ML Preparation Notes

April 26, 2021

## **Basic Statistics**

In this section, we review some basic concepts from Statistics. The material in this section is based on [4].

#### 1.1 Covariance and Correlation

The correlation between two sets of data is a measure of the strength of the relationship between them. In particular, Pearson's correlation coefficient is a measure of linear relationship between two sets of data. Let X and Y be two random variables. Then Pearson's correlation coefficient  $\rho(X,Y)$  is defined as:

$$\rho(X,Y) = \frac{\text{Cov}(X,Y)}{\sigma_X \, \sigma_Y} \tag{1.1}$$

Two important facts about the Pearson's correlation coefficient [3]:

- 1.  $-1 \le \rho(X, Y) \le 1$
- 2.  $|\rho(X,Y)| = 1$  iff there exists  $a \neq 0$  and b such that Y = aX + b.

#### 1.2 Tail Inequalities and the Law of Large Numbers

We first look at Markov and Chebyshev inequalities and then apply them to random samples.

The Markov inequality tells us how much probability mass can be at large values given the mean of the distribution.

**Theorem 1.1.** Let X be a random variable that takes on positive values only. Then for t > 0,

$$P(X \ge t) \le \frac{E[X]}{t}.$$
(1.2)

*Proof.* Let us assume that X is a continuous r.v. with pdf  $f_X$ . By definition,

$$E[X] = \int_0^\infty x f_X(x) dx$$
$$= \int_0^t x f_X(x) dx + \int_t^\infty x f_X(x) dx$$
$$\ge t \int_t^\infty f_X(x) dx.$$

Since t > 0, dividing both sides of the last inequality yields the Markov inequality.

This inequality is useful when t > E[X]; when  $t \le E[X]$ , it merely bounds  $P(X \ge t)$  by 1.

The Chebyschev inequality involves both the mean and the variance of the distribution. It bounds the probability of how far a random variable can be from its mean as a function of the variance.

**Theorem 1.2.** Let X be a random variable for which Var(X) exists. Then

$$P(|X - E[X]| > t) \le \frac{Var(X)}{t^2}.$$
(1.3)

*Proof.* Define  $Y := (X - E[X])^2$  so that E[Y] = Var(X). Applying the Markov inequality to the r.v. Y, we obtain:

$$P(Y \ge t^2) \le \frac{Var(X)}{t^2}.$$

But  $P(Y \ge t^2) = P(|X - E[X]| > t)$  and so this proves the Chebyschev inequality too.

To talk about the law of large numbers, one has to talk about the notion of a sequence of random variables converging to a real number. We say that a sequence  $Z_1, Z_2, \ldots$  of random variables converges to the number b if the distribution of  $Z_n$  as  $n \to \infty$  becomes more and more concentrated around this single number.

**Definition 1.1** (Convergence in Probability). A sequence  $Z_1, Z_2, ...$  of random variables converges in probability to the number b if for every  $\epsilon > 0$ , the following condition holds:

$$\lim_{n \to \infty} P(|Z_n - b| < \epsilon) = 1. \tag{1.4}$$

This fact is denoted by  $Z_n \stackrel{p}{\longrightarrow} b$ .

**Theorem 1.3** (The Law of Large Numbers). Let  $X_1, \ldots, X_n$  be a random sample from a distribution with mean  $\mu$  and finite variance  $\sigma^2$ . Then the sample mean  $\bar{X}_n$  converges in probability to  $\mu$ .

*Proof.* We have  $E[\bar{X}_n] = \mu$  and  $Var(\bar{X}_n) = \sigma^2/n$ . Use the Chebyschev inequality to obtain:

$$P(|\bar{X}_n - \mu| \le t) \ge 1 - \frac{\sigma^2}{nt^2}.$$

For every fixed t > 0, take limits as  $n \to \infty$  to obtain:  $\lim_{n \to \infty} P(|\bar{X}_n - \mu| \le t) = 1$ . Hence  $\bar{X} \stackrel{p}{\longrightarrow} \mu$ .  $\square$ 

If we know that a sequence of random variables converges in probability, what can we say about continuous functions of that sequence of random variables?

**Theorem 1.4** (Continuous Functions of Random Variables). Suppose a sequence  $Z_1, Z_2, \ldots$  converges in probability to b and that g is a function continuous at b. Then  $g(Z_n) \stackrel{p}{\longrightarrow} g(b)$ .

Proof. We have to show  $\lim_{n\to\infty} P(|g(Z_n)-g(b)|<\epsilon)=1$ . Since g is continuous at b, given  $\epsilon>0$  there exists  $\delta>0$  such that  $|z_n-b|<\delta$  implies  $|g(z_n)-g(b)|<\epsilon$ . Now consider the events  $\mathcal{E}_1=\{|Z_n-b|<\delta\}$  and  $\mathcal{E}_2=\{|g(Z_n)-g(b)|<\epsilon\}$ . By the continuity of g,  $\mathcal{E}_1$  implies  $\mathcal{E}_2$ , and we have that  $P(\mathcal{E}_2)\geq P(\mathcal{E}_1)$ . Since  $\lim_{n\to\infty} P(\mathcal{E}_1)=1$ , we must have  $\lim_{n\to\infty} P(\mathcal{E}_2)=1$ .

**Theorem 1.5** (Chernoff Bounds). Let X be a random variable with moment generating function  $\psi_X$ . Then for any real t,

$$P(X \ge t) \le \min_{s>0} e^{-st} \psi_X(s). \tag{1.5}$$

*Proof.* Fix s > 0. Since the map  $x \longrightarrow e^x$  is 1-1 and increasing, we have that  $X \ge t$  if and only if  $e^{sX} \ge e^{st}$ . Therefore by the Markov inequality, we obtain:

$$P(X \ge t) = P(e^{sX} \ge e^{st}) \le \frac{E[e^{sX}]}{e^{st}} = e^{-st}\psi_X(s).$$

This holds for any s > 0 and, in particular, when we minimize  $e^{-st}\psi_X(s)$  subject to the condition that s > 0.

**Example 1.1.** For each positive integer n, let  $X_n$  be a nonnegative random variable with finite mean  $\mu_n$ . Show that if  $\lim_{n\to\infty}\mu_n=0$ , then  $X_n\stackrel{p}{\longrightarrow}0$ .

Solution. We have to show that for every  $\epsilon > 0$ ,  $\lim_{n \to \infty} P(X_n \ge \epsilon) = 0$ . To this end, fix  $\epsilon > 0$  and  $\delta > 0$ . Since  $\lim_{n \to \infty} \mu_n = 0$ , there exists  $n_{\epsilon \delta} \in \mathbb{N}$  such that for all  $n \ge n_{\epsilon \delta}$ ,  $\mu_n < \epsilon \cdot \delta$ . Combining this with the Markov inequality, we obtain that for all  $n \ge n_{\epsilon \delta}$ 

$$P(X_n \ge \epsilon) \le \frac{\mu_n}{\epsilon} < \delta.$$

This is precisely what we need to show in order to prove that  $\lim_{n\to\infty} P(X_n \ge \epsilon) = 0$ .

**Example 1.2.** Let  $Z_1, Z_2, ...$  be a sequence of random variables, and suppose that, for n = 1, 2, ..., the distribution of  $Z_n$  is as follows:

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$$P(Z_n = n^2) = \frac{1}{n}$$
 and  $P(Z_n = 0) = 1 - \frac{1}{n}$ .

Show that  $\lim_{n\to\infty} E[Z_n] = \infty$  but  $Z_n \stackrel{p}{\longrightarrow} 0$ .

Solution. Now  $\mathrm{E}[Z_n] = n^2 \cdot \frac{1}{n} = n$  and clearly  $\mathrm{E}[Z_n] \to \infty$  as  $n \to \infty$ . Since  $\lim_{n \to \infty} \mathrm{P}(Z_n = 0) = 1$ , it follows that for any  $\epsilon > 0$ ,  $\lim_{n \to \infty} \mathrm{P}(Z_n > \epsilon) = 0$ . This shows that  $Z_n \xrightarrow{p} 0$ .

#### 1.3 The Central Limit Theorem

The CLT is a formal statement of how normal distributions can approximate distributions of general sums or averages of iid random variables. To start with, consider the case when  $X_1, \ldots, X_n$  are iid random variables with  $X_i \sim N(\mu_i, \sigma_i^2)$  for  $1 \le i \le n$ . Let  $Y = a_1 X_1 + \cdots + a_n X_n + b$ , where  $a_i \ne 0$  for at least one  $1 \le i \le n$ . Then Y is normally distributed as the following shows.

**Theorem 1.6.** Let  $X_1, \ldots, X_n$  be iid normally distributed random variables with  $X_i$  having mean  $\mu_i$  and variance  $\sigma_i^2$ , for  $1 \le i \le n$ . Define  $Y = a_1X_1 + \cdots + a_nX_n + b$ , where  $a_i \ne 0$  for at least one  $1 \le i \le n$ . Then Y is normally distributed with mean  $a_1\mu_1 + \cdots + a_n\mu_n + b$  and variance  $a_1^2\mu_1^2 + \cdots + a_n^2\mu_n^2$ .

*Proof.* Let  $\psi_i(t)$  be the mgf of  $X_i$  for  $1 \le i \le n$ , and let  $\psi(t)$  denote the mgf of  $a_1X_1 + \cdots + a_nX_n + b$ . Then

$$\psi(t) = \mathbf{E}[e^{t(a_1X_1 + \dots + a_nX_n + b)}]$$

$$= e^{tb} \cdot \prod_{i=1}^n \mathbf{E}[e^{a_iX_it}]$$

$$= e^{tb} \cdot \prod_{i=1}^n \psi_i(a_it)$$

$$= e^{tb} \cdot \prod_{i=1}^n \exp\left(a_i\mu_it + \frac{1}{2}a_i^2\sigma_i^2t^2\right)$$

$$= \exp\left(\left(b + \sum_i \mu_i\right)t + \frac{1}{2}\left(\sum_i a_i^2\sigma_i^2\right)t^2\right).$$

Note that  $\psi(t)$  is the mgf of a normal distribution with mean  $b + \sum_i \mu_i$  and variance  $\sum_i a_i^2 \sigma_i^2$ . Hence  $Y \sim N(b + \sum_i \mu_i, \sum_i a_i^2 \sigma_i^2)$ .

**Theorem 1.7** (Central Limit Theorem for IID RVs). Let  $X_1, \ldots, X_n$  be iid random variables from a distribution with mean  $\mu$  and variance  $\sigma^2 < \infty$ . Then for each fixed number x,

$$\lim_{n \to \infty} P\left(\frac{\bar{X}_n - \mu}{\sigma/\sqrt{n}} \le x\right) = \Phi(x), \tag{1.6}$$

where  $\Phi$  denotes the cdf of the standard normal distribution.

We can interpret Equation 1.6 as follows: If we have a large random sample  $X_1, \ldots, X_n$  from an arbitrary distribution (whether discrete or continuous), the random variable  $(\bar{X}_n - \mu)/(\sigma/\sqrt{n})$  is distributed approximately as the standard normal. Consequently,  $\bar{X}_n$  is distributed approximately as  $N(\mu, \sigma^2/n)$  and  $\sum_{i=1}^n X_i$  is distributed approximately as  $N(n\mu, n\sigma^2)$ .

#### 1.4 Maximum Likelihood Estimation

Let the random variables  $X_1, \ldots, X_n$  form a random sample from a distribution with pdf  $f(x \mid \theta)$ . Recall that this means that  $X_i \stackrel{\text{iid}}{\sim} f(\cdot \mid \theta)$  for all  $1 \leq i \leq n$ . Let  $f_n(\vec{x} \mid \theta)$  denote the value of the joint pdf of the random vector  $(X_1, \ldots, X_n)'$  at the point  $\vec{x} = (x_1, \ldots, x_n)'$ . The *likelihood function* is the joint pdf of the observations of a random sample viewed as a function of  $\theta$  for a given set of values of the sample. The maximum likelihood estimate of  $\theta$  is that value of  $\theta$  for which  $f_n(\vec{x} \mid \theta)$  is maximized.

**Example 1.3.** Suppose that  $X_1, \ldots, X_n$  form a random sample from a distribution with pdf  $f(x \mid \theta)$  defined as follows:

$$f(x \mid \theta) = \begin{cases} \theta x^{\theta - 1} & \text{for } 0 < x < 1 \\ 0 & \text{otherwise.} \end{cases}$$

Let us assume that  $\theta > 0$ . The joint distribution of the vector  $(X_1, \ldots, X_n)'$  at the point  $\vec{x} = (x_1, \ldots, x_n)'$  is

$$f_n(\vec{x} \mid \theta) = \prod_{i=1}^n \theta x_i^{\theta-1},$$

where we assume that  $0 < x_i < 1$  for all *i*. Taking logs, we obtain that  $\log f_n(\vec{x} \mid \theta) = n \log \theta + (\theta - 1) \sum_i \log x_i$ . Take the derivative wrt  $\theta$  and set to 0:

$$\frac{\partial \log f_n(\vec{x} \mid \theta)}{\partial \theta} = \frac{n}{\theta} + \sum_i \log x_i \stackrel{\text{set}}{=} 0$$

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to obtain that  $\theta = n / \sum_{i} \log \frac{1}{x_i}$ .

**Example 1.4.** Suppose that  $X_1, \ldots, X_n$  form a random sample from a distribution with pdf  $f(x \mid \theta)$  defined as follows:

 $f(x \mid \theta) = \frac{1}{2}e^{-|x-\theta|}$  for  $-\infty < x < \infty$ .

Suppose that  $\theta$  is unknown and that  $-\infty < \theta < \infty$ . In this case, the joint probability distribution is easily seen to be:

$$f_n(\vec{x} \mid \theta) = \frac{1}{2^n} e^{-\sum_i |x_i - \theta|}.$$

Take logs to obtain:  $\log f_n(\vec{x} \mid \theta) = \log \frac{1}{2^n} - \sum_i |x_i - \theta|$ . Maximizing  $f_n$  is equivalent to minimizing  $\sum_i |x_i - \theta|$ . This is equivalent to obtaining a point on the real line that minimizes the sum of the distances to the points  $x_1, \ldots, x_n$ . This happens when  $\theta$  is the median of  $x_1, \ldots, x_n$ .

**Example 1.5.** Suppose that  $X_1, \ldots, X_n$  form a random sample from the uniform distribution on the interval  $[\theta_1, \theta_2]$ , where both  $\theta_1$  and  $\theta_2$  are unknown  $(-\infty < \theta_1 < \theta_2 < \infty)$ . In this case, the log pdf of the joint distribution of  $(X_1, \ldots, X_n)'$  is given by

$$\log f_n(\vec{x} \mid \theta_1, \theta_2) = \log \prod_{i=1}^n \frac{1}{\theta_2 - \theta_1} = -n \log(\theta_2 - \theta_1).$$

Maximizing the likelihood is equivalent to minimizing  $\log(\theta_2 - \theta_1)$ . The minimum possible value of  $\theta_2$  is  $\max\{x_1, \ldots, x_n\}$  and the maximum possible value of  $\theta_1$  is  $\min\{x_1, \ldots, x_n\}$ .

**Example 1.6.** Suppose that a certain large population contains k different types of individuals  $(k \ge 2)$ , and let  $\theta_i$  denote the proportion of people of type i, for  $1 \le i \le k$ . Here,  $0 \le \theta_i \le 1$  and  $\sum_{i=1}^k \theta_i = 1$ . Suppose also that in a random sample of n individuals from this population there are exactly  $n_i$  individuals of type i so that  $n = n_1 + \dots + n_k$ .

In this setting, for  $1 \le i \le k$ , define  $X_i$  to be the number of individuals of type i in a random sample of size n. Then the probability that  $\bigwedge_{i=1}^k X_i = n_i$  is given by  $\theta_1^{n_1} \cdots \theta_k^{n_k}$ . The log pdf of the joint distribution is given by:

$$\log f_k((n_1,\ldots,n_k)\mid \theta_1,\ldots,\theta_k,n) = \sum_{i=1}^k n_i \log \theta_i.$$

Note that there are actually k-1 variables here since we may write  $\theta_k = 1 - \sum_{i=1}^{k-1} \theta_i$ . Differentiating wrt  $\theta_i$  for  $1 \le i \le k-1$ , we obtain:

$$\frac{\partial \log f_k}{\partial \theta_i} = \frac{n_i}{\theta_i} - \frac{n_k}{\theta_k}.$$

Setting this to 0, we get that  $\theta_i/\theta_k = n_i/n_k$ . Sum this up from  $1 \le i \le k-1$ , to obtain:

$$\frac{1-\theta_k}{\theta_k} = \frac{n-n_k}{n_k},$$

which yields  $\theta_k = n_k/n$ . Substitute this in  $\theta_i/\theta_k = n_i/n_k$  to obtain  $\theta_i = n_i/n$ .

**Example 1.7** (Nonexistence of an MLE). An obvious disadvantage of the technique of maximum likelihood estimation is when the maximum does not exist. Consider again Example 1.5 where we let  $\theta_1 = 0$  and  $\theta = \theta_2$  The pdf of the uniform distribution is defined as:

$$f(x \mid \theta) = \begin{cases} \frac{1}{\theta} & 0 \le x \le \theta \\ 0 & \text{otherwise.} \end{cases}$$

Let us modify the definition of the pdf so that we use strict inequalities  $0 < x < \theta$  above. Given a sample  $x_1, \ldots, x_n$ , the log pdf is  $-n \log \theta$  as before and the MLE technique would require us to minimize  $\log \theta$ . However, in this case, there is no  $\theta > \max\{x_1, \ldots, x_n\}$  that minimizes  $\log \theta$  and the MLE does not exist.

**Example 1.8** (Non-uniqueness of an MLE). Consider a random sample  $X_1, \ldots, X_n$  from the uniform distribution over the interval  $[\theta, \theta + 1]$ . In this case, the joint pdf is given by:

$$f_n(\vec{x} \mid \theta) = \begin{cases} 1 & \text{for } \theta \le x_i \le \theta + 1 & (1 \le i \le n) \\ 0 & \text{otherwise.} \end{cases}$$

In this case, the condition  $\theta \leq x_i \leq \theta + 1$  for  $1 \leq i \leq n$  may be written using the two conditions:

$$\theta \le \min\{x_1, \dots, x_n\}$$
 and  $\max\{x_1, \dots, x_n\} - 1 \le \theta$ .

Any value of  $\theta$  in the interval  $[\max\{x_1,\ldots,x_n\}-1,\min\{x_1,\ldots,x_n\}]$  is valid, but there is no unique value of  $\theta$ .

#### 1.5 Bayesian Statistics and MCMC

This section is based on Chapters 12–15 from [6]. Bayes' rule gives us a recipe for calculating the posterior probability density.

$$P(\Theta \mid data) = \frac{P(data \mid \Theta) \cdot P(\Theta)}{P(data)}.$$
(1.7)

Consider a case in which we have a sample of N data points  $x_1, \ldots, x_N$ . We assume that the likelihood is a Poisson distribution with mean  $\lambda$  and that the prior for  $\lambda$  is a log-normal(1,1) distribution. To calculate the probability of the data P(data), we must evaluate the integral:

$$P(\text{data}) = \int_0^\infty \prod_{i=1}^N \frac{\lambda^{x_i} e^{-\lambda}}{x_i!} \cdot \frac{1}{\sqrt{2\pi}\lambda} e^{-\frac{1}{2}(\log \lambda - 1)^2} d\lambda.$$
 (1.8)

While this integral is not too difficult, it explains the problem of calculating posteriors analytically. As the number of parameters (the length of  $\Theta$ ) increases, calculating the probability of the data requires evaluating integrals in higher dimensional spaces. This is why we use alternative methods to derive approximate versions of the posterior.

In certain special cases, the posterior distribution can be easily derived. This occurs, for example, when the prior distribution is from a so-called *conjugate prior* family. A conjugate prior family is a set of distributions defined wrt a likelihood function. If one chooses the prior to be from this family, then the posterior is guaranteed to be from the same family. In such cases, one can dispense with computing integrals to find out the posterior.

**Example 1.9** (Beta-Binomial). Let  $X_1, \ldots, X_n \stackrel{\text{iid}}{\sim} \text{Bernoulli}(\theta)$  and suppose that  $\theta \sim \text{Beta}(\alpha, \beta)$ . Suppose that one observes  $X_i = x_i$  for  $1 \leq i \leq n$ . Then the posterior distribution of  $\theta$  is given by:

$$f(\theta \mid X_1, \dots, X_n) = \frac{f(X_1, \dots, X_n \mid \theta) \cdot f(\theta)}{f(X_1, \dots, X_n)}.$$

In the above expression,  $f(X_1, ..., X_n)$  is a constant for a given set of values of the random variables  $X_i$ . The likelihood, when  $X_i = x_i$  for  $1 \le i \le n$ , is given by:

$$f(x_1, \dots, x_n \mid \theta) = \prod_{i=1}^n \theta^{x_i} (1 - \theta)^{1 - x_i}$$
$$= \theta^{\sum_i x_i} \cdot (1 - \theta)^{n - \sum_i x_i}.$$

Now we may write:

$$f(\theta \mid X_1, \dots, X_n) = \frac{f(X_1, \dots, X_n \mid \theta) \cdot f(\theta)}{f(X_1, \dots, X_n)}$$
$$\propto \theta^{\sum_i x_i} \cdot (1 - \theta)^{n - \sum_i x_i} \cdot \theta^{\alpha - 1} \cdot (1 - \theta)^{\beta - 1}$$
$$\propto \theta^{\alpha + \sum_i x_i - 1} \cdot (1 - \theta)^{\beta + n - \sum_i x_i - 1}.$$

The last expression is the kernel of the beta distribution with parameters  $\alpha + \sum_i x_i$  and  $\beta + n - \sum_i x_i$ . Since the expression on the right must be a valid probability distribution, the proportionality constant must be that of the beta distribution with these parameters, that is,

$$\frac{\Gamma(\alpha+\beta+n)}{\Gamma(\alpha+\sum_{i}x_{i})\cdot\Gamma(\beta+n-\sum_{i}x_{i})}.$$

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Thus if the prior is a beta distribution and the likelihood is a bernoulli distribution, then the posterior is a beta distribution also. Thus the set of beta distributions is a conjugate family wrt the bernoulli likelihood.

#### **Timeseries**

The material in the section follows [2]. A time series is a sequence of random variables  $\{X_1, X_2, \ldots\}$ . A complete model of such a time series would require specifying the joint probability distributions of all random vectors  $(X_1, \ldots, X_n)'$  for all  $n \geq 1$ . In practice, this might be impossible to do unless the time series is generated by some simple well-understood mechanism. Instead one typically specifies the first-and second-order moments of the joint distributions, that is,  $E[X_t]$  and  $E[X_tX_{t+h}]$  for all  $t \geq 1$  and for all  $h \geq 0$ .

The next important concepts are those of stationarity and the auto-correlation function. Roughly speaking, a time series  $\{X_t\}_{t=-\infty}^{\infty}$  is stationary if its "statistical properties" are similar to those of the time-shifted series  $\{X_{t+h}\}_{t=-\infty}^{\infty}$  for every integer "lag" h. By statistical properties, we mean the first- and second-order moments of  $\{X_t\}$ .

Formally, a time series  $\{X_t\}$  is weakly stationary if

- 1.  $E[X_t]$  is independent of t
- 2.  $Cov(X_t, X_{t+h})$  is independent of t for every fixed lag h.

In contrast, a time series  $\{X_t\}$  is strictly stationary if the random vectors  $(X_{t_1}, \ldots, X_{t_n})'$  and  $(X_{t_1+h}, \ldots, X_{t_n+h})'$  have the same joint distributions for all sets of indices  $\{t_1, \ldots, t_n\}$ , for all  $h \geq 0$  and all  $n \geq 1$ . This is written as:

$$(X_{t_1}, \dots, X_{t_n})' \stackrel{d}{=} (X_{t_1+h}, \dots, X_{t_n+h})'$$

Strict stationarity implies the following:

- 1. The random variables  $X_t$  are identically distributed.
- 2. Pairs of random variables  $(X_t, X_{t+h})$  have the same distribution as  $(X_1, X_{1+h})$  (set n = 2).
- 3. Strict stationarity implies weak stationarity:  $E[X_t] = E[X_1]$  and  $Cov(X_t, X_{t+h}) = Cov(X_1, X_{1+h})$  for all  $t \ge 1$  and all  $h \ge 0$ . Both terms are independent of t.
- 4. Weak stationarity does not imply strong stationarity. We show this by an example. Let  $Z_i \stackrel{\text{iid}}{\sim} N(0,1)$  for all i. Define  $X_t$  as:

$$X_t = \begin{cases} Z_t & \text{if } t \text{ is even} \\ 2Z_t & \text{if } t \text{ is odd.} \end{cases}$$

Then  $E[X_t] = 0$  for all t. Also,

$$Cov(X_t, X_{t+h}) = \begin{cases} 0 & \text{if } h > 0\\ 1 & \text{if } h = 0. \end{cases}$$

This follows from the assumed independence of  $X_t$  and  $X_{t+h}$  when h > 0. However,  $X_t$ s do not have the same distribution, a requirement of strict stationarity.

The autocovariance function of a stationary time series  $\{X_t\}$  at lag h is defined as  $Cov(X_{t+h}, X_t) = Cov(X_h, X_0)$ . The autocorrelation of  $\{X_t\}$  at lag h is defined as

$$\frac{\operatorname{Cov}(X_{t+h}, X_t)}{\operatorname{Var}(X_t)} = \frac{\operatorname{Cov}(X_h, X_0)}{\operatorname{Var}(X_0)}.$$

**Example 2.1** (iid Noise). Suppose  $\{X_t\}$  is iid noise with zero mean and  $\mathrm{E}[X_t^2] = \sigma^2 < \infty$ . Then  $\mathrm{E}[X_t] = 0$  for all t and for every fixed lag h and all t:

$$Cov(X_t, X_{t+h}) = \begin{cases} 0 & \text{if } h > 0 \\ \sigma^2 & \text{if } h = 0. \end{cases}$$

Thus iid noise is stationary. We denote such a series as  $\{X_t\} \sim \text{IID}(0, \sigma^2)$ .

**Example 2.2** (White Noise). A sequence  $\{X_t\}$  of uncorrelated random variables with zero mean and finite second moment  $\sigma^2$  is called white noise. Clearly, the covariance function at lag h is the same as that of iid noise and, as such, white noise is stationary. We denote such a series as  $\{X_t\} \sim \text{WN}(0, \sigma^2)$ . Unlike iid noise, the components of white noise need not be independent (recall that independence implies zero correlation but not the other way around). In particular, every  $\text{IID}(0, \sigma^2)$  sequence is a  $\text{WN}(0, \sigma^2)$  sequence but not the other way around.

**Example 2.3** (Random Walk). A random walk  $\{S_t\}$  is a sequence obtained by cumulatively summing iid random variables. A random walk with zero mean is obtained by defining  $S_0 = 0$  and

$$S_t = X_1 + \cdots + X_t$$

for all t > 0, where  $\{X_t\} \sim \text{IID}(0, \sigma^2)$ . In this case,  $\text{E}[S_t] = 0$  and  $\text{E}[S_t^2] = t\sigma^2 < \infty$  for all t. For all lags  $h \ge 0$  and all t,

$$Cov(S_t, S_{t+h}) = Cov(S_t, S_t + X_{t+1} + \dots + X_{t+h})$$
  
=  $Cov(S_t, S_t) + Cov(S_t, X_{t+1}) + \dots + Cov(S_t, X_{t+h})$   
=  $t\sigma^2$ .

The last equality follows since  $Cov(S_t, X_{t+i}) = 0$  for all  $i \ge 1$ . Hence  $Cov(S_t, S_{t+h})$  depends on t and  $\{S_t\}$  is not stationary.

**Example 2.4** (First-Order Moving Average). Consider the series defined by

$$X_t = Z_t + \theta Z_{t-1}, \quad t = 0, \pm 1, \pm 2, \dots,$$
 (2.1)

where  $Z_t \sim \text{WN}(0, \sigma^2)$  and  $\theta$  is a real-valued constant. Now  $\text{E}[X_t] = 0$  and

$$E[X_t^2] = E[(Z_t + \theta Z_{t-1})^2]$$

$$= E[Z_t^2 + 2\theta Z_t Z_{t-1} + \theta^2 Z_{t-1}^2]$$

$$= \sigma^2 + 2\theta E[Z_t Z_{t-1}] + \theta^2 \sigma^2$$

$$= (1 + \theta^2)\sigma^2.$$

The last equality follows since,  $Z_t$  and  $Z_{t-1}$  being uncorrelated, satisfy  $Cov(Z_t, Z_{t-1}) = 0$ . Recall that  $Cov(Z_t, Z_{t-1}) = E[Z_t Z_{t-1}] - E[Z_t] E[Z_{t-1}]$  and that being uncorrelated is sufficient for the expectation

of the product of two random variables to be equal to the product of their expectations. One can easily verify that for all t

$$\operatorname{Cov}(X_t, X_{t+h}) = \begin{cases} (1+\theta^2)\sigma^2 & \text{if } h = 0\\ \theta\sigma^2 & \text{if } h = \pm 1\\ 0 & \text{if } |h| \ge 2. \end{cases}$$
 (2.2)

Thus the conditions of weak stationarity hold and the sequence  $\{X_t\}$  is weakly stationary. The autocorrelation function is given by:

$$\rho(h) = \frac{\text{Cov}(X_0, X_h)}{\text{Var}(X_0)} = \begin{cases} 1 & \text{if } h = 0\\ \frac{\theta}{1 + \theta^2} & \text{if } h = \pm 1\\ 0 & \text{if } |h| \ge 2. \end{cases}$$
 (2.3)

**Example 2.5** (First-Order Autoregression). Let  $\{X_t\}$  be a stationary series satisfying the equation:

$$X_t = \phi X_{t-1} + Z_t \quad t = 0, \pm 1, \pm 2, \dots,$$
 (2.4)

where  $\{Z_t\} \sim WN(0, \sigma^2)$ ,  $|\phi| < 1$  and  $Z_t$  is uncorrelated with  $X_s$  for all s < t. As can be seen,  $E[X_t] = 0$  and for h > 0:

$$Cov(X_{t+h}, X_t) = Cov(\phi X_{t+h-1} + Z_{t+h}, X_t)$$
  
=  $\phi Cov(X_{t+h-1}, X_t) + Cov(Z_{t+h}, X_t)$   
=  $\phi Cov(X_{t+h-1}, X_t)$ .

From this, one can show that  $Cov(X_{t+h}, X_t) = \phi^h Cov(X_t, X_t)$ . By assumption,  $\{X_t\}$  is stationary and hence  $Cov(X_{t+h}, X_t) = \phi^h Cov(X_0, X_0)$  Next suppose that h < 0. Let s = t + h = t - |h| so that t = s + |h|. Then  $Cov(X_{t+h}, X_t) = Cov(X_s, X_{s+|h|}) = \phi^{|h|} Cov(X_s, X_s) = \phi^{|h|} Cov(X_0, X_0)$ . The autocorrelation function at lag h is

$$\rho(h) = \frac{\phi^{|h|} \operatorname{Cov}(X_0, X_0)}{\operatorname{Cov}(X_0, X_0)} = \phi^{|h|}.$$

One can also obtain a closed-form expression for  $Cov(X_0, X_0)$ . Since  $Cov(X_0, X_0) = Cov(X_t, X_t)$  and

$$Cov(X_t, X_t) = Cov(\phi X_{t-1} + Z_t, \phi X_{t-1} + Z_t) = \phi^2 Cov(X_0, X_0) + Cov(Z_t, Z_t),$$

we obtain that  $Cov(X_0, X_0) = \sigma^2/(1 - \phi^2)$ .

Trees, Boosting and Random Forests

## Gradient Boosted Regression

We first consider gradient boosting for classification problems. Suppose that we are given a dataset  $\{(\mathbf{x}_i, y_i)_{i=1}^n\}$ , where  $\mathbf{x}_i \in \mathbf{R}^p$  and  $y_i \in \{0, 1\}$ . The problem is to find a mapping  $f : \mathbf{x} \to y$ . As in logistic regression, we modify the problem slightly and do not work directly with the class labels  $y_i$ . Instead, we consider the log(odds) of the event  $P\{y = 1 \mid \mathbf{x}\}$ . Since  $-\infty < \log(\text{odds}) < +\infty$ , this restatement allows us to focus on functions  $F : \mathbf{R}^p \to \mathbf{R}$  rather than from  $\mathbf{R}^p \to \{0, 1\}$ . This is helpful because gradient boosting requires differentiable loss functions.

Our next step is to build an appropriate loss function. Let  $p = P\{y = 1 \mid \mathbf{x}\}$ . Then we may write

$$P\{y \mid \mathbf{x}\} = p^y \cdot (1-p)^{1-y}.$$

Consequently, the likelihood of  $y_1, \ldots, y_n$  given  $\mathbf{x}_1, \ldots, \mathbf{x}_n$  assuming that the data  $\{(\mathbf{x}_i, y_i)_{i=1}^n\}$  are independent is

$$P\{y_1, \dots, y_n \mid \mathbf{x}_1, \dots, \mathbf{x}_n\} = \prod_{i=1}^n p_i^{y_i} \cdot (1 - p_i)^{1 - y_i}.$$

The log-likelihood is  $\sum_{i} [y_i \log p_i + (1 - y_i) \log (1 - p_i)]$ . If we were to fit a model such a logistic regression, we would search for those model parameters for which the log-likelihood is a maximum. If we were to use a transform of the log-likelihood as a loss function, we would want to minimize that transform. The easiest transform is the negative log-likelihood. Minimizing the negative log-likelihood is equivalent to maximizing the log-likelihood (which is what we want to do). This negative log-likelihood will then be our loss function and this also goes by the name of cross entropy.

We re-write the loss function in terms of  $\log(\text{odds})$  because this is what our gradient boosting model will output. To simplify the resulting expressions, we consider only one data point and omit the data index i. We may then write:

$$\begin{split} -\left[y\log p + (1-y)\log(1-p)\right] &= -y\log p - (1-y)\log(1-p) \\ &= -y\log p + y\log(1-p) - \log(1-p) \\ &= -y\log\frac{p}{1-p} - \log\left(1-\sigma\left(\log\frac{p}{1-p}\right)\right). \end{split}$$

In the last step, we used the fact that  $p = \sigma(p/(1-p))$ , where  $\sigma$  is the sigmoid function. This last step can be simplified by expanding out the sigmoid function and the loss function can then be written as:

$$-y\log\frac{p}{1-p} + \log(1 + e^{\log\frac{p}{1-p}}).$$

Note the positive sign before the logarithm.

Finally, we note that the gradient boosting engine gives us a function  $F(\mathbf{x})$  that represents the  $\log(\text{odds})$ . With this, we can write the loss function more clearly as:

$$L(y, F(\mathbf{x})) = -yF(\mathbf{x}) + \log(1 + e^{F(\mathbf{x})}).$$

This function is differentiable wrt  $F(\mathbf{x})$  and with some manipulation, one can show that:

$$\frac{\mathrm{d} L}{\mathrm{d} F(\mathbf{x})} = -y + \sigma(F(\mathbf{x})).$$

Now this has a nice interpretation. The term  $\sigma(F(\mathbf{x}))$  is the predicted probability that y = 1. If we were to interpret the label  $y \in \{0,1\}$  as a probability, then the derivative of the loss function wrt  $F(\mathbf{x})$  is the negative of the difference of the actual probability and the predicted probability.

# Neural Networks

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