# **Appendix B**

# **Overview in Probability and Random Processes**

**Po-Ning Chen, Professor**

**Institute of Communications Engineering**

**National Chiao Tung University**

**Hsin Chu, Taiwan 30010, R.O.C.**

**Definition B.1** ( $\sigma$ -**Fields)** Let F be a collection of subsets of a non-empty set Ω. Then F is called <sup>a</sup> <sup>σ</sup>*-field* (or <sup>σ</sup>*-algebra*) if the following hold:

1.  $\Omega \in \mathcal{F}$ .

2. F is closed under complementation: If  $A \in \mathcal{F}$ , then  $A^c \in \mathcal{F}$ , where  $A^c =$  $\{\omega\in\Omega\colon\omega\not\in A\}.$ 

3. F is closed under countable unions: If  $A_i \in \mathcal{F}$  for  $i = 1, 2, 3, \ldots$ , then  $\bigcup_{i=1}^{\infty} A_i \in \mathcal{F}.$ 

• It directly follows that the empty set  $\emptyset$  is also an element of  $\mathcal F$  (as  $\Omega^c = \emptyset$ ) and that  $\mathcal F$  is closed under countable intersection since

$$
\bigcap_{i=1}^{\infty} A_i^c = \left(\bigcup_{i=1}^{\infty} A_i\right)^c.
$$

- The largest  $\sigma$ -field of subsets of a given set  $\Omega$  is the collection of all subsets of  $\Omega$  (i.e., its powerset), while the smallest  $\sigma$ -field is given by  $\{\Omega, \emptyset\}$ .
- Also, if A is a proper (strict) non-empty subset of  $\Omega$ , then the smallest  $\sigma$ -field containing A is given by  $\{\Omega, \emptyset, A, A^c\}.$

**Definition B.2 (Probability space)** A *probability space* is <sup>a</sup> triple  $(\Omega, \mathcal{F}, P)$ , where  $\Omega$  is a given set called *sample space* containing all possible outcomes (usually observed from an experiment),  $\mathcal F$  is a  $\sigma$ -field of subsets of  $Ω$ , and *P* is a probability measure  $P: F → [0, 1]$  on the σ-field satisfying the following:

- 1.  $0 \leq P(A) \leq 1$  for all  $A \in \mathcal{F}$ .
- 2.  $P(\Omega) = 1$ .
- 3. *Countable additivity:* If  $A_1, A_2, \ldots$  is a sequence of disjoint sets (i.e.,  $A_i \cap$  $A_j = \emptyset$  for all  $i \neq j$  in F, then

$$
P\left(\bigcup_{k=1}^{\infty} A_k\right) = \sum_{k=1}^{\infty} P(A_k).
$$

- It directly follows from Properties 1-3 of the above definition that  $P(\emptyset) = 0$ .
- Usually, the  $\sigma$ -field  $\mathcal F$  is called the *event space* and its elements (which are subsets of Ω satisfying the properties of Definition B.1) are called *events*.

- The Borel  $\sigma$ -field of R, denoted by  $\mathscr{B}(\mathbb{R})$ , is the smallest  $\sigma$ -field of subsets of R containing all open intervals in R.
- The elements of  $\mathscr{B}(\mathbb{R})$  are called Borel sets.
- For any random variable X, we use  $P_X$  to denote the probability distribution on  $\mathscr{B}(\mathbb{R})$  induced by X, given by

 $P_X(B) := \Pr[X \in B] = P(w \in \Omega : X(w) \in B), \qquad B \in \mathscr{B}(\mathbb{R}).$ 

Note that the quantities  $P_X(B)$ ,  $B \in \mathscr{B}(\mathbb{R})$ , fully characterize the random variable  $X$  as they determine the probabilities of all events that concern  $X$ .

## B.2 Random variables and random processes I: b-5

• A random variable X defined over probability space  $(\Omega, \mathcal{F}, P)$  is a real-valued function  $X: \Omega \to \mathbb{R}$  that is *measurable* (or F-measurable), i.e., satisfying the property that

$$
X^{-1}((-\infty, t]) := \{\omega \in \Omega : X(\omega) \le t\} \in \mathcal{F}
$$

for each real  $t$ .

• A random process (or random source) is <sup>a</sup> collection of random variables that arise from the same probability space. It can be mathematically represented by the collection

 $\{X_t, t \in I\},\$ 

where  $X_t$  denotes the  $t<sup>th</sup>$  random variable in the process, and the index t runs over an index set  $I$  which is arbitrary.

# B.2 Random variables and random processes I: b-6

- The index set I can be uncountably infinite (e.g.,  $I = \mathbb{R}$ ), in which case we are dealing with <sup>a</sup> continuous-time process.
- Except for <sup>a</sup> brief interlude with the continuous-time (waveform) Gaussian channel in Chapter 5, we will consider discrete-time communication systems throughout the lectures.

To be precise, we will only consider the following cases of index set I: *case a)* I consists of one index only. *case b)* I is finite. *case c)* I is countably infinite.

# Why define random variables based on  $(\Omega, \mathcal{F}, P)$ ? I: b-7

Answer 1:  $(\Omega, \mathcal{F}, P)$  is what truly occurs internally,

but is possibly **non-observable**.

- **–** In order to infer which of the *non-observable* <sup>ω</sup> occurs, an experiment is performed resulting in an observable x that is a function of  $\omega$ .
- **–** Such experiment yields the random variable X whose probability is defined over the probability space  $(\Omega, \mathcal{F}, P)$ .

Answer 2: With the underlying probability space, any finite dimensional distribution of  $\{X_t, t \in I\}$  is well-defined.

**–** For example,

$$
\Pr[X_1 \le x_1, X_5 \le x_5, X_9 \le x_9] \n= P\left(\{\omega \in \Omega : X_1(\omega) \le x_1, X_5(\omega) \le x_5, X_9(\omega) \le x_9\}\right)
$$

- In many applications, we are perhaps more interested in the distribution functions of random variables than the underlying probability space on which they are defined.
- It can be proved [Billingsley, Thm. 14.1] that <sup>g</sup>iven <sup>a</sup> real-valued non-negative function  $F(\cdot)$  that is non-decreasing and right-continuous and satisfies

$$
\lim_{x \downarrow -\infty} F(x) = 0 \quad \text{and} \quad \lim_{x \uparrow \infty} F(x) = 1,
$$

there exist <sup>a</sup> random variable and an underlying probability space such that the cumulative distribution function (cdf) of the random variable,  $Pr[X \le x]$  =  $P_X ((-\infty, x])$ , defined over the probability space is equal to  $F(\cdot)$ .

- This result releases us from the burden of referring to <sup>a</sup> probability space before defining the random variable. In other words, we can define <sup>a</sup> random variable X directly by its cdf,  $F_X(x) = Pr[X \leq x]$ , without bothering to refer to its underlying probability space.
- Nevertheless, it is important to keep in mind that, formally, random variables are defined over underlying probability spaces.

### Generalization of random variables I: b-9

• The definition of a random variable  $X$  can be generalized by allowing it to take values that are not real numbers:

**Definition** A random variable over the probability space  $(\Omega, \mathcal{F}, P)$  is a function  $X: \Omega \to \mathcal{X}$  satisfying the property that for every  $F \in \mathcal{F}_X$ ,

$$
X^{-1}(F) := \{ w \in \Omega : X(w) \in F \} \in \mathcal{F},
$$

where the alphabet X is a general set and  $\mathcal{F}_X$  is a  $\sigma$ -field of subsets of X [R. M. Gray 2010, P. C. Shields 1991].

• Contrary to the standard definition of a random variable (by taking  $\mathcal{X} = \mathbb{R}$ ), the elements in  $\mathcal X$  may not have a pre-defined ordering; thus, the cdf,

$$
\Pr[X \le x] = P(\{w \in \Omega : X(w) \le x\}),
$$

needs to be explicitly defined.

• Note that this extension definition of a random variable allows  $\mathcal X$  to be an arbitrary (often finite) set so that <sup>a</sup> random source taking values from, e.g., English alphabet, can now be regarded as <sup>a</sup> random variable.

- Statistical evolution in time is an important factor for <sup>a</sup> random source.
- In particular, <sup>a</sup> "time-shift" property should be noted first.

**Definition.** An event E is said to be  $\mathbb{T}$ -invariant with respect to the left-shift (or shift transformation)  $\mathbb{T} \colon \mathcal{X}^{\infty} \to \mathcal{X}^{\infty}$  if

$$
\mathbb{T} E \subseteq E,
$$

where

$$
\mathbb{T}E := \{ \mathbb{T}\boldsymbol{x} \colon \boldsymbol{x} \in E \} \quad \text{and} \quad \mathbb{T}\boldsymbol{x} := \mathbb{T}(x_1, x_2, x_3, \ldots) = (x_2, x_3, \ldots).
$$

• In other words,  $\mathbb T$  is equivalent to "chopping the first component."

**Example.** Applying T onto an event E defined below,

$$
E := \{(x_1 = 1, x_2 = 1, x_3 = 1, x_4 = 1, \ldots), (x_1 = 0, x_2 = 1, x_3 = 1, x_4 = 1, \ldots),
$$
  
\n
$$
(x_1 = 0, x_2 = 0, x_3 = 1, x_4 = 1, \ldots)\},
$$
  
\n
$$
= \{\underbrace{1111 \ldots}_{all \text{ one}}\}, \underbrace{0111 \ldots}_{all \text{ one but the first}}\}
$$
  
\n(B.3.1)

yields

$$
\mathbb{T}E = \{(x_1 = 1, x_2 = 1, x_3 = 1, \ldots), (x_1 = 1, x_2 = 1, x_3 = 1 \ldots),
$$
  
\n
$$
(x_1 = 0, x_2 = 1, x_3 = 1, \ldots)\}
$$
  
\n
$$
= \{(x_1 = 1, x_2 = 1, x_3 = 1, \ldots), (x_1 = 0, x_2 = 1, x_3 = 1, \ldots)\}
$$
  
\n
$$
= \{\underbrace{1111 \ldots}_{all \text{ one}}\}, \underbrace{0111 \ldots}_{all \text{ one but the first}}\}
$$

We then have  $\mathbb{T}E \subseteq E$ , and hence E is T-invariant.

<sup>E</sup> will ge<sup>t</sup> smaller and smaller (more condensed) as time evolves.

 $\Box$ 

- It can be proved (cf. the textbook) that if  $\mathbb{T}E \subseteq E$ , then  $\mathbb{T}^2E \subseteq \mathbb{T}E$ .
- By induction, we can further obtain

$$
\cdots \subseteq \mathbb{T}^3 E \subseteq \mathbb{T}^2 E \subseteq \mathbb{T} E \subseteq E.
$$

- Thus, if an element say  $(1,0,0,1,0,0,\ldots)$  is in a T-invariant set E, then all its left-shift counterparts (i.e.,  $(0, 0, 1, 0, 0, 1...)$  and  $(0, 1, 0, 0, 1, 0, ...)$ ) should be contained in E.
- As a result, for a  $\mathbb{T}$ -invariant set E, an element and all its left-shift counterparts are either all in  $E$  or all outside  $E$ , but cannot be partially inside  $E$ .
- Hence, a "T-invariant group" such as one containing

 $(1, 0, 0, 1, 0, 0, \ldots), (0, 0, 1, 0, 0, 1, \ldots)$  and  $(0, 1, 0, 0, 1, 0, \ldots)$ 

should be treated as an indecomposable group in T-invariant sets.

- Although we are in particular interested in these "T-invariant indecomposable groups" (especially when defining an ergodic random process), it is possible that some single "transient" element, such as  $(0, 0, 1, 1, \ldots)$  in  $(B.3.1)$ , is included in <sup>a</sup> T-invariant set, and will be excluded after applying left-shift operation T.
- This however can be resolved by introducing the inverse operation  $\mathbb{T}^{-1}$ .
- Note that  $\mathbb T$  is a many-to-one mapping, so its inverse operation does not exist in general.
- Similar to taking the closure of an open set, the definition adopted below [P. C. Shields 1991, p. 3] allows us to "enlarge" the T-invariant set such that all right-shift counterparts of the single "transient" element are included:

$$
\mathbb{T}^{-1}E:=\{\boldsymbol{x}\in\mathcal{X}^{\infty}\colon\mathbb{T}\boldsymbol{x}\in E\}\,.
$$

• We then notice from the above definition that if

$$
\mathbb{T}^{-1}E = E,\tag{B.3.2}
$$

then

$$
\mathbb{T} E = \mathbb{T} (\mathbb{T}^{-1} E) = E,
$$

and hence  $E$  is constituted only by the  $\mathbb{T}$ -invariant groups because

$$
\cdots = \mathbb{T}^{-2}E = \mathbb{T}^{-1}E = E = \mathbb{T}E = \mathbb{T}^2E = \cdots.
$$

• The sets that satisfy (B.3.2) are sometimes referred to as *ergodic sets* because as time goes by (the left-shift operator T can be regarded as <sup>a</sup> shift to <sup>a</sup> future time), the set always stays in the state that it has been before.

- As the textbook only deals with one-sided random processes, the discussion on T-invariance only focuses on sets of one-sided sequences.
- When a two-sided random process  $\ldots, X_{-2}, X_{-1}, X_0, X_1, X_2, \ldots$  is considered, the left-shift operation T of <sup>a</sup> two-sided sequence actually has <sup>a</sup> unique inverse. Hence,  $\mathbb{T}E \subseteq E$  implies  $\mathbb{T}E = E$ . Also,  $\mathbb{T}E = E$  iff  $\mathbb{T}^{-1}E = E$ . Ergodicity for two-sided sequences can therefore be directly defined using  $\mathbb{T}E = E$ .

We now classify several useful statistical properties of (one-sided) random process

$$
\mathbf{X} = \{X_1, X_2, \ldots\}.
$$

- *Memoryless***:** A random process or <sup>a</sup> source *X* is said to be *memoryless* if the sequence of random variables  $X_i$  is *independent* and *identically distributed*  $(i.i.d.).$
- *Stationary process***:** A process is said to be *stationary* (or *strictly stationary*) if the probability of every sequence or event is unchanged by <sup>a</sup> left (time) shift.
- *Ergodic process***:** A process is said to be *ergodic* if any ergodic set (satisfying  $(B.3.2)$ ) in  $\mathcal{F}_X$  has probability either 1 or 0. This definition is not very intuitive, but some interpretations and examples may shed some light.
	- Observe that the definition has nothing to do with stationarity. It simply states that events that are unaffected by time-shifting (both left- and rightshifting) must have probability either zero or one.
	- Ergodicity implies that all convergen<sup>t</sup> time averages converge to <sup>a</sup> constant (but not necessarily to the ensemble average or statistical expectation).

Below is an example that can be used to explain the idea.

**Example.** Suppose  $X_1, X_2, \cdots, X_n, \cdots$  is an ergodic process, where each  $X_n$  takes values in  $\{0,1\}$ . Let  $\Omega$  be the set of all one-sided zero-one sequences. Define for  $\alpha \in [0,1],$ 

$$
\bar{E}_n(\alpha) := \left\{ \boldsymbol{x} \in \{0, 1\}^\infty \; : \; \alpha \le \limsup_{m \to \infty} \frac{x_1 + \dots + x_m}{m} < \alpha + \frac{1}{n} \right\}
$$

and

$$
\underline{E}_n(\alpha) := \left\{ \boldsymbol{x} \in \{0,1\}^\infty \; : \; \alpha \leq \liminf_{m \to \infty} \frac{x_1 + \dots + x_m}{m} < \alpha + \frac{1}{n} \right\}.
$$

Then it can be verified that both  $\bar{E}_n(\alpha)$  and  $\underline{E}_n(\alpha)$  are ergodic sets, i.e.,

$$
\bar{E}_n(\alpha) = \mathbb{T}^{-1} \bar{E}_n(\alpha)
$$
 and  $\underline{E}_n(\alpha) = \mathbb{T}^{-1} \underline{E}_n(\alpha)$ .

Observe that

$$
\Omega = \bigcup_{k=0}^{n} \bar{E}_n\left(\frac{k}{n}\right) = \bigcup_{k=0}^{n} \underline{E}_n\left(\frac{k}{n}\right)
$$

and

$$
\bar{E}_n\left(\frac{k}{n}\right)\bigcap \bar{E}_n\left(\frac{\ell}{n}\right)=\underline{E}_n\left(\frac{k}{n}\right)\bigcap \underline{E}_n\left(\frac{\ell}{n}\right)=\emptyset \text{ for } k\neq \ell.
$$

The definition of ergodicity implies the existence of  $k$  and  $\ell$  such that

$$
\Pr\left[\boldsymbol{X} \in \bar{E}_n\left(\frac{k}{n}\right)\right] = \Pr\left[\boldsymbol{X} \in \underline{E}_n\left(\frac{\ell}{n}\right)\right] = 1.
$$

If  $\frac{X_1 + \dots + X_n}{n}$  converges with probability one, then  $k = \ell$ .

In other words,

$$
\frac{k}{n} \le \limsup_{m \to \infty} \frac{X_1 + \dots + X_m}{m} < \frac{k+1}{n} \text{ with probability 1}
$$

and

$$
\frac{k}{n} \le \liminf_{m \to \infty} \frac{X_1 + \dots + X_m}{m} < \frac{k+1}{n} \text{ with probability 1.}
$$

As <sup>a</sup> result,

$$
\left|\limsup_{m\to\infty}\frac{X_1+\dots+X_m}{m}-\liminf_{m\to\infty}\frac{X_1+\dots+X_m}{m}\right|<\frac{1}{n}\text{ with probability 1.}
$$

Ergodicity implies that all convergen<sup>t</sup> time averages converge to <sup>a</sup> constant.

- It needs to be pointed out that in the above example, ergodicity does not guarantee that the ensemble average lies in  $\lfloor k/n, (k+1)/n \rfloor$ .
- A quick example is that

$$
Pr{(x_1 = 0, x_2 = 1, x_3 = 0, x_4 = 1, ...)} = 0.2
$$

and

$$
Pr{(x_1 = 1, x_2 = 0, x_3 = 1, x_4 = 0, \ldots)} = 0.8
$$

assure the validity of ergodicity, but

$$
\frac{X_1 + \dots + X_n}{n} \to \frac{1}{2}
$$
 with probability 1.

which is not equal to  $E[X_i]$  for any i.

- In principle,
	- **–** ergodicity implies that all convergen<sup>t</sup> sample averages converge to <sup>a</sup> constant (but not necessarily to the statistical expectation), and
	- **–** stationarity assures that the time average converges to <sup>a</sup> random variable;

hence, it is reasonable to expect that they jointly imply the ultimate time average equals the ensemble average. This is validated by the well-known *ergodic theorem* by Birkhoff and Khinchin.

**Theorem B.4 (Pointwise ergodic theorem)** Consider <sup>a</sup> discrete-time stationary random process,  $\mathbf{X} = \{X_n\}_{n=1}^{\infty}$ . For real-valued function  $f(\cdot)$  on R with finite mean (i.e.,  $|E[f(X_n)]| < \infty$ ), there exists a random variable Y such that

$$
\lim_{n \to \infty} \frac{1}{n} \sum_{k=1}^{n} f(X_k) = Y
$$
 with probability 1.

If, in addition to stationarity, the process is also ergodic, then

$$
\lim_{n \to \infty} \frac{1}{n} \sum_{k=1}^{n} f(X_k) = E[f(X_1)] \text{ with probability 1.}
$$

# Operational meaning of stationary ergodic assumption I: b-21

- Stationary ergodic random source
	- **–** One of the important consequences that the pointwise ergodic theorem indicates is that the time average can ultimately replace the statistical average, which is <sup>a</sup> useful result.
	- **–** Hence, with stationarity and ergodicity, one, who observes

$$
X_1^{30} = 154326543334225632425644234443
$$

from the experiment of rolling <sup>a</sup> dice, can draw the conclusion that the true distribution of rolling the dice can be well approximated by:

$$
\Pr{X_i = 1} \approx \frac{1}{30} \qquad \Pr{X_i = 2} \approx \frac{6}{30} \qquad \Pr{X_i = 3} \approx \frac{7}{30}
$$
\n
$$
\Pr{X_i = 4} \approx \frac{9}{30} \qquad \Pr{X_i = 5} \approx \frac{4}{30} \qquad \Pr{X_i = 6} \approx \frac{3}{30}
$$

- **–** Such result is also known by the *law of large numbers*. The relation between ergodicity and the law of large numbers will be further explored later.
- Non-stationary or non-ergodic source
	- **–** Empirical distribution (relative frequency) cannot necessarily be used to approximate the true distribution.

# Operational meaning of stationary ergodic assumption I: b-22

- In communications theory, one may assume that *the source is stationary* or *the source is stationary ergodic*. But it is not common to see the assumption of *the source being ergodic but non-stationary*. Why?
	- **–** This is perhaps because an ergodic but non-stationary source in general does not facilitate the analytical study of communications problems.
- This, to some extent, justifies that the *ergodicity* assumption usually comes after *stationarity* assumption. A specific example is the pointwise ergodic theorem, where the random processes considered is presumed to be stationary.

We continue to classify useful statistical properties of (one-sided) random process

$$
\mathbf{X} = \{X_1, X_2, \ldots\}.
$$

 $\Diamond$  Markov chain for three random variables:

Three random variables  $X, Y$  and  $Z$  are said to form a Markov chain if

$$
P_{X,Y,Z}(x,y,z) = P_X(x) \cdot P_{Y|X}(y|x) \cdot P_{Z|Y}(z|y); \tag{B.3.3}
$$

i.e.,

$$
P_{Z|X,Y}(z|x,y) = P_{Z|Y}(z|y).
$$

This is usually denoted by

$$
X \to Y \to Z.
$$

• " $X \to Y \to Z$ " is sometimes read as "X and Z are conditionally independent given  $Y$ " because it can be shown that  $(B.3.3)$  is equivalent to

$$
P_{X,Z|Y}(x,z|y) = P_{X|Y}(x|y) \cdot P_{Z|Y}(z|y).
$$

• Therefore, " $X \to Y \to Z$ " is equivalent to " $Z \to Y \to X$ ". Accordingly, the Markovian notation is sometimes expressed as " $X \leftrightarrow Y \leftrightarrow Z$ ".

 $\diamond$  *k*th-order Markov sources:

The sequence of random variables  $\{X_n\}_{n=1}^{\infty} = X_1, X_2, X_3, \ldots$  with common finite-alphabet  $\mathcal X$  is said to form a k-th order Markov chain (or k-th order Markov source or process) if for all  $n > k$ ,  $x_i \in \mathcal{X}$ ,  $i = 1, \ldots, n$ ,

$$
\Pr[X_n = x_n | X_{n-1} = x_{n-1}, \dots, X_1 = x_1] \\
= \Pr[X_n = x_n | X_{n-1} = x_{n-1}, \dots, X_{n-k} = x_{n-k}].
$$
\n(B.3.4)

Each  $x^{n-1}$  $x_{n-k}^{n-1} := (x_{n-k}, x_{n-k+1}, \ldots, x_{n-1}) \in \mathcal{X}^k$  is called the *state* of the Markov chain at time  $n$ .

• **Irreducible:** A Markov chain is *irreducible* if with some (non-zero) probability, we can go from any state in  $\mathcal{X}^k$  to another state in a finite number of steps, i.e., for all  $x^k, y^k \in \mathcal{X}^k$  there exists  $j \geq 1$  such that

$$
\Pr\left\{X_j^{k+j-1} = x^k \middle| X_1^k = y^k \right\} > 0.
$$

• **Time-invariant:** A Markov chain is said to be *time-invariant* or *homogeneous*, if for every  $n > k$ ,

$$
\Pr[X_n = x_n | X_{n-1} = x_{n-1}, \dots, X_{n-k} = x_{n-k}]
$$
  
= 
$$
\Pr[X_{k+1} = x_{k+1} | X_k = x_k, \dots, X_1 = x_1].
$$

**–** Therefore, <sup>a</sup> homogeneous first-order Markov chain can be defined through its transition probability:

$$
[Pr{X_2 = x_2 | X_1 = x_1}]_{|\mathcal{X}| \times |\mathcal{X}|},
$$

and its initial state distribution  $P_{X_1}(x)$ .

#### • **Aperiodic:**

 $-$ In a first-order Markov chain, the *period*  $d(x)$  of state  $x \in \mathcal{X}$  is defined by

$$
d(x) := \gcd\left\{n \in \{1, 2, 3, \ldots\} : \Pr\{X_{n+1} = x | X_1 = x\} > 0\right\},\
$$

where gcd denotes the greatest common divisor; in other words, if the Markov chain starts in state x, then the chain cannot return to state x at any time that is not a multiple of  $d(x)$ .

- $-$  If  $Pr{X_{n+1} = x | X_1 = x} = 0$  for all n, we say that state x has an infinite period and write  $d(x) = \infty$ .
- $-$  We also say that *state* x *is aperiodic* if  $d(x) = 1$  and *periodic* if  $d(x) >$ 1.
- **–** The first-order Markov chain is called **aperiodic** if all its states are aperiodic. In other words, the first-order Markov chain is aperiodic if

$$
\gcd\{n \in \{1, 2, 3, \ldots\}: \Pr\{X_{n+1} = x | X_1 = x\} > 0\} = 1 \quad \forall x \in \mathcal{X}.
$$

**Property.** In an irreducible first-order Markov chain, all states have the same period. Hence, if one state in such <sup>a</sup> chain is aperiodic, then the entire Markov chain is aperiodic.

**• Stationarity:** A distribution  $\pi(\cdot)$  on X is said to be a *stationary* distribution for <sup>a</sup> homogeneous (i.e., time-invariant) first-order Markov chain, if for every  $y \in \mathcal{X},$ 

$$
\pi(y) = \sum_{x \in \mathcal{X}} \pi(x) \Pr\{X_2 = y | X_1 = x\}.
$$

#### **Properties.**

- 1. For a finite-alphabet homogeneous first-order Markov chain,  $\pi(\cdot)$  always exists.
- 2.  $\pi(\cdot)$  is unique if the Markov chain is irreducible.
- 3. For <sup>a</sup> finite-alphabet homogeneous first-order Markov chain that is both irreducible and aperiodic,

$$
\lim_{n \to \infty} \Pr\{X_{n+1} = y | X_1 = x\} = \pi(y)
$$

for all states x and y in  $\mathcal{X}$ .

If the initial state distribution is equal to <sup>a</sup> stationary distribution, then the homogeneous first-order Markov chain becomes <sup>a</sup> stationary process.



# B.4 Convergence of sequences of random variables  $\qquad$ <sub>I: b-29</sub>

• Relation of five modes of convergence

$$
X_n \xrightarrow{\text{p.w.}} X
$$
\n
$$
\downarrow
$$
\n
$$
X_n \xrightarrow{a.s.} X \xrightarrow{\text{Thm. B.10}} X_n \xrightarrow{L_r} X \ (r \ge 1)
$$
\n
$$
X_n \xrightarrow{p} X
$$
\n
$$
\downarrow
$$
\n
$$
X_n \xrightarrow{d} X
$$

## B.4 Convergence of sequences of random variables  $\qquad$ <sub>I: b-30</sub>

• Pointwise convergence and almost surely convergence

**Example B.7** Give a probability space

 $(\Omega = \{0, 1, 2, 3\}, 2^{\Omega}, P(0) = P(1) = P(2) = 1/3).$ 

 $-$  A random variable  $X_n$  is a mapping from a probability space to  $\mathbb{R}$ . Let the mapping be

$$
X_n(\omega) = \frac{\omega}{n} \implies \Pr\{X_n = 0\} = \Pr\left\{X_n = \frac{1}{n}\right\} = \Pr\left\{X_n = \frac{2}{n}\right\} = \frac{1}{3}.
$$

**–** (*Pointwise convergence*) Observe that

$$
(\forall \omega \in \Omega) \; X_n(\omega) \to X(\omega),
$$

where  $X(\omega) = 0$  for every  $\omega \in \Omega$ . So

$$
X_n \xrightarrow{p.w} X.
$$

 $-$  (*Almost surely convergence*) Let  $\tilde{X}(\omega) = 0$  for  $\omega = 0, 1, 2$  and  $\tilde{X}(\omega) = 1$ for  $\omega = 3$ . Then both of the following statements are true:

$$
X_n \xrightarrow{a.s.} X \quad \text{and} \quad X_n \xrightarrow{a.s.} \tilde{X},
$$

(since

$$
\Pr\left\{\lim_{n\to\infty}X_n=\tilde{X}\right\}=\sum_{\omega=0}^3P(\omega)\cdot\mathbf{1}\left\{\lim_{n\to\infty}X_n(\omega)=\tilde{X}(\omega)\right\}=1.
$$

# B.4 Convergence of sequences of random variables I: b-31

• Almost surely convergence (with probability 1) and convergence in probability

$$
X_n \xrightarrow{a.s.} X \equiv \Pr\left\{ \lim_{n \to \infty} X_n = X \right\} = 1
$$
  

$$
X_n \xrightarrow{p} X \equiv (\forall \gamma > 0) \lim_{n \to \infty} \Pr\left\{ |X_n - X| < \gamma \right\} = 1
$$

• Convergence in  $r$ th mean

$$
X_n \xrightarrow{L_r} X \equiv \lim_{n \to \infty} E\left[ |X_n - X|^r \right] = 0
$$

• Convergence in distribution

$$
X_n \xrightarrow{d} X \equiv \lim_{n \to \infty} F_{X_n}(x) = F_X(x)
$$
 for every continuous point of  $F_X(x)$ 

# B.4 Convergence of sequences of random variables I: b-32

The next observation facilitates the finding of limiting random variable.

#### **Observation B.8 (Uniqueness of convergence)**

1. If  $X_n \xrightarrow{p.w.} X$  and  $X_n \xrightarrow{p.w.} Y$ , then  $X = Y$  pointwisely. I.e.,

$$
(\forall \omega \in \Omega) \quad X(\omega) = Y(\omega).
$$

2. If 
$$
X_n \xrightarrow{a.s.} X
$$
 and  $X_n \xrightarrow{a.s.} Y$   
\n(or  $X_n \xrightarrow{p} X$  and  $X_n \xrightarrow{p} Y$ )  
\n(or  $X_n \xrightarrow{L_r} X$  and  $X_n \xrightarrow{L_r} Y$ ),  
\nthen  $X = Y$  with probability 1. I.e.,

$$
\Pr\{X = Y\} = 1.
$$

3. 
$$
X_n \xrightarrow{d} X
$$
 and  $X_n \xrightarrow{d} Y$ ,  
then  $F_X(x) = F_Y(x)$  for all x.

# B.4 Convergence of sequences of random variables I: b-33

**Theorem B.9 (Monotone convergence theorem)**

(i) 
$$
X_n \xrightarrow{a.s.} X
$$
  
\n(ii)  $(\forall n) Y \le X_n \le X_{n+1}$   $\rightarrow X_n \xrightarrow{L_1} X \Rightarrow E[X_n] \rightarrow E[X].$   
\n(iii)  $E[|Y|] < \infty$ 

**Theorem B.10 (Dominated convergence theorem)**

(i) 
$$
X_n \xrightarrow{a.s.} X
$$
  
\n(ii)  $(\forall n) |X_n| \le Y$   $\Rightarrow X_n \xrightarrow{L_1} X \Rightarrow E[X_n] \rightarrow E[X].$   
\n(iii)  $E[|Y|] < \infty$ 

The implication of  $X_n \stackrel{L_1}{\longrightarrow} X$  to  $E[X_n] \to E[X]$  can be easily seen from

$$
|E[X_n] - E[X]| = |E[X_n - X]| \le E[|X_n - X|].
$$

**Theorem B.13 (Weak law of large numbers)** Let  $\{X_n\}_{n=1}^{\infty}$  be a sequence of uncorrelated random variables with common mean  $E[X_i] = \mu$ . If the variables also have common variance, or more generally,

$$
\lim_{n \to \infty} \frac{1}{n^2} \sum_{i=1}^n \text{Var}[X_i] = 0, \quad \text{(equivalently, } \frac{X_1 + \dots + X_n}{n} \xrightarrow{L_2} \mu)
$$

then

$$
\frac{X_1 + \dots + X_n}{n} \xrightarrow{p} \mu.
$$

**proof**: By Chebyshev's inequality,

$$
\Pr\left\{\left|\frac{1}{n}\sum_{i=1}^n X_i - \mu\right| \ge \varepsilon\right\} \le \frac{1}{n^2 \varepsilon^2} \sum_{i=1}^n \text{Var}[X_i].
$$

 $\Box$ 

**Note:**  $X_n \stackrel{\mathcal{L}_2}{\longrightarrow} X$  implies  $X_n \stackrel{p}{\longrightarrow} X$ .

 $\bf{Theorem~B.14}$   $(\bf{Kolmogorov's\, strong\, law\, of\, large\, numbers})$  Let  $\{X_n\}_{n=1}^\infty$ be an independent sequence of random variables with common mean  $E[X_n] = \mu$ . If either

- 1.  $X_n$ 's are identically distributed; or
- 2.  $X_n$ 's are square-integrable with

$$
\sum_{i=1}^{\infty} \frac{\text{Var}[X_i]}{i^2} < \infty,
$$

Then

$$
\frac{X_1 + \dots + X_n}{n} \xrightarrow{a.s.} \mu.
$$

**Note**: The difference of *weak* and *strong* laws of large number is that the former is convergence *in probability*, while the latter is *almost sure* convergence.

- After the introduction of Kolmogorov's strong law of large numbers, one may find that the pointwise ergodic theorem (Theorem B.4) actually indicates <sup>a</sup> similar result.
	- **–** In fact, the pointwise ergodic theorem can be viewed as another version of strong law of large numbers, which states that *for stationary and ergodic processes, time averages converge with probability* 1 *to the ensemble expectation.*
- The notion of ergodicity is often misinterpreted, since the definition is not very intuitive. Some engineering texts may provide <sup>a</sup> definition that <sup>a</sup> stationary process satisfying the ergodic theorem is also ergodic.

# $B.5$  Ergodicity and law of large numbers  $\qquad \qquad$

Let us try to clarify the notion of ergodicity by the following remarks.

- The concept of ergodicity does not require stationarity. In other words, <sup>a</sup> non-stationary process can be ergodic.
- Many perfectly good models of <sup>p</sup>hysical processes are not ergodic, ye<sup>t</sup> they have <sup>a</sup> form of law of large numbers. In other words, non-ergodic processes can be perfectly good and useful models.
- There is no finite-dimensional equivalent definition of ergodicity as there is for stationarity. This fact makes it more difficult to describe and interpret ergodicity.
- I.i.d. processes are ergodic; hence, ergodicity can be thought of as <sup>a</sup> (kind of) generalization of i.i.d.
- As mentioned earlier, stationarity and ergodicity imply the time average converges with probability 1 to the ensemble mean. Now if <sup>a</sup> process is stationary but not ergodic, then the time average still converges, but possibly not to the ensemble mean.

## $B.5$  Ergodicity and law of large numbers  $\qquad \qquad$

**Example.** Let  $\{A_n\}_{n=1}^{\infty}$  $\sum_{n=-\infty}^{\infty}$  and  $\{B_n\}_{n=1}^{\infty}$  $\sum_{n=-\infty}^{\infty}$  be two i.i.d. binary 0-1 random variables with

$$
Pr{A_n = 0} = Pr{B_n = 1} = 1/4.
$$

Suppose that

$$
X_n = \begin{cases} A_n, & \text{if } U = 1 \\ B_n, & \text{if } U = 0, \end{cases}
$$

where U is equiprobable binary random variable, and  $\{A_n\}_{n=1}^{\infty}$ ,  $\{B_n\}_{n=1}^{\infty}$  and U are independent.

Then  $\{X_n\}_{n=1}^{\infty}$  is stationary.

Is the process ergodic? The answer is negative.

If the stationary process were ergodic, then from the pointwise ergodic theorem (Theorem B.4), its relative frequency would converge to <sup>a</sup> constant!

## $B.5$  Ergodicity and law of large numbers  $\qquad \qquad$

However, one should observe that the outputs of  $(X_1,\ldots,X_n)$  form a Bernoulli process with relative frequency of 1's being either 3 /<sup>4</sup> or <sup>1</sup> /4, depending on the value of U. Therefore,

$$
\lim_{n \to \infty} \frac{1}{n} \sum_{i=1}^{n} X_n \xrightarrow{a.s.} Y,
$$

where  $Pr(Y = 1/4) = Pr(Y = 3/4) = 1/2$ , which contradicts to the ergodic theorem.  $\Box$ 

• *Ergodic decomposition theorem:* Under fairly genera<sup>l</sup> assumptions, any (not necessarily ergodic) stationary process is <sup>a</sup> mixture of stationary ergodic processes, and hence one always observes <sup>a</sup> stationary ergodic outcome. As in the above example, one always observe either  $A_1, A_2, A_3, \ldots$  or  $B_1, B_2, B_3, \ldots,$  depending on the value of  $U$ , for which both sequences are stationary ergodic (i.e., the time-stationary observation  $X_n$  satisfies

$$
X_n = U \cdot A_n + (1 - U) \cdot B_n.
$$

- The previous remark implies that ergodicity is not required for the strong law of large numbers to be useful.
- The next question is whether or not stationarity is required. Again the answer is negative !

- In fact, what is needed in this course is the **law of large numbers**, which results the convergence of sample averages to its ensemble expectation.
	- **–** It should be reasonable to expect that random processes could exhibit transient behavior that violates the stationarity definition, ye<sup>t</sup> the sample average still converges. One can then introduce the notion of *asymptotically stationary* to achieve the law of large numbers.

## B.6 Central limit theorem I: b-41

**Theorem B.15 (Central limit theorem)** If  $\{X_n\}_{n=1}^{\infty}$  is a sequence of i.i.d. random variables with finite common marginal mean  $\mu$  and variance  $\sigma^2$ , then

$$
\frac{1}{\sqrt{n}}\sum_{i=1}^{n}(X_i-\mu)\stackrel{d}{\longrightarrow}Z\sim\mathcal{N}(0,\sigma^2),
$$

where the convergence is in distribution (as  $n \to \infty$ ) and  $Z \sim \mathcal{N}(0, \sigma^2)$  is a Gaussian distributed random variable with mean 0 and variance  $\sigma^2$ .

## B.7 Convexity, concavity and Jensen's inequality I: b-42

**Definition B.16 (Convexity)** Consider a convex set  $\mathcal{O} \subset \mathbb{R}^m$ , where m is a fixed positive integer. Then a function  $f: \mathcal{O} \to \mathbb{R}$  is said to be *convex* over  $\mathcal{O}$  if for every <u>x</u>, y in  $\mathcal{O}$  and  $0 \leq \lambda \leq 1$ ,

$$
f\left(\lambda \underline{x} + (1-\lambda)\underline{y}\right) \leq \lambda f(\underline{x}) + (1-\lambda)f(\underline{y}).
$$

Furthermore, a function f is said to be *strictly convex* if equality holds only when  $\lambda = 0$  or  $\lambda = 1$ .

• A set  $\mathcal{O} \subset \mathbb{R}^m$  is said to be *convex* if for every  $\underline{x} = (x_1, x_2, \cdots, x_m)^T$  and  $y = (y_1, y_2, \dots, y_m)^T$  in  $\mathcal{O}$  (where T denotes transposition), and every  $0 \leq$  $\lambda \leq 1, \lambda \underline{x} + (1 - \lambda)y \in \mathcal{O}$ ; in other words, the "convex combination" of any two "points"  $\underline{x}$  and  $y$  in  $\mathcal O$  also belongs to  $\mathcal O$ .

**Definition B.17 (Concavity)** A function f is *concave* if − f is convex.

Jensen's inequality I: b-43

**Theorem B.18** (Jensen's inequality) If  $f : \mathcal{O} \to \mathbb{R}$  is convex over a convex set  $\mathcal{O} \subset \mathbb{R}^m$ , and  $\underline{X} = (X_1, X_2, \cdots, X_m)^T$  is an *m*-dimensional random vector with alphabet  $\mathcal{X} \subset \mathcal{O}$ , then

$$
E[f(\underline{X})] \ge f(E[\underline{X}]).
$$

Moreover, if  $f$  is strictly convex, then equality in the above inequality immediately implies  $\underline{X} = E[\underline{X}]$  with probability 1.

Optimization of a function  $f(\boldsymbol{x})$  over  $\boldsymbol{x} = (x_1, \dots, x_n) \in \mathcal{X} \subseteq \mathbb{R}^n$ 

subject to

$$
\begin{cases}\text{inequality constraints } g_i(\boldsymbol{x}) \leq 0 \text{ for } 1 \leq i \leq m, \text{ and} \\
\text{equality constraints } h_j(\boldsymbol{x}) = 0 \text{ for } 1 \leq j \leq \ell\n\end{cases}
$$

is <sup>a</sup> center technique to problems in information theory.

Mathematically, the problem can be formulated as:

$$
\min_{\boldsymbol{x}\in\mathcal{Q}}f(\boldsymbol{x}),\tag{B.8.1}
$$

where

$$
\mathcal{Q} := \{ \boldsymbol{x} \in \mathcal{X} : g_i(\boldsymbol{x}) \leq 0 \text{ for } 1 \leq i \leq m \text{ and } h_i(\boldsymbol{x}) = 0 \text{ for } 1 \leq j \leq \ell \}.
$$

# B.8 Lagrange multipliers tech.  $&$  KKT conditions  $\qquad$ <sub>I: b-45</sub>

- In most cases, solving the constrained optimization problem defined in  $(B.8.1)$ is hard due to the constraints.
- Instead, one may introduce <sup>a</sup> **dual** optimization problem without constraints:

$$
L(\lambda, \nu) := \min_{\mathbf{x} \in \mathcal{X}} \left( f(\mathbf{x}) + \sum_{i=1}^{m} \lambda_i g_i(\mathbf{x}) + \sum_{j=1}^{\ell} \nu_j h_j(\mathbf{x}) \right) = \min_{\mathbf{x} \in \mathcal{X}} L(\mathbf{x}; \lambda, \nu).
$$
\n(B.8.2)

- In the literature,  $\boldsymbol{\lambda} = (\lambda_1, \ldots, \lambda_m)$  and  $\boldsymbol{\nu} = (\nu_1, \ldots, \nu_\ell)$  are usually referred to as **Lagrange multipliers**, and <sup>L</sup>(*λ*, *<sup>ν</sup>*) is called the **Lagrange dual function**.
	- $-$  Note that  $L(\lambda, \nu)$  is a concave function of  $\lambda$  and  $\nu$  since it is the minimization of affine functions of *λ* and *ν*.

• It can be verified that when  $\lambda_i \geq 0$  for  $1 \leq i \leq m$ ,

$$
L(\lambda, \nu) \le \min_{\mathbf{x} \in \mathcal{Q}} \left( f(\mathbf{x}) + \sum_{i=1}^{m} \lambda_i g_i(\mathbf{x}) + \sum_{j=1}^{\ell} \nu_j h_j(\mathbf{x}) \right) \le \min_{\mathbf{x} \in \mathcal{Q}} f(\mathbf{x}). \quad (B.8.3)
$$

• We are however interested in when the above inequality becomes equality (i.e., when the so-called *strong duality* holds) because if there exist non-negative  $\tilde{\lambda}$ and  $\tilde{\nu}$  that equate (B.8.3), then

$$
f(\boldsymbol{x}^*) = \min_{\boldsymbol{x} \in \mathcal{Q}} f(\boldsymbol{x})
$$
  
\n
$$
= L(\tilde{\boldsymbol{\lambda}}, \tilde{\boldsymbol{\nu}}) = \min_{\boldsymbol{x} \in \mathcal{X}} \left( f(\boldsymbol{x}) + \sum_{i=1}^m \tilde{\lambda}_i g_i(\boldsymbol{x}) + \sum_{j=1}^\ell \tilde{\nu}_j h_j(\boldsymbol{x}) \right)
$$
  
\n
$$
\leq f(\boldsymbol{x}^*) + \sum_{i=1}^m \tilde{\lambda}_i g_i(\boldsymbol{x}^*) + \sum_{j=1}^\ell \tilde{\nu}_j h_j(\boldsymbol{x}^*)
$$
  
\n
$$
\leq f(\boldsymbol{x}^*), \qquad (B.8.4)
$$

where (B.8.4) follows because the minimizer *<sup>x</sup>*<sup>∗</sup> of (B.8.1) lies in Q.

# B.8 Lagrange multipliers tech.  $&$  KKT conditions  $\qquad$ <sub>E</sub>: b-47

• Hence, if the strong duality holds, the same *<sup>x</sup>*<sup>∗</sup> achieves both

$$
\min_{\bm{x}\in\mathcal{Q}}f(\bm{x})
$$

and

$$
L(\tilde{\boldsymbol{\lambda}},\tilde{\boldsymbol{\nu}}),
$$

and  $\tilde{\lambda}_i g_i(\boldsymbol{x}^*) = 0$  for  $1 \leq i \leq m$ .<sup>1</sup>

- The strong duality does not in general hold.
- A situation that guarantees the validity of the strong duality has been determined by William Karush [212] (1936), and separately Harold W. Kuhn and Albert W. Tucker [235] (1951).
- In particular, when  $f(\cdot)$  and  $\{g_i(\cdot)\}_{i=1}^m$  are both convex, and  $\{h_j(\cdot)\}_{j=1}^{\ell}$  are affine, and these functions are all differentiable, they found that the strong duality holds if, and only if, the KKT condition is satisfied [56, p. 258].
	- **–** Again, we are free to choose *λ* and *<sup>ν</sup>* that satisfy the KKT condition (cf. Definition B.19).

 $^{1}$ Equating (B.8.4) implies  $\sum_{i=1}^{m} \tilde{\lambda}_{i} g_{i}(x^{*}) = 0$ . It can then be easily verified from  $\tilde{\lambda}_{i} g_{i}(x^{*}) \leq 0$  for every  $1 \leq i \leq m$  that  $\tilde{\lambda}_i g_i(\boldsymbol{x}^*) = 0$  for  $1 \leq i \leq m$ .

## B.8 Lagrange multipliers tech.  $&$  KKT conditions  $\qquad$ <sub>I: b-48</sub>

 $\bf{Definition~B.19}$   $(\bf{Karush-Kuhn-Tucker}$   $(\bf{KKT})$   $\bf{condition})$   $\rm{Point}$   $\boldsymbol{x}$  =  $(x_1,$ ...,  $x_n$ ) and multipliers  $\boldsymbol{\lambda} = (\lambda_1, \ldots, \lambda_m)$  and  $\boldsymbol{\nu} = (\nu_1, \ldots, \nu_\ell)$  are said to satisfy the KKT condition if

$$
\begin{cases}\ng_i(\boldsymbol{x}) \leq 0, & \lambda_i \geq 0, & \lambda_i g_i(\boldsymbol{x}) = 0 & i = 1, \dots, m \\
h_j(\boldsymbol{x}) = 0 & j = 1, \dots, \ell \\
\frac{\partial L}{\partial x_k}(\boldsymbol{x}; \boldsymbol{\lambda}, \boldsymbol{\nu}) = \frac{\partial f}{\partial x_k}(\boldsymbol{x}) + \sum_{i=1}^m \lambda_i \frac{\partial g_i}{\partial x_k}(\boldsymbol{x}) + \sum_{j=1}^\ell \nu_j \frac{\partial h_j}{\partial x_k}(\boldsymbol{x}) = 0 & k = 1, \dots, n\n\end{cases}
$$

- Note that when  $f(\cdot)$  and constraints  $\{g_i(\cdot)\}_{i=1}^m$  and  $\{h_j(\cdot)\}_{j=1}^{\ell}$  are arbitrary functions, the KKT condition is only <sup>a</sup> necessary condition for the validity of the strong duality.
- In other words, for <sup>a</sup> non-convex optimization, we can only claim that if the strong duality holds, then the KKT condition is satisfied but not vice versa.

- A case that is particularly useful in information theory is when  $x$  is restricted to be <sup>a</sup> probability distribution.
- In such case, apart from other problem-specific constraints, we have additionally

 $\begin{cases} n \text{ inequality constraints } g_{m+i}(\boldsymbol{x}) = -x_i \leq 0 \text{ for } 1 \leq i \leq n, \text{ and} \\ \text{one equality constraint } h_{\ell+1}(\boldsymbol{x}) = \sum_{k=1}^n x_k - 1 = 0. \end{cases}$ one equality constraint  $h_{\ell+1}(\boldsymbol{x}) = \sum_{k=1}^{n} h_k(\boldsymbol{x})$  $\int_{k=1}^{n} x_k - 1 = 0.$ 

The above relation is the mostly seen form of the KKT condition when it is used in problems of information theory.

**Example B.20** Suppose for non-negative  $\{q_{i,j}\}_{1 \leq i \leq n, 1 \leq j \leq n'}$  with  $\sum_{j=1}^{n'}$  $\sum_{j=1}^n q_{i,j} = 1,$ 

$$
\begin{cases}\nf(\boldsymbol{x}) = -\sum_{i=1}^{n} \sum_{j=1}^{n'} x_i q_{i,j} \log \frac{q_{i,j}}{\sum_{i'=1}^{n} x_{i'} q_{i',j}} \\
g_i(\boldsymbol{x}) = -x_i \leq 0 & i = 1, ..., n \\
h(\boldsymbol{x}) = \sum_{i=1}^{n} x_i - 1 = 0\n\end{cases}
$$

Then

$$
L(\boldsymbol{x}; \boldsymbol{\lambda}, \boldsymbol{\nu}) := \left(f(\boldsymbol{x}) + \sum_{i=1}^n \lambda_i g_i(\boldsymbol{x}) + \nu h(\boldsymbol{x})\right).
$$

Then the KKT condition implies

$$
\begin{cases}\nx_i \ge 0, & \lambda_i \ge 0, \quad \lambda_i x_i = 0 \\
\sum_{i=1}^n x_i = 1 \\
\frac{\partial L}{\partial x_k}(\boldsymbol{x}; \boldsymbol{\lambda}, \boldsymbol{\nu}) = \left(1 - \sum_{j=1}^{n'} q_{k,j} \log \frac{q_{k,j}}{\sum_{i'=1}^n x_{i'} q_{i',j}}\right) - \lambda_k + \nu = 0 \quad k = 1, \dots, n\n\end{cases}
$$

which further implies that (we can choose)

$$
\lambda_k = \begin{cases}\n1 - \sum_{j=1}^{n'} q_{k,j} \log \frac{q_{k,j}}{\sum_{i'=1}^{n} x_{i'} q_{i',j}} + \nu = 0 & x_k > 0 \\
1 - \sum_{j=1}^{n'} q_{k,j} \log \frac{q_{k,j}}{\sum_{i'=1}^{n} x_{i'} q_{i',j}} + \nu \ge 0 & x_k = 0\n\end{cases}
$$

By this, the input distributions that achieve the channel capacities of some channels such as BSC and BEC can be identified.  $\Box$ 

# Key Notes I: b-52

- Definitions of (weakly, strictly) stationarity, ergodicity and Markovian (irreducible, homogeneous)
- Mode of convergences (almost surely or with probability 1, in probability, in distribution, in  $L_r$  mean)
- Laws of large numbers
- Central limit theorem
- Jensen's inequality (convexity and concavity)
- KKT conditions