infty | X_0 = i) = 1.$$

2. If i is a transient state, then

$$V|X_0 = i \sim Geometric(1 - f_{ii}).$$

- The **period** of a state i is the largest integer d satisfying the following property: $p_{ii}^{(n)} = 0$ whenever n is not divisible by d. The period of i is shown as d(i). If $p_{ii}^{(n)} = 0$ for all n > 0, we let $d(i) = \infty$
 - If d(i) < 1, we say that state i is **periodic**
 - If d(i) = 1, we say that state i is aperiodic
- The states in the same communicating class have the same period
 - a class therefor can be periodic or aperiodic

If
$$i \leftrightarrow j$$
, then $d(i) = d(j)$

- Consider a finite irreducible Markov chain X_n
 - 1. If there is a self-transition in the chain ($p_{ii} > 0$ for some i), then the chain is aperiodic.
 - 2. Suppose that you can go from state i to state i in l steps i.e. $p_{ii}^{(l)} > 0$. Also suppose that $p_{ii}^{(m)} > 0$ for some m. If gcd(l, m) = 1, then state i is aperiodic.
 - 3. The chain is aperiodic if and only if there exists a positive integer n such that all the elements of the matrix P^n are strictly positive, i.e.

$$p_{ij}^{(n)} > 0$$
, for all $i, j \in S$.

9.4 Using the Law of Total Probability with Recursion

9.4.1 Absorption Probabilities

• Consider a finite Markov chain $\{X_n, n = 0, 1, 2, \cdots\}$ with state space $S = \{0, 1, 2, \cdots r\}$. Suppose that all states are either absorbing or transient. Let $l \in S$ be an absorbing state. Define

$$a_i = P(\text{absorption in } l | X_0 = i), \text{ for all } i \in S.$$

By the above definition, we have $a_l = 1$, and $a_j = 0$ if j is another absorbing state. To find unknown values of a_i 's we can use the following equations

$$a_i = \sum_k a_k p_{ik}$$
, for $i \in S$.

9.4.2 Mean Hitting Times

• Consider a finite Markov chain $\{X_n, n = 0, 1, 2, \dots\}$ with state space $S = \{0, 1, 2, \dots r\}$. Let $A \subset S$ be a set of states. Let T be the first time the chain visits a state in A. For all $i \in S$ define

$$t_i = E[T|X_0 = i].$$

• By the above definition we have, we have $t_j = 0$, for all $j \in A$. To find the unknown values of t_i 's we can use the following equations

$$t_i = 1 + \sum_k t_k p_{ik}$$
, for $i \in S - A$.

9.4.3 Mean Return Times

• Consider a finite Markov chain $\{X_n, n = 0, 1, 2, \dots\}$ with state space $S = \{0, 1, 2, \dots r\}$. Let $l \in S$ be a state. Let r_l be the **mean return time** to state l. Then

$$r_l = 1 + \sum_k t_k p_{lk},$$

where t_k is the expected time until the chain hits state l given $X_0 = k$. Specifically

$$t_l = 0,$$

 $t_k = 1 + \sum_j t_j p_{kj}, \quad \text{for } k \neq l.$

9.5 Limiting and stationary distributions

• The probability distribution $\pi = [\pi_0, \pi_1, \pi_2, \cdots]$ is called the **limiting distribution** of the Markov chain X_n if

$$\pi_j = \lim_{n \to \infty} P(X_n = j | X_0 = i)$$

for all $i, j \in S$ we have

$$\sum_{j\in S}\pi_j=1.$$

- Consider a finite Markov chain $\{X_n, n = 0, 1, 2, \dots\}$ where $X_n \in S = \{1, 2, \dots, r\}$. Assume that the chain is irreducible and aperiodic. Then
- 1. The set of equations

$$\pi = \pi P,$$

$$\sum_{j \in S} \pi_j = 1$$

has a unique solution

2. The unique solution to the above equations is the limiting distribution of the Markov chain i.e.

$$\pi_j = \lim_{n \to \infty} P(X_n = j | X_0 = i),$$

for all $i, j \in S$

3. We have

$$r_j = \frac{1}{\pi_j}$$
, for all $j \in S$,

where r_i is the mean return time to state j

9.6 Countably Infinite Markov Chains

• Let i be a recurrent state. Assuming $A_0 = i$, let R_i be the number for transitions needed to return to i, i.e.

$$R_i = \min\{n \geq 1 : X_n = i\}.$$

If $r_i = E[R_i|X_0 = i] < \infty$, then i is said to be **positive recurrent**. If $E[R_i|X_0 = i] = \infty$, then i is said to be **null recurrent**

Theorem 11.2

- Consider an infinite Markov chain $\{X_n, n = 0, 1, 2, ...\}$ where $X_n \in S = \{0, 1, 2, ...\}$. Assume that the chain is irreducible and and aperiodic. The one of the following cases can occur:
 - 1. All states are transient and

$$\lim_{n\to\infty} P(X_n = j|X_0 = i) = 0, \text{ for all } i, j.$$

2. All states are null recurrent

$$\lim_{n \to \infty} P(X_n = j | X_0 = i) = 0$$
, for all *i*, *j*.

3. All states are positive recurrent. In this case, there exists a limiting distribution, $\pi = [\pi_0, \pi_1, \cdots]$, where

$$\pi_j = \lim_{n \to \infty} P(X_n = j | X_0 = i) > 0,$$

for all $i, j \in S$. The limiting distribution in unique solution to the equations

$$\pi_j = \sum_{k=0}^{\infty} \pi_k P_{kj}, \quad \text{for } j = 0, 1, 2, \cdots,$$

$$\sum_{j=0}^{\infty} \pi_j = 1.$$

We also have

$$r_j = \frac{1}{\pi_j}$$
, for all $j = 0, 1, 2, \cdots$,

where r_j is the mean return time to state j