

## Basic Stats

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# Preface: Motivation

All the notes I have done here are about basic stats. While I have tried my best, probably there are still some typos and errors. Please feel free to let me know in case you find one. Thank you!



# Chapter 1

## 443

### 1.1 Some basic concepts

#### 1.1.1 Permutation

An ordered arrangement of a set of objects is known as a permutation.

e.g., The number of permutations of  $n$  distinguishable objects is  $n!$ . e.g., The number of permutations of  $n$  distinct objects taken  $r$  at a time is

$${}_nP_r = \frac{n!}{(n-r)!}$$

#### 1.1.2 Combinations

If the order of the objects is not important, then one may simply be interested in the number of combinations.

$$\binom{n}{r} = \frac{n!}{r!(n-r)!}$$

#### 1.1.3 Partitioning

The number of ways of partitioning a set of  $n$  objects into  $k$  cells with  $r_1$  objects into the first cell,  $r_2$  in the second cell, and so forth is

$$\frac{n!}{r_1!r_2!\dots r_k!}$$

## 1.2 Discrete Random Variables

### 1.2.1 Binomial

$$X \sim \text{BIN}(n, p)$$

$$\binom{n}{x} P^x (1 - P)^{n-x}$$

mean:  $np$

variance:  $npq$

(Note that, Bernoulli is written as  $\text{BIN}(1, p)$ )

### 1.2.2 Poisson

$$X \sim \text{POI}(\mu)$$

$$\frac{e^{-\mu} \mu^x}{x!}$$

mean:  $\mu$

variance:  $\mu$

## 1.3 Continuous Random Variables

### 1.3.1 Uniform

$$X \sim \text{UNIF}(a, b)$$

$$\frac{1}{b - a}$$

Mean:  $\frac{a+b}{2}$

Variance:  $\frac{(b-a)^2}{12}$



### 1.3.2 Exponential

$$X \sim EXP(\theta)$$

$$\frac{1}{\theta} e^{-x/\theta}$$

Mean:  $\theta$

Variance:  $\theta^2$

### 1.3.3 Normal

$$X \sim N(\mu, \sigma^2)$$

$$\frac{1}{\sqrt{2\pi}\sigma} e^{-\frac{1}{2}\left(\frac{x-\mu}{\sigma}\right)^2}$$

Mean:  $\mu$

Variance:  $\sigma^2$

## 1.4 Large Sample Theory

### 1.4.1 Convergence in distribution

[https://en.wikipedia.org/wiki/Law\\_of\\_large\\_numbers](https://en.wikipedia.org/wiki/Law_of_large_numbers)

$$\bar{X} \rightarrow \mu \quad (n \rightarrow \infty)$$

$$Var(\bar{X}) = Var\left(\frac{1}{n}(X_1 + \dots + X_n)\right) = \frac{1}{n^2} Var(X_1 + \dots + X_n) = \frac{n\sigma^2}{n^2} = \frac{\sigma^2}{n}$$

### 1.4.2 Weak law

There are two different versions of the Law of Large Numbers: Strong law of large numbers and Weak law of large numbers.

The weak law of large numbers: The sample average converges in probability towards the expected value.

$$\bar{X}_n \xrightarrow{p} \mu \quad (n \rightarrow \infty)$$

This, for any positive number  $\epsilon$

$$\lim_{n \rightarrow \infty} Pr(|\bar{X}_n - \mu| > \epsilon) = 0$$

### 1.4.3 Strong law

$$\bar{X}_n \xrightarrow{a.s.} \mu \quad (n \rightarrow \infty)$$

This is,

$$Pr(\lim_{n \rightarrow \infty} \bar{X}_n = \mu) = 1$$

### 1.4.4 Central limit theorem

If  $X_1, \dots, X_n$  is a random sample from a distribution with mean  $\mu$  and variance  $\sigma^2 < \infty$ , then the limiting distribution of

$$Z_n = \frac{\sum_{i=1}^n X_i - n\mu}{\sqrt{n}\sigma}$$

is the standard normal,  $Z_n \xrightarrow{d} Z \sim N(0, 1)$  as  $n \rightarrow \infty$ .

#### 1.4.4.1 Bernoulli law of large number

$\hat{p}_n$  converges stochastically to  $p$  as  $n$  approaches infinity. For example, if a coin is tossed repeatedly, and  $A = \{H\}$ , then the successive relative frequencies of  $A$  correspond to a sequence of random variables that will converge stochastically to  $p = 1/2$ .

#### 1.4.4.2 Normal approximation to Binomial

$$Z_n = \frac{Y_n - np}{\sqrt{npq}} \xrightarrow{d} Z \sim N(0, 1)$$

Example: The probability that a basketball player hits a shot is  $p = 0.5$ . If he takes 20 shots, what is the probability that he hits at least 9?

$$\begin{aligned}
P[Y_{20} \geq 9] &= 1 - P[Y_{20} < 8] \\
&= 1 - \sum_{y=0}^8 \binom{20}{y} 0.5^y 0.5^{20-y} \\
&= 0.7483
\end{aligned}$$

A normal approximation is

$$\begin{aligned}
P[Y_{20} \geq 9] &= 1 - P[Y_{20} < 8] \\
&= 1 - \Phi\left(\frac{8-10}{\sqrt{5}}\right) \\
&= 0.8133
\end{aligned}$$

#### 1.4.4.3 Normal approximation to Poisson

$$\begin{aligned}
P[10 \leq Y_{20} \leq 30] &= P[Y_{20} \leq 30] - P[Y_{20} \leq 10] \\
&= \Phi\left[\frac{30.5-20}{\sqrt{20}}\right] - \Phi\left[\frac{9.5-20}{\sqrt{20}}\right] \\
&= 0.981
\end{aligned}$$

#### 1.4.5 Poisson approximation to binomial

We know that the mean for binomial is

$$\mu = np \rightarrow p = \frac{\mu}{n}$$

The moment generating function for Binomial is

$$\begin{aligned}
M_n(t) &= (1-p+pe^t)^n = \left(1 + \frac{\mu(e^t-1)}{n}\right)^n \\
\lim_{n \rightarrow \infty} M_n(t) &= e^{\mu(e^t-1)}
\end{aligned}$$

Note that the MGF for Poisson is as follows.

$$POI(\lambda) : e^{\lambda(e^t-1)}$$

Thus,

$$Y_n \rightarrow Y \sim POI(\mu)$$



# Chapter 2

## 556

### 2.1 Statistics and Sampling Distributions

#### 2.1.1 Statistics

##### 2.1.1.1 Definition of Statistic

*P.264*

A function of observable random variables,  $T = t(X_1, \dots, X_n)$ , which does not depend on any unknown parameters is called statistic.

For example, let  $X_1, \dots, X_n$  represent a random sample from a population with *pdf*  $f(x)$ . The sample mean provides an example of a statistic with the function

$$t(x_1, \dots, x_n) = (x_1 + \dots + x_n)/n$$

This statistic usually is denoted by

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

When a random sample is observed, the value of  $\bar{X}$ , computed from the data, usually is denoted by lower case  $\bar{x}$ .

$$\bar{x} = \sum_{i=1}^n \frac{x_i}{n}$$

### 2.1.1.2 Sample and parameters

*P.265*

If  $X_1, \dots, X_n$  denotes a random sample from  $f(x)$  with  $E(X) = \mu$  and  $var(X) = \sigma^2$ , then

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

$$Var(\bar{X}) = \frac{\sigma^2}{n}$$

For example, a random sample of size  $n$  from a Bernoulli distribution  $X_i \sim BIN(1, p)$ . We know Bernoulli has  $\mu = p$  and  $\sigma^2 = pq$ . In this case, the sample mean is

$$\bar{X} = Y/n = \hat{p}$$

Thus,

$$E(\hat{p}) = p$$

$$Var(\hat{p}) = \frac{pq}{n}$$

Thus, sample mean is the unbiased estimate for the population mean. However, you can not use sample mean's variance to estimate population variance. That lead to definition of sample variance.

*P.266*

Sample variance:

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

$$E(S^2) = \sigma^2$$

**2.1.2**  $\chi^2, t, F, \text{beta}$ **2.1.2.1**  $\chi^2$ 

Always square from standard normal, and the standardization can be using  $\mu$  or  $\bar{X}$ .

$$\frac{(n-1)S_n^2}{\sigma^2} \sim \chi^2(n-1)$$

(Thus, we can see this is a bit weird, as the numerator is  $\bar{X}$  is from the sample, whereas  $\sigma^2$  is from the population.)

Thus, we can

$$\frac{\sum_{i=1}^n (X_i - \bar{X})^2}{\sigma^2} \sim \chi^2(n-1)$$

(You can compare  $\bar{X}$  with  $\mu$ , we can see the only difference is that the  $\chi^2$  has one less degree of freedom because we use this degree of freedom to calculate the mean.)

For the mean and variance of  $\chi^2$ :

Assume that

$$X \sim \chi^2(v)$$

$$\text{mean} : v$$

$$\text{variance} : 2v$$

**2.1.2.2**  $t$ 

Definition

$$t(k) = \frac{N(0,1)}{\sqrt{\frac{\chi^2(k)}{k}}}$$

**Property 1**

$t$  distribution is symmetrical

Given that  $t$  distribution is symmetrical, we can get

$$H(-c) = 1 - H(c)$$

**Property 2**

$t$  distribution has heavier tails than the normal.

My note:  $t$  distribution only has a parameter of  $k$ , which is determined by the  $\chi^2$ 's degree of freedom. Of course,  $\chi^2$  also only has one parameter, namely the degree of freedom.

**2.1.2.3  $F$** 

If  $V_1 \sim \chi^2(v_1)$  and  $V_2 \sim \chi^2(v_2)$  are independent, then the random variable

$$\frac{V_1/v_1}{V_2/v_2} \sim F(v_1, v_2)$$

**2.1.2.4 Beta**

If  $X \sim F(v_1, v_2)$

$$Y = \frac{(v_1/v_2)X}{1 + (v_1/v_2)X} \sim \text{Beta}(\alpha, \beta)$$

**2.1.3 Large-sample approximations**

*P.280*

If  $Y_v \sim x^2(x)$ , then

$$Z_v = \frac{Y_v - v}{\sqrt{2v}} \xrightarrow{d} Z \sim N(0, 1)$$

(The proof is based on CLT.)

**2.2 Point Estimation****2.2.1 Method of moments estimators****2.2.1.1 Definition about moments (chapter 2)**

*P.73*

The  $k$ th moment about the origin of a random variable  $X$  is



$$\mu_k' = E(X^k)$$

and the **kth moment about the mean** is

$$\mu_k = E[X - E(X)]^k = E(X - \mu)^k$$

Thus,  $\mu_k = E(X^k)$  may be considered as the  $k$ th moment of  $X$  or the first moment of  $X^k$ .

The first moment about the mean is zero,

$$\mu_1 = E[X - E(X)] = E(X) - E(X) = 0$$

The second moment about the mean is the variance,

$$\mu_2 = E[(X - \mu)^2] = \sigma^2$$

Note that the definition of variance:

*P.73*

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

### 2.2.1.2 Definition

Based on the last chapter (i.e., Chapter 8), sample mean  $\bar{X}$  is an estimator of the population mean  $\mu$ . A more general approach, which produced estimators known as the **method of moments estimators (MMEs)**, can be developed.

If  $X_1, \dots, X_n$  is a random sample from  $f(x; \theta_1, \dots, \theta_k)$ , the first  $k$  sample moments are given by

$$M_j' = \frac{\sum_{i=1}^n X_i^j}{n}$$

where,

$$j = 1, 2, \dots, k$$

### Example 1

*P.291*

Consider a random sample from a distribution with two unknown parameters, the mean  $\mu$  and the variance  $\sigma^2$ . We know from earlier considerations that  $\mu = \mu'_1$  and  $\sigma^2 = E(X^2) - \mu^2 = \mu'_2 - (\mu'_1)^2$ .

Thus,

$$\hat{\sigma}^2 = \mu'_2 - (\mu'_1)^2 = \frac{\sum_{i=1}^n X_i^2}{n} - \frac{\sum_{i=1}^n X_i}{n} = \frac{\sum_{i=1}^n X_i^2}{n} - \bar{X} = \sum_{i=1}^n \frac{(X_i - \bar{X})^2}{n}$$

(Thus, we can see that the estimation of  $\sigma^2$  is not the same as the definition of sample variance  $S^2$ .  $\hat{\sigma}^2 = \frac{n-1}{n} S^2$ .)

### Example 2

*P.292*

If a sample is from a Gamma distribution  $X_i \sim GAM(\theta, k)$ , and we want to estimate the  $\theta$  and  $k$ .

We know that for Gamma distribution, the mean is  $k\theta$ , and the variance is  $k\theta^2$ .

We also know that  $\mu'_1 = \mu = k\theta$  and  $\mu'_2 = \sigma^2 + \mu^2 = k\theta^2 + k^2\theta^2 = k\theta^2(1 + k)$ .

Thus, we can get

$$\bar{X} = k\theta$$

$$\sum \frac{X_i^2}{n} = k\theta^2(1 + k)$$

Thus, we can get

$$\hat{\theta} = \sum_{i=1}^n \frac{(X_i - \bar{X})^2}{n\bar{X}} = \frac{(n-1)/nS^2}{\bar{X}}$$

$$\hat{k} = \frac{\bar{X}}{\hat{\theta}}$$

#### 2.2.1.3 Property

The joint MGF of  $(X_1, \dots, X_n)$  is defined as  $M(t_1, \dots, t_n) = E(e^{\sum_{i=1}^n t_i X_i})$

When  $X_1, \dots, X_n$  are independent if and only if

$$M(t_1, \dots, t_n) = \prod_{i=1}^n M_{X_i}(t_i)$$

where  $M_{X_i}(t_i)$  is the MGF of  $X_i$

**2.2.1.4 Well-known MGF**

- (1) Bernoulli with success probability  $p$ :  $1 - p + pe^t$
- (2) Binomial  $\text{Bin}(n, p)$ :  $(1 - p + pe^t)^n$
- (3) Poisson  $\text{POI}(\lambda)$ :  $e^{\lambda(e^t - 1)}$
- (4) Normal  $N(\mu, \sigma^2)$ :  $e^{\mu t + \frac{1}{2}\sigma^2 t^2}$
- (5) Gamma  $\text{GAM}(\theta, k)$ :  $(1 - \theta t)^{-k}$

Two special cases:

- (6) Chi-square  $\chi^2(v) = \text{GAM}(2, \frac{v}{2})$ :  $(1 - 2t)^{-\frac{v}{2}}$
- (7) Exponential  $\text{EXP}(\theta) = \text{GAM}(\theta, 1)$ :  $(1 - \theta t)^{-1}$

**2.2.2 least squares estimators****2.2.3 likelihood function and maximum likelihood estimators****2.2.3.1 Likelihood function**

*P.293*

The joint density function of  $n$  random variables  $X_1, \dots, X_n$  evaluated at  $x_1, \dots, x_n$ , say  $f(x_1, \dots, x_n; \theta)$ , is referred to as a *likelihood function*.

**2.2.3.2 Maximum likelihood estimators**

*P.294*

Let  $L(\theta) = f(x_1, \dots, x_n; \theta)$ ,  $\theta \in \Omega$ , be the joint pdf of  $X_1, \dots, X_n$ . For a given set of observations,  $(x_1, \dots, x_n)$ , a value  $\hat{\theta}$  in  $\Omega$  at which  $L(\theta)$  is a maximum is called a *maximum likelihood estimate (MLE)* of  $\theta$ . That is  $\hat{\theta}$  is a value of  $\theta$  that satisfies

$$f(x_1, \dots, x_n; \theta) = \text{MAX}_{\theta \in \Omega} f(x_1, \dots, x_n; \theta)$$

### 2.2.4 Invariance property of MLEs

*P.296*

If  $\hat{\theta}$  is the MLE of  $\theta$  and if  $u(\theta)$  is a function of  $\theta$ , then  $u(\hat{\theta})$  is an MLE of  $u(\theta)$ .

#### Example 1

We know that the *pdf* of exponential distribution ( $X \sim EXP(\theta)$ ) is as follows:

$$\frac{1}{\theta} e^{-\frac{x}{\theta}}$$

Thus, its likelihood function is as follows

$$L(\theta) = \frac{1}{\theta^n} e^{-\frac{\sum X_i}{\theta}}$$

Thus, log-likelihood is as follows.

$$\ln L(\theta) = -n \ln(\theta) - \frac{\sum X_i}{\theta}$$

Thus,

$$\frac{d}{d\theta} \ln L(\theta) = -n \frac{1}{\theta} + \frac{\sum X_i}{\theta^2}$$

Thus, we can get the *MLE* for  $\theta$  is  $\hat{\theta} = \bar{x}$ .

If we want to estimate  $\tau(\theta) = P(X \geq 1)$ :

$$\tau(\theta) = P(X \geq 1) = \int_1^{\infty} \frac{1}{\theta} e^{-\frac{x}{\theta}} dx = - \int_1^{\infty} e^{-\frac{x}{\theta}} d\left(-\frac{x}{\theta}\right) = -[e^{-\frac{x}{\theta}}]_1^{\infty} = -[0 - e^{-\frac{1}{\theta}}] = e^{-\frac{1}{\theta}}$$

Thus, based on the invariance property, we know that the *MLE* for  $\tau(\theta)$  is as follows.

$$e^{-\frac{1}{\bar{x}}}$$

#### Example 2: MLE vs. MME

*P.296*

Assume a random sample from a two-parameter exponential distribution,  $X_i \sim EXP(1)$ . Thus, the *pdf* is  $e^{-(x-\eta)}$ . Thus, the likelihood function is

$$L(\eta) = e^{-\sum (x_i - \eta)}$$

Thus, the log likelihood,

$$\ln L(\eta) = -\sum (x_i - \eta) = n\eta - n\bar{X}$$

Thus, we know that as  $\eta$  increases, the log likelihood increases accordingly. Thus, we want to find the maximum  $\eta$ . Note that a two-parameter exponential distribution has the support of  $x_i \geq \eta$ . Thus, all  $\eta$  are smaller than any  $X_i$ . Thus, we can get the ML estimator is the first order statistic

$$\hat{\eta} = X_{1:n}$$

**Note that** the estimators above is based on ML. What would be the answer if using MME?

We know that for a two-parameter exponential distribution, its mean is  $\mu = 1 + \eta$ . And, we know that based on MME,  $\mu = \bar{X}$ . Thus, we can get the following,

$$\hat{\eta} = \bar{X} - 1$$

### Conclusion

We can see that ML and MME have different estimators for the same  $\eta$  for a two-parameter exponential distribution.

### 2.2.5 Unbiased estimators

An estimator  $T$  is said to be an unbiased estimator of  $\tau(\theta)$  if

$$E(T) = \tau(\theta)$$

for all  $\theta \in \Omega$ . Otherwise, we said that  $T$  is biased estimator of  $\tau(\theta)$ .

For instance, if we want to estimate a percentile, say the 95th percentile of  $N(\mu, 9)$ . Note that the percentiles that we know are about standardized normal (i.e.,  $N(0, 1)$ ). Thus, we need to have some calculation to get the non-standard one.

$$\frac{X_{95 \text{ percentile}} - \mu}{\sigma} = 1.645$$

Thus, we can get

$$X_{95 \text{ percentile}} = 1.645 \times \sigma + \mu$$

We know that  $\bar{X}$  is the unbiased estimate for  $\mu$ . Thus, we can get

$$X_{95 \text{ percentile}} = 1.645 \times \sigma + \mu = 4.94 + \mu$$

We know that

$$E(T) = E(\bar{X} + 4.94) = \mu + 4.94$$

Thus,  $T = \bar{X} + 4.94$  is the unbiased estimator of  $\tau(\mu) = \mu + 4.94$ .

### 2.2.6 Unbiased estimators vs. Invariance property of MLEs

**Do not apply “Invariance property of MLEs” to the Unbiased estimators.**

( **Note that:** You can apply the Invariance property to “Unbiased estimators” when it is a linear combination. In that case,  $E(a\theta + b) = aE(\theta) + b$ . Thus, if you find a  $T$  that is a unbiased estimator for  $\theta$ , it should be unbiased estimator for  $a\theta + b$  as well. Thus, note that  $\frac{1}{\theta}$  is not a linear combination of  $\theta$ , thus  $\frac{1}{\theta}$  has a very different estimator, compared to  $\theta$ .)

*P.303*

For example, consider a random sample of size  $n$  from an exponential distribution,  $X_i \sim EXP(\theta)$ , with parameter  $\theta$ . We know that,  $\bar{X}$  is unbiased for  $\theta$  (which is the mean of an exponential distribution).

If we want to estimate  $\tau(\theta) = \frac{1}{\theta}$ , then by the invariance property of MLE is  $T_1 = \frac{1}{\bar{X}}$ .

However,  $T_1$  is a biased estimators of  $\frac{1}{\theta}$ . Specifically,

We know that if  $X \sim Gam(\theta, k)$ , then  $Y = \frac{2X}{\theta} \sim x^2(2k)$ . We know that exponential distributions are a special case of Gamma distribution,  $EXP(\theta) = Gam(\theta, 1)$ , thus we get the following,

$$Y = \frac{2n\bar{X}}{\theta} = \frac{2n}{\theta} \frac{\sum_{i=1}^n X_i}{n} = \frac{\sum_{i=1}^n 2X_i}{\theta} \sim \sum_{i=1}^n x^2(2 \cdot 1) = \sum_{i=1}^n x^2(2) = x^2(2n)$$

We further know that if  $Y \sim x^2(v)$ ,  $E(Y^r) = 2^r \frac{\Gamma(v/2+r)}{\Gamma(v/2)}$ . Thus, we know that,

$$E(Y^{-1}) = 2^{-1} \frac{\Gamma(2n/2 - 1)}{\Gamma(2n/2)} = \frac{1}{2} \cdot \frac{1}{n-1}$$

Thus,

$$E(Y^{-1}) = E\left(\frac{\theta}{2n\bar{X}}\right) = \frac{\theta}{2n} E\left(\frac{1}{\bar{X}}\right) = \frac{1}{2} \cdot \frac{1}{n-1}$$

Thus,

$$E\left(\frac{1}{\bar{X}}\right) = \frac{1}{n-1} \frac{n}{\theta} = \frac{n}{n-1} \frac{1}{\theta}$$

Thus,

$$E\left(\frac{n-1}{n} \frac{1}{\bar{X}}\right) = \frac{1}{\theta}$$

**Conclusion:**

$\frac{1}{\bar{X}}$  is not the unbiased estimator for  $\frac{1}{\theta}$ . However, we can adjust it to  $\frac{n-1}{n} \frac{1}{\bar{X}}$ , which is the unbiased estimator for  $\frac{1}{\theta}$ . When the sample size is big enough, we know that  $\frac{n-1}{n}$  will be close to 1.

## 2.2.7 UMVUE and Cramer-Rao lower bound

### 2.2.7.1 UMVUE

Let  $X_1, X_2, \dots, X_n$  be a random sample of size  $n$  from  $f(x; \theta)$ . An estimator  $T^*$  of  $\tau(\theta)$  is called a *uniformly minimum variance unbiased estimator* (UMVUE) of  $\tau(\theta)$  if

1.  $T^*$  is unbiased for  $\tau\theta$
2. For any other unbiased estimator  $T$  of  $\tau(\theta)$ ,  $Var(T^*) \leq Var(T)$  for all  $\theta \in \Omega$ .

### 2.2.7.2 Cramer-Rao lower bound

If  $T$  is an unbiased estimator of  $\tau(\theta)$ , then the Cramer-Rao lower bound (CRLB), based on a random sample, is

$$Var(T) = \frac{[\tau'(\theta)]^2}{nE\left[\frac{\partial}{\partial\theta} \ln f(X; \theta)\right]^2}$$

**Example:**

Consider a random sample from an exponential distribution,  $X_i \sim \text{Exp}(\theta)$ . Because

$$f(x; \theta) = \frac{1}{\theta} e^{-\frac{x}{\theta}}$$

Thus,

$$\ln(f(x; \theta)) = -\frac{x}{\theta} - \ln \theta$$

$$\frac{\partial}{\partial \theta} \ln(f(X; \theta)) = \frac{X}{\theta^2} - \frac{1}{\theta} = \frac{X - \theta}{\theta^2}$$

Thus,

$$E\left[\frac{\partial}{\partial \theta} \ln(f(X; \theta))\right]^2 = E\left[\frac{(X - \theta)^2}{\theta^4}\right] = \frac{1}{\theta^4} E(X - \theta)^2 = \frac{1}{\theta^4} \text{Var}(X) = \frac{\theta^2}{\theta^4} = \frac{1}{\theta^2}$$

Thus, the CRLB for  $\tau(\theta) = \theta$  is as follows.

$$\text{Var}(T) \geq \frac{[\tau'(\theta)]^2}{nE\left[\frac{\partial}{\partial \theta} \ln f(X; \theta)\right]^2} = \frac{\left[\frac{\partial}{\partial \theta} \theta\right]^2}{n \frac{1}{\theta^2}} = \frac{1^2}{\frac{n}{\theta^2}} = \frac{\theta^2}{n}$$

We know that the variance for exponential distribution

$$\text{Var}(\bar{X}) = \frac{\text{Var}(X)}{n} = \frac{\theta^2}{n}$$

We know that  $\theta$  is the mean for exponential distribution. We also know that sample mean  $\bar{X}$  is the unbiased estimator for population mean.

Thus,

$\bar{X}$  is the UMVUE of  $\theta$ .

### 2.2.8 Best linear unbiased estimation (BLUE or MVLUE)

### 2.2.9 Consistency, asymptotic unbiasedness

**Simple consistency:**

*P.311*

Let  $\{T_n\}$  be a sequence of estimators of  $\tau(\theta)$ . These estimators are said to be **consistent** estimators of  $\tau(\theta)$  if for every  $\varepsilon > 0$ ,



$$\lim_{n \rightarrow \infty} P[|T_n - \tau(\theta)| < \varepsilon] = 1$$

for every  $\theta \in \Omega$ .

In the terminology of Chapter 7,  $T_n$  converges stochastically to  $\tau(\theta)$ ,  $T_n \xrightarrow{P} \tau(\theta)$  as  $n \rightarrow \infty$ . Sometimes this also is referred to as **simple consistency**.

One interpretation of consistency is that for large sample size the estimator tends to be more concentrated about  $\tau(\theta)$ , and by making  $n$  sufficiently large  $T_n$  can be made as concentrated as desired.

**MSE consistency:**

If  $\{T_n\}$  is a sequence of estimator of  $\tau(\theta)$ , then they are called **mean squared error consistent** if

$$\lim_{n \rightarrow \infty} E[T_n - \tau(\theta)]^2 = 0$$

for every  $\theta \in \Omega$ .

**Asymptotic Unbiased**

A sequence  $\{T_n\}$  is said to be **asymptotically unbiased** for  $\tau(\theta)$  if

$$\lim_{n \rightarrow \infty} E(T_n) = \tau(\theta)$$

for every  $\theta \in \Omega$ .

**Example:**

*P.313*

For a sample from  $X_i \sim EXP(\theta)$ , we know that  $T_n = 1/\bar{X}$  is an MLE estimator for  $\tau(\theta) = 1/\theta$ . However,  $T_n$  is not a unbiased estimator for  $\tau(\theta)$ , as

$$E(T_n) = \frac{n}{n-1} \cdot \frac{1}{\theta}$$

We also know that

$$Var(T_n) = \frac{(\frac{n}{n-1})^2}{(n-2)\theta^2}$$

Thus, while  $T_n$  is not unbiased, it is asymptotically unbiased and MSE consistent for  $\tau(\theta) = \frac{1}{\theta}$ .

**Note that,** MES consistency is a stronger property than simple consistency. Thus, if a sequenc  $\{T_n\}$  is mean squared error consistent, it is also simply consistent.

### 2.2.10 Efficiency

*P.308*

The relative efficiency of an unbiased estimator  $T$  of  $\tau(\theta)$  to another unbiased estimator  $T^*$  of  $\tau(\theta)$  is given by

$$re(T, T^*) = \frac{Var(T^*)}{Var(T)}$$

An unbiased estimator  $T^*$  of  $\tau(\theta)$  is said to be efficient if  $re(T, T^*) \leq 1$  for all unbiased estimators  $T$  of  $\tau(\theta)$ , and all  $\theta \in \Omega$ . The efficiency of an unbiased estimator  $T$  of  $\tau(\theta)$  is given by

$$e(T) = re(T, T^*)$$

if  $T^*$  is an efficient estimator of  $\tau(\theta)$ .

Thus, **an effecient estimator is just a UMVUE.**

**Example:**

*P.233 Example 7.2.2*

*P.303 Example 9.3.2*

*P.309 Example 9.3.7*

Let  $X_1, X_2, \dots, X_n$  sample from  $X_i \sim EXP(\theta)$ . We know that

$$X_{1:n} = EXP(\theta/n)$$

Thus,

$$E(nX_{1:n}) = nE(X_{1:n}) = n\frac{\theta}{n} = \theta$$

Thus,

$$nX_{1:n}$$

is the unbiased estimator for  $\theta$ .

Thus,

$$re(T, T^*) = \frac{Var(T^*)}{Var(T)} = \frac{Var(\bar{X})}{Var(nX_{1:n})} = \frac{\theta^2/n}{n^2 Var(X_{1:n})} = \frac{\theta^2/n}{n^2(\theta/n)^2} = \frac{\theta^2/n}{\theta^2} = \frac{1}{n}$$

Thus,  $T^* = \bar{X}$  is a more efficient estimator for  $\theta$  than  $T = nX_{1:n}$ .

### 2.2.11 Asymptotic efficiency

Let  $\{T_n\}$  and  $\{T_n^*\}$  be the two asymptotically unbiased sequences of estimators for  $\tau(\theta)$ . The **asymptotic relative efficiency** of  $T_n$  relative to  $T_n^*$  is given by

$$are(T_n, T_n^*) = \lim_{n \rightarrow \infty} \frac{Var(T_n^*)}{T_n}$$

The sequence  $\{T_n^*\}$  is said to be asymptotically efficient if  $are\{T_n, T_n^*\} \leq 1$  for all other asymptotically unbiased sequences  $\{T_n\}$ , and all  $\theta \in \Omega$ .

### 2.2.12 Asymptotic properties of MLEs

*P.316*

If certain regularity conditions are satisfied, then the solutions,  $\hat{\theta}$ , of the MLE have the following properties:

- (1)  $\hat{\theta}_n$  exists and is unique.
- (2)  $\hat{\theta}_n$  is a consistent estimator of  $\theta$ .
- (3)  $\hat{\theta}_n$  is asymptotically normal with asymptotic mean  $\theta$  and variance

$$\frac{1}{2} E \left[ \frac{\partial}{\partial \theta} \ln f(X; \theta) \right]^2$$

- (4)  $\hat{\theta}$  is asymptotically efficient.

Note that the asymptotic efficiency of  $\hat{\theta}$  follows from the fact that the asymptotic variance is the same as the **CRLB** for unbiased estimators of  $\theta$ . Thus, for large  $n$ , approximately

$$\hat{\theta}_n \sim N(\theta, CRLB)$$

#### Example

*P.317*

From Example 9.2.7, we know the MLE of the mean  $\theta$  of an exponential distribution is the sample mean,  $\hat{\theta}_n = \bar{X}$ .

We can infer the same asymptotic properties either from the discussion above or from the Central Limit Theorem. That is,  $\hat{\theta}_n$  is asymptotically normal with

asymptotic mean  $\theta$  and variance  $\theta^2/n$ . From example 9.3.4, we know that  $CRLB = \theta^2/n$ .

Thus,

$$\sqrt{n} \frac{\hat{\theta}_n - \theta}{\theta} \sim N(0, 1)$$

## 2.3 Sufficient and completeness

### 2.3.1 Sufficiency and minimal sufficiency

*P.335*

#### Sufficient statistic:

A statistic  $S$  will be considered a “sufficient” statistic for a parameter  $\theta$  if the conditional distribution of any other statistic  $T$  given the value of  $S$  does not involve  $\theta$ .

#### Jointly sufficient statistics:

Let  $X = (X_1, \dots, X_n)$  have joint pdf  $f(x, \theta)$ , and let  $S = (S_1, \dots, S_k)$  be a  $k$ -dimensional statistic.

Then,  $S_1, \dots, S_k$  is a set of **jointly sufficient statistics** for  $\theta$  if for any other vector of statistics,  $T$ , the conditional pdf of  $T$  given  $S = s$ , denoted by  $f_{T|s}(t)$ , does not depend on  $\theta$ .

In the one-dimension case, we simply say that  $S$  is a sufficient statistic for  $\theta$ .

If  $k$  unknown parameters are present in the model, then quite often there will exist a set of  $k$  sufficient statistics. In some cases, the number of sufficient statistics will exceed the number of parameters.

#### Minimal sufficient:

A set of statistics is called a **minimal sufficient** set if the members of the set are jointly sufficient for the parameters and if they are a function of every other set of jointly sufficient statistics.

**In other words, once the value of a sufficient statistic is known, the observed value of any other statistic does not contain any further information about the parameter.**

#### Example:

*P.339*

Assume a random sample from an exponential distribution,  $X_i \sim EXP(\theta)$ . It follows that

$$f(x_1, \dots, x_n; \theta) = \frac{1}{\theta^n} e^{-\frac{\sum x_i}{\theta}}$$

which suggests checking the statistic  $S = \sum X_i$ . We know that  $S \sim GAM(\theta, n)$ , thus

$$f_S(s; \theta) = \frac{1}{\theta^n \Gamma(n)} s^{n-1} e^{-\frac{s}{\theta}}$$

If  $s = \sum x_i$ , then,

$$\frac{f(x_1, \dots, x_n; \theta)}{f_S(s; \theta)} = \frac{\Gamma(n)}{s^{n-1}}$$

which is free of  $\theta$ , and thus by definition  $S$  is sufficient for  $\theta$ .

### 2.3.2 Neyman factorization theorem, minimal sufficiency of MLEs

#### Factorization criterion:

If  $X_1, \dots, X_n$  have joint pdf  $f(x_1, \dots, x_n; \theta)$ , and if  $S = (S_1, \dots, S_n)$ , then  $S_1, \dots, S_k$  are jointly sufficient for  $\theta$  if and only if

$$f(x_1, \dots, x_n; \theta) = g(s; \theta) h(x_1, \dots, x_n)$$

where  $g(s; \theta)$  does not depend on  $x_1, \dots, x_n$ , except through  $s$ , and  $h(x_1, \dots, x_n)$  does not involve  $\theta$ .

#### Example 1:

P.340

$X_i \sim \text{Bin}(1, \theta)$ . We can use factorization criterion to check its sufficient statistic.

$$f(x_1, \dots, x_n; \theta) = \theta^{\sum x_i} (1 - \theta)^{n - \sum x_i} = \theta^s (1 - \theta)^{n-s} = g(s; \theta) h(x_1, \dots, x_n)$$

where  $s = \sum X_i$  and  $h(x_1, \dots, x_n) = 1$ . Thus,  $s$  is sufficient for  $\theta$ .

#### Example 2:

**It is important to specify the regions of zero probability. The following shows that care must be exercised in this matter.**

P.340

$X_i \sim \text{UNIF}(0, \theta)$ , where  $\theta$  is unknown. We get the joint pdf of  $X_1, \dots, X_n$  is

$$f(x_1, \dots, x_n; \theta) = \frac{1}{\theta^n}$$

where,

$$0 < x_i < \theta$$

It is easier to specify this *pdf* in terms of the minimum,  $x_{1:n}$ , and maximum,  $x_{n:n}$ , of  $x_1, \dots, x_n$ . In particular,

$$0 < x_{1:n} \leq x_{n:n} < \theta$$

Thus,  $x_{n:n}$  is sufficient for  $\theta$ .

**Example 3:**

*P.341*

Consider a random sample from a normal distribution,  $X_i \sim N(\mu, \sigma^2)$ .

We know the *pdf* for normal is

$$f = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{1}{2\sigma^2}(x-\mu)^2}$$

Thus, the joint *pdf* is as follows.

$$f(x_1, \dots, x_n; \mu, \sigma^2) = \frac{1}{(2\pi\sigma^2)^{\frac{n}{2}}} e^{-\frac{1}{2\sigma^2} \sum (x_i - \mu)^2} = \frac{1}{(2\pi\sigma^2)^{\frac{n}{2}}} e^{-\frac{1}{2\sigma^2} \sum (x_i^2 + \mu^2 - 2x_i\mu)} = \frac{1}{(2\pi\sigma^2)^{\frac{n}{2}}} e^{-\frac{1}{2\sigma^2} (\sum x_i^2 + n\mu^2 - 2\mu \sum x_i)}$$

Thus,  $S_1 = \sum X_i$  and  $S_2 = \sum X_i^2$  are jointly sufficient for  $\theta = \mu, \sigma^2$ .

**Note that,**

*P.341*

based on *P.298*, Example 9.2.10, ML estimation results in

$$\hat{\mu} = \bar{x}$$

$$\hat{\sigma}^2 = \frac{\sum (x_i - \bar{x})^2}{n}$$

Further, note that

$$\hat{\mu} = \bar{x} = \frac{S_1}{n}$$

$$\begin{aligned}
\hat{\sigma}^2 &= \frac{\sum (x_i - \bar{x})^2}{n} = \frac{\sum (x_i^2 + \bar{x}^2 - 2x_i\bar{x})}{n} \\
&= \frac{\sum x_i^2 + n\bar{x}^2 - 2\bar{x} \sum x_i}{n} \\
&= \frac{\sum x_i^2 + n\bar{x}^2 - 2n\bar{x}^2}{n} \\
&= \frac{\sum x_i^2 - n\bar{x}^2}{n} \\
&= \frac{\sum x_i^2}{n} - \bar{x}^2 \\
&= \frac{\sum x_i^2}{n} - \left(\frac{\sum x_i}{n}\right)^2 \\
&= \frac{S_2}{n} - \left(\frac{S_1}{n}\right)^2
\end{aligned}$$

Thus, we can see that  $\hat{\mu}$  and  $\hat{\sigma}^2$  correspond to a one-to-one transformation of  $S_1$  and  $S_2$ , thus  $\hat{\mu}$  and  $\hat{\sigma}^2$  also are jointly sufficient for  $\mu$  and  $\sigma^2$ .

### 2.3.3 Rao-Blackwell theorem

Let  $X_1, \dots, X_n$  have joint *pdf*  $f(x_1, \dots, x_n; \theta)$ , and let  $S = (S_1, \dots, S_k)$  be a vector of jointly sufficient statistics for  $\theta$ . If  $T$  is any unbiased estimator of  $\tau(\theta)$ , and if  $T^* = E(T|S)$ , then

- (1)  $T^*$  is an unbiased estimator of  $\tau(\theta)$ ,
- (2)  $T^*$  is a function of  $S$ , and
- (3)  $Var(T^*) \leq Var(T)$  for every  $\theta$ , and  $Var(T^*) < Var(T)$  for some  $\theta$  unless  $T^* = T$  with probability 1.

#### Discussion:

*P.345*

"It is clear from the Rao-Blackwell theorem that if we are searching for an unbiased estimator with small variance, we may as well restrict attention to function of sufficient statistics. If any unbiased estimator exists, then there will be one that is a function of sufficient statistics, namely  $E(T|S)$ , which also is unbiased and has variance at least as small or smaller. In particular, we still are interested in knowing how to find a UMVUE for a parameter, and the above theorem narrows our problem down somewhat.

For instance, consider a one-parameter model  $f(x; \theta)$ , and assume that a single sufficient statistic,  $S$ , exists. We know that we must consider only unbiased

functions of  $S$  in searching for a UMVUE. in some cases it may be possible to show that only one function of  $S$  is unbiased, and in that case we would know that it is a UMVUE.

The concept of “completeness” is helpful in determining unique unbiased estimators, and this concept is defined in the next section.”

### 2.3.4 completeness

#### Completeness:

*P.345*

A family of density functions  $\{f_T(t, \theta); \theta \in \Omega\}$ , is called **complete** if  $E[u(T)] = 0$  for all  $\theta \in \Omega$  implies  $u(T) = 0$  with probability 1 for all  $\theta \in \Omega$ .

*P.345*

“This sometimes is expressed by saying that there are no nontrivial unbiased estimators of zero. In particular, it means that two different functions of  $T$  cannot have the same expected value.”

For example, if  $E[u_1(T)] = \tau(\theta)$  and  $E[u_2(T)] = \tau(\theta)$ , which implies  $u_1(T) - u_2(T) = 0$ , or  $u_1(T) = u_2(T)$  with probability 1, if the family of density functions is complete. That is, any unbiased estimator is unique in this case.

We mainly are interested in knowing that the family of density functions of a sufficient statistic is complete, because in that case an unbiased function of the sufficient statistic will be unique, and it must be a UMVUE by the Rao-Blackwell theorem.

### 2.3.5 Lehmann-Scheffe completeness theorem

Let  $X_1, \dots, X_n$  have joint pdf  $f(x_1, \dots, x_n; \theta)$  and let  $S$  be a vector of jointly complete sufficient statistics for  $\theta$ . If  $T^* = t^*(S)$  is a statistic that is unbiased for  $\tau(\theta)$  and a function of  $S$ , then  $T^*$  is a UMVUE of  $\tau(\theta)$ .

#### Example:

*P.346*

### 2.3.6 Exponential class, complete sufficient statistics

A density function is said to be a member of the **regular exponential class** if it can be expressed in the form

$$f(x; \theta) = c(\theta)h(x)e^{\sum_{j=1}^k q_j(\theta)t_j(x)}$$



where,

$$x \in A$$

And zero otherwise, where  $\theta = (\theta_1, \dots, \theta_k)$  is a vector of  $k$  unknown parameters, if the parameter space has the form

$$\Omega = \{\theta | a_i \leq \theta_i \leq b_i, i = 1, \dots, k\}$$

(note that  $a_i = -\infty$  and  $b_i = \infty$  are permissible values), and if it satisfies regularity conditions 1, 2, and 3a or 3b given by:

- (1) The set  $A = \{x : f(x; \theta) > 0\}$  does not depend on  $\theta$ .
- (2) The functions  $q_j(\theta)$  are nontrivial, functionally independent, continuous function of the  $\theta_i$ .

(3a) For a continuous random variable, the derivatives  $t'_k(x)$  are linearly independent continuous functions of  $x$  over  $A$ .

(3b) For a discrete random variable, the  $t_k(x)$  are nontrivial functions of  $x$  on  $A$ , and none is a linear function of the others.

### Example

*P.348*

Consider Bernoulli distribution  $X \sim \text{Bin}(1, p)$ . It follows that

$$\begin{aligned} f(x; p) &= p^x (1-p)^{1-x} \\ &= P^x (1-p) (1-p)^{-x} \\ &= (1-p) (1-p)^{-x} P^x \\ &= (1-p) e^{\ln(1-p)^{-x} P^x} \\ &= (1-p) e^{x \ln(\frac{p}{1-p})} \end{aligned}$$

Compared to the definition of **Exponential Class** defined above, we can get the following.

$$\begin{aligned} c(\theta) &= (1-p) \\ h(x) &= 1 \\ q_1(\theta) &= \ln\left(\frac{p}{1-p}\right) \\ t_1(x) &= x \end{aligned}$$

### Theorem 10.4.2

P.348

If  $X_1, \dots, X_n$  is a random sample from a member of the regular exponential class  $REC(q_1, \dots, q_k)$ , then the statistics

$$S_1 = \sum_{i=1}^n t_1(X_i), \dots, S_k = \sum_{i=1}^n t_k(X_i)$$

are a minimal set of complete sufficient statistics for  $\theta_1, \dots, \theta_k$ .

**Example 1:**

P.348

Again, consider Bernoulli distribution  $X \sim Bin(1, p)$ . For a random sample of size  $n$ ,  $t(x_i) = x_i$  and thus based on **Theorem 10.4.2** (see above)  $S = \sum_{i=1}^n X_i$  is a complete sufficient statistic for  $p$ .

**How about we want to find UMVUE for  $Var(X) = p(1-p)$ ?**

We might try  $\bar{X}(1 - \bar{X})$ .

$$\begin{aligned} E[\bar{X}(1 - \bar{X})] &= E(\bar{X}) - E(\bar{X}^2) \\ &= p - (p^2 + var(\bar{X})) \\ &= p - p^2 - \frac{p(1-p)}{n} \\ &= p(1-p)(1 - \frac{1}{n}) \end{aligned}$$

Thus,

$$E[\frac{1}{1 - \frac{1}{n}} \bar{X}(1 - \bar{X})] = p(1-p)$$

Thus,  $\frac{1}{1 - \frac{1}{n}} \bar{X}(1 - \bar{X})$  is the UMVUE of the  $p(1-p)$ .

but **why?**

**Example 2:**

P.349

If  $X \sim N(\mu, \sigma^2)$ , then

$$\begin{aligned} f(x; \mu, \sigma) &= \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}} \\ &= \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{1}{2\sigma^2}(x^2 + \mu^2 - 2x\mu)} \\ &= \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{x^2}{2\sigma^2} - \frac{\mu^2}{2\sigma^2} + \frac{x\mu}{\sigma^2}} \end{aligned}$$

Thus,  $S_1 = \sum X_i^2$  and  $S_2 = \sum X_i$  are jointly complete and sufficient statistics of  $\mu$  and  $\sigma^2$ . **Refer to the normal example in the section of “Neyman factorization theorem, minimal sufficiency of MLEs”**

**Theorem 10.4.3**

*P.349*

If a CRLB estimator  $T$  exists for  $\tau(\theta)$ , then a single sufficient statistic exists, and  $T$  is a function of the sufficient statistic. Conversely, if a single sufficient statistic exists and the CRLB exists, then a CRLB estimator exists for some  $\tau(\theta)$ .



## Chapter 3

**530\_\_533**

**3.1 Definition of the general linear model - full rank, less than full rank**

**3.2 Simple Linear Regression**

**3.2.1 MLE vs. least squares, propoerties of the regression parameters**

**3.2.2 Assumptions, checking assumptions**

**3.3 Bootstrapping used in linear models**

**3.4 Generalized linear models**

**3.4.1 Definition, similarities and differences from general linear models**

**3.4.2 Advantage and disadvantages**

**3.4.3 logistic and poisson regression**



## Chapter 4

# 512

- 4.1 Completely randomized designs
- 4.2 Randomized complete block designs
- 4.3 incomplete block designs
- 4.4 Split-plot designs
- 4.5 Row-column design (Latin square designs)





## Chapter 5

# Logit and Probit

### 5.1 Logit

$$f(x) = \log\left(\frac{p(y=1)}{1-p(y=1)}\right)$$

The basic idea of logistic regression:

$$p(y=1) = \frac{1}{1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)}} = \frac{e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}{1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}$$

Thus,  $\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n$  can be from  $-\infty$  to  $+\infty$ , and  $p(y=1)$  will be always within the range of  $(0, 1)$ .

```
f<-function(x){exp(x)/(1+exp(x))}  
data<-seq(-10,10,1)  
plot(data,f(data),type = "b")
```



We can also write the function into another format as follows:

$$\log \frac{p(y=1)}{1-p(y=1)} = \beta_0 + \beta_1 x_1 + \dots + \beta_n x_n$$

Thus, we know that the regression coefficients of  $\beta_i$  actually change the “log-odds” of the event. Of course, note that the magnitude of  $\beta_i$  is dependent upon the units of  $x_i$ .

The following is an example testing whether that home teams are more likely to win in NFL games. The results show that the odd of winning is the same for both home and away teams.

```
mydata = read.csv(url('https://raw.githubusercontent.com/nfl-football-ops/Big-Data-Bow
mydata$result_new<-ifelse(mydata$HomeScore>mydata$VisitorScore,1,0)
summary(mydata$result_new)
```

```
##      Min. 1st Qu.  Median    Mean 3rd Qu.    Max.
## 0.0000 0.0000 0.0000 0.4945 1.0000 1.0000
```

```
mylogit1 = glm(result_new~1, family=binomial, data=mydata)
summary(mylogit1)
```

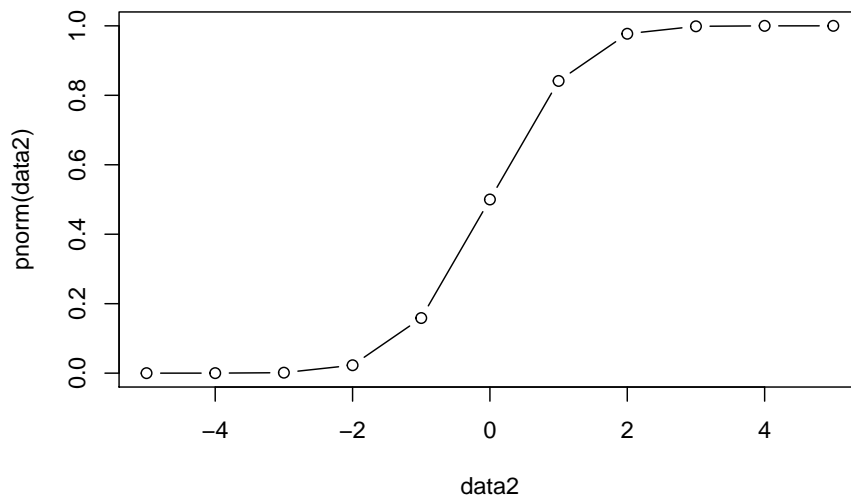
```
##
## Call:
```

```
## glm(formula = result_new ~ 1, family = binomial, data = mydata)
##
## Deviance Residuals:
##      Min       1Q   Median       3Q      Max
## -1.168  -1.168  -1.168   1.187   1.187
##
## Coefficients:
##              Estimate Std. Error z value Pr(>|z|)
## (Intercept) -0.02198    0.20967  -0.105   0.917
##
## (Dispersion parameter for binomial family taken to be 1)
##
##      Null deviance: 126.14  on 90  degrees of freedom
## Residual deviance: 126.14  on 90  degrees of freedom
## AIC: 128.14
##
## Number of Fisher Scoring iterations: 3
```

## 5.2 Probit

As noted above, logit  $f(x) = \log\left(\frac{p(y=1)}{1-p(y=1)}\right)$  provides the resulting range of (0, 1). Another way to provide the same range is through the cdf of normal distribution. The following R code is used to illustrate this process.

```
data2<-seq(-5,5,1)
plot(data2,pnorm(data2),type = "b")
```



Thus, the cdf of normal distribution can be used to indicate the probability of  $p(y = 1)$ .

$$\Phi(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n) = p(y = 1)$$

Similar to logit model, we can also write the inverse function of the cdf to get the function that can be from  $-\infty$  to  $+\infty$ .

$$\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n = \Phi^{-1}(p(y = 1))$$

Thus, for example, if  $X\beta = -2$ , based on  $\Phi(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n) = p(y = 1)$  we can get that the  $p(y = 1) = 0.023$ .

In contrast, if  $X\beta = 3$ , the  $p(y = 1) = 0.999$ .

```
pnorm(-2)
```

```
## [1] 0.02275013
```

```
pnorm(3)
```

```
## [1] 0.9986501
```

Let's assume that there is a latent variable called  $Y^*$  such that

$$Y^* = X\beta + \epsilon, \epsilon \sim N(0, \sigma^2)$$

You could think of  $Y^*$  as a kind of “proxy” between  $X\beta + \epsilon$  and the observed  $Y(1 \text{ or } 0)$ . Thus, we can get the following. Note that, it does not have to be zero, and can be any constant.

$$Y^* = \begin{cases} 0 & \text{if } y_i^* \leq 0 \\ 1 & \text{if } y_i^* > 0 \end{cases}$$

Thus,

$$y_i^* > 0 \Rightarrow \beta' X_i + \epsilon_i > 0 \Rightarrow \epsilon_i > -\beta' X_i$$

Thus, we can write it as follows. Note that  $\frac{\epsilon_i}{\sigma} \sim N(0, 1)$

$$p(y = 1|x_i) = p(y_i^* > 0|x_i) = p(\epsilon_i > -\beta' X_i) = p\left(\frac{\epsilon_i}{\sigma} > \frac{-\beta' X_i}{\sigma}\right) = \Phi\left(\frac{\beta' X_i}{\sigma}\right)$$

We thus can get:

$$p(y = 0|x_i) = 1 - \Phi\left(\frac{\beta' X_i}{\sigma}\right)$$

For  $p(y = 1|x_i) = \Phi\left(\frac{\beta' X_i}{\sigma}\right)$ , we can not really estimate both  $\beta$  and  $\sigma$  as they are in a ratio. We can assume  $\sigma = 1$ , then  $\epsilon \sim N(0, 1)$ . We know  $y_i$  and  $x_i$  since we observe them. Thus, we can write it as follows.

$$p(y = 1|x_i) = \Phi(\beta' X_i)$$



## Chapter 6

# Normal distribution

### 6.1 Basics

$\mu$  and  $\sigma$  determine the center and spread of the distribution.

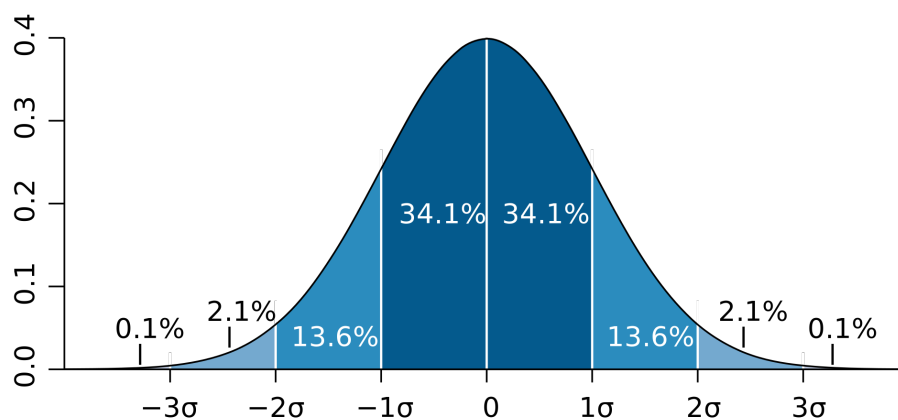


Figure 6.1: Normal

The empirical rule holds for all normal distributions:

- (1) 68% of the area under the curve lies between  $(\mu - \sigma, \mu + \sigma)$ .
- (2) 95% of the area under the curve lies between  $(\mu - 2\sigma, \mu + 2\sigma)$ .
- (3) 99.7% of the area under the curve lies between  $(\mu - 3\sigma, \mu + 3\sigma)$ .

## 6.2 Confidence intervals for normal distributions

$$\bar{X} \pm Z \frac{\sigma}{\sqrt{n}}$$

where,

$\bar{X}$  is the mean

$Z$  is the Z value (see the table below)

$\sigma$  is the standard deviation

$n$  is the number of observations

(We can see the connection between this formula and information shown in the *Basics* section.)

<i>Confidence Levels</i>	<i>Z</i>
80	1.282
85	1.440
90	1.645
95	1.960
99	2.576
99.5	2.807
99.9	3.291

## 6.3 Percentile

A percentile is a measure used in statistics indicating the value below which a given percentage of observations in a group of observations falls.

For example, the 20th percentile is the value (or score) below which 20% of the observations may be found.

For normal distribution,

$-3 \sigma$  is the 0.13th percentile (i.e.,  $\frac{100-99.7}{2} = 0.15$ );

$-2 \sigma$  is the 2.28th percentile ((i.e.,  $\frac{100-95}{2} = 2.50$ ));

$-1\sigma$  is the 15.87th percentile (i.e.,  $\frac{100-68}{2} = 16$ );

$0 \sigma$  is 50th percentile.

$+2 \sigma$  is the 97.72nd percentile (i.e.,  $100 - \frac{100-95}{2} = 100 - 2.5 = 97.50$ );

$+3 \sigma$  is the 99.87th percentile (i.e.,  $100 - \frac{100-99.70}{2} = 100 - 0.15 = 99.85$ ).

This is related to the 68-95-99.7 rule or the three-sigma rule.



(Note that, it is *related*, not *direct* 68-95-99.7 rule, which is about symmetric situations. See the figure above)

<i>Percentile</i>	<i>Z</i>
90	1.282
—	1.440
95	1.645
—	1.960
—	2.576
—	2.807
99.9	3.000



# Chapter 7

## MLE

### 7.1 Basic idea of MLE

Suppose that we flip a coin,  $y_i = 0$  for tails and  $y_i = 1$  for heads. If we get  $p$  heads from  $n$  trials, we can get the proportion of heads is  $p/n$ , which is the sample mean. If we do not do any further calculation, this is our best guess.

Suppose that the true probability is  $\rho$ , then we can get:

$$\mathbf{L}(y_i) = \begin{cases} \rho & y_i = 1 \\ 1 - \rho & y_i = 0 \end{cases}$$

Thus, we can also write it as follows.

$$\mathbf{L}(y_i) = \rho^{y_i} (1 - \rho)^{1-y_i}$$

Thus, we can get:

$$\prod \mathbf{L}(y_i|\rho) = \rho^{\sum y_i} (1 - \rho)^{\sum (1-y_i)}$$

Further, we can get a log-transformed format.

$$\log(\prod \mathbf{L}(y_i|\rho)) = \sum y_i \log \rho + \sum (1 - y_i) \log(1 - \rho)$$

To maximize the log-function above, we can calculate the derivative with respect to  $\rho$ .

$$\frac{\partial \log(\prod \mathbf{L}(y_i|\rho))}{\partial \rho} = \sum y_i \frac{1}{\rho} - \sum (1 - y_i) \frac{1}{1 - \rho}$$

Set the derivative to zero and solve for  $\rho$ , we can get

$$\begin{aligned}
& \sum y_i \frac{1}{\rho} - \sum (1 - y_i) \frac{1}{1 - \rho} = 0 \\
& \Rightarrow (1 - \rho) \sum y_i - \rho \sum (1 - y_i) = 0 \\
& \Rightarrow \sum y_i - \rho \sum y_i - n\rho + \rho \sum y_i = 0 \\
& \Rightarrow \sum y_i - n\rho = 0 \\
& \Rightarrow \rho = \frac{\sum y_i}{n} = \frac{p}{n}
\end{aligned}$$

Thus, we can see that the  $\rho$  maximizing the likelihood function is equal to the sample mean.

## 7.2 Coin flip example, probit, and logit

In the example above, we are not really trying to estimate a lot of regression coefficients. What we are doing actually is to calculate the sample mean, or intercept in the regression sense. What does it mean? Let's use some data to explain it.

Suppose that we flip a coin 20 times and observe 8 heads. We can use the R's `glm` function to estimate the  $\rho$ . If the result is consistent with what we did above, we should observe that the *cdf* of the estimate of  $\beta_0$  (i.e., intercept) should be equal to  $8/20 = 0.4$ .

```
coins<-c(rep(1,times=8),rep(0,times=12))
table(coins)
```

```
## coins
##  0  1
## 12  8
```

```
coins<-as.data.frame(coins)
```

### 7.2.1 Probit

```
probitresults <- glm(coins ~ 1, family = binomial(link = "probit"), data = coins)
probitresults
```

```
##
## Call:  glm(formula = coins ~ 1, family = binomial(link = "probit"),
##       data = coins)
##
## Coefficients:
## (Intercept)
##      -0.2533
##
## Degrees of Freedom: 19 Total (i.e. Null);  19 Residual
## Null Deviance:      26.92
## Residual Deviance: 26.92      AIC: 28.92
```

```
pnorm(probitresults$coefficients)
```

```
## (Intercept)
##           0.4
```

As we can see the intercept is  $-0.2533$ , and thus  $\Phi(-0.2533471) = 0.4$

## 7.2.2 Logit

We can also use logit link to calculate the intercept as well. Recall that

$$p(y = 1) = \frac{1}{1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)}} = \frac{e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}{1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}$$

Thus,

$$p(y = 1) = \frac{e^{\beta_0}}{1 + e^{\beta_0}}$$

```
logitresults <- glm(coins ~ 1, family = binomial(link = "logit"), data = coins)
logitresults$coefficients
```

```
## (Intercept)
##      -0.4054651
```

```
exp(logitresults$coefficients)/(1+exp(logitresults$coefficients))
```

```
## (Intercept)
##           0.4
```

Note that, the default link for the binomial in the glm function is logit.

### 7.3 Further on logit

The probability of  $y = 1$  is as follows:

$$p = p(y = 1) = \frac{1}{1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)}} = \frac{e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}{1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}}$$

Thus, the likelihood function is as follows:

$$\begin{aligned} L &= \prod p^{y_i} (1-p)^{1-y_i} = \prod \left( \frac{1}{1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)}} \right)^{y_i} \left( \frac{1}{1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}} \right)^{1-y_i} \\ &= \prod (1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)})^{-y_i} (1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n})^{-(1-y_i)} \end{aligned}$$

Thus, the log-likelihood is as follows:

$$\log L = \sum (-y_i \cdot \log(1 + e^{-(\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n)}) - (1 - y_i) \cdot \log(1 + e^{\beta_0 + \beta_1 x_1 + \dots + \beta_n x_n}))$$

Typically, optimisers minimize a function, so we use negative log-likelihood as minimising that is equivalent to maximising the log-likelihood or the likelihood itself.

*#Source of R code: <https://www.r-bloggers.com/logistic-regression/>*

```
mle.logreg = function(fmla, data)
{
  # Define the negative log likelihood function
  logl <- function(theta,x,y){
    y <- y
    x <- as.matrix(x)
    beta <- theta[1:ncol(x)]

    # Use the log-likelihood of the Bernoulli distribution, where p is
    # defined as the logistic transformation of a linear combination
    # of predictors, according to logit(p)=(x%*%beta)
    loglik <- sum(-y*log(1 + exp(-(x%*%beta))) - (1-y)*log(1 + exp(x%*%beta)))
    return(-loglik)
  }

  # Prepare the data
  outcome = rownames(attr(terms(fmla),"factors"))[1]
  dfrTmp = model.frame(data)
  x = as.matrix(model.matrix(fmla, data=dfrTmp))
}
```

```

y = as.numeric(as.matrix(data[,match(outcome,colnames(data))]))

# Define initial values for the parameters
theta.start = rep(0,(dim(x)[2]))
names(theta.start) = colnames(x)

# Calculate the maximum likelihood
mle = optim(theta.start,logl,x=x,y=y, method = 'BFGS', hessian=T)
out = list(beta=mle$par,vcov=solve(mle$hessian),ll=2*mle$value)
}

mydata = read.csv(url('https://stats.idre.ucla.edu/stat/data/binary.csv'))
mylogit1 = glm(admit~gre+gpa+as.factor(rank), family=binomial, data=mydata)

mydata$rank = factor(mydata$rank) #Treat rank as a categorical variable
fmla = as.formula("admit~gre+gpa+rank") #Create model formula
mylogit2 = mle.logreg(fmla, mydata) #Estimate coefficients

print(cbind(coef(mylogit1), mylogit2$beta))

##                [,1]      [,2]
## (Intercept)    -3.989979073 -3.772676422
## gre            0.002264426  0.001375522
## gpa            0.804037549  0.898201239
## as.factor(rank)2 -0.675442928 -0.675543009
## as.factor(rank)3 -1.340203916 -1.356554831
## as.factor(rank)4 -1.551463677 -1.563396035

```

## 7.4 References

[http://www.columbia.edu/~so33/SusDev/Lecture\\_9.pdf](http://www.columbia.edu/~so33/SusDev/Lecture_9.pdf)





## Chapter 8

# Score, Gradient and Jacobian

### 8.1 Score

The score is the gradient (the vector of partial derivatives) of  $\log L(\theta)$ , with respect to an  $m$ -dimensional parameter vector  $\theta$ .

$$S(\theta) = \frac{\partial \ell}{\partial \theta}$$

Typically, they use  $\nabla$  to denote the partial derivative.

$$\nabla \ell$$

Such differentiation will generate a  $m \times 1$  row vector, which indicates the sensitivity of the likelihood.

Quote from Steffen Lauritzen's slides: "Generally the solution to this equation must be calculated by iterative methods. One of the most common methods is the Newton–Raphson method and this is based on successive approximations to the solution, using Taylor's theorem to approximate the equation."

For instance, using logit link, we can get the first derivative of log likelihood logistic regression as follows. We can not really find  $\beta$  easily to make the equation to be 0.

$$\begin{aligned}\frac{\partial \ell}{\partial \beta} &= \sum_{i=1}^n x_i^T \left[ y_i - \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}} \right] \\ &= \sum_{i=1}^n x_i^T [y_i - \hat{y}_i]\end{aligned}$$

## 8.2 Fisher scoring

[I will come back to this later.]

<https://www2.stat.duke.edu/courses/Fall00/sta216/handouts/diagnostics.pdf>

<https://stats.stackexchange.com/questions/176351/implement-fisher-scoring-for-linear-regression>

## 8.3 Gradient and Jacobian

**Remarks:** This part discusses gradient in a more general sense.

When  $f(x)$  is only in a single dimension space:

$$\mathbb{R}^n \rightarrow \mathbb{R}$$

$$\nabla f(x) = \left[ \frac{\partial f}{\partial x_1}, \frac{\partial f}{\partial x_2}, \dots, \frac{\partial f}{\partial x_n} \right]$$

When  $f(x)$  is only in a m-dimension space (i.e., Jacobian):  $\mathbb{R}^n \rightarrow \mathbb{R}^m$

$$Jac(f) = \begin{bmatrix} \frac{\partial f_1}{\partial x_1} & \frac{\partial f_1}{\partial x_2} & \frac{\partial f_1}{\partial x_3} & \dots & \frac{\partial f_1}{\partial x_n} \\ \frac{\partial f_2}{\partial x_1} & \frac{\partial f_2}{\partial x_2} & \frac{\partial f_2}{\partial x_3} & \dots & \frac{\partial f_2}{\partial x_n} \\ \dots & \dots & \dots & \dots & \dots \\ \frac{\partial f_m}{\partial x_1} & \frac{\partial f_m}{\partial x_2} & \frac{\partial f_m}{\partial x_3} & \dots & \frac{\partial f_m}{\partial x_n} \end{bmatrix}$$

For instance,

$$\mathbb{R}^n \rightarrow \mathbb{R}:$$

$$f(x, y) = x^2 + 2y$$

$$\nabla f(x, y) = \left[ \frac{\partial f}{\partial x}, \frac{\partial f}{\partial y} \right] = [2x, 2]$$

$$\mathbb{R}^n \rightarrow \mathbb{R}^m$$

$$f(x, y) = (x^2 + 2y, x^3)$$

$$Jac(f) = \begin{bmatrix} 2x & 2 \\ 2x^2 & 0 \end{bmatrix}$$

## 8.4 Hessian and Fisher Information

Hessian matrix or Hessian is a square matrix of second-order partial derivatives of a scalar-valued function, or scalar field.

$$\mathbb{R}^n \rightarrow \mathbb{R}$$

$$Hessian = \nabla^2(f) = \begin{bmatrix} \frac{\partial^2 f}{\partial x_1^2} & \frac{\partial^2 f}{\partial x_1 \partial x_2} & \frac{\partial^2 f}{\partial x_1 \partial x_3} & \cdots & \frac{\partial^2 f}{\partial x_1 \partial x_n} \\ \frac{\partial^2 f}{\partial x_2 \partial x_1} & \frac{\partial^2 f}{\partial x_2^2} & \frac{\partial^2 f}{\partial x_2 \partial x_3} & \cdots & \frac{\partial^2 f}{\partial x_2 \partial x_n} \\ \frac{\partial^2 f}{\partial x_3 \partial x_1} & \frac{\partial^2 f}{\partial x_3 \partial x_2} & \frac{\partial^2 f}{\partial x_3^2} & \cdots & \frac{\partial^2 f}{\partial x_3 \partial x_n} \\ \cdots & \cdots & \cdots & \cdots & \cdots \\ \frac{\partial^2 f}{\partial x_n \partial x_1} & \frac{\partial^2 f}{\partial x_n \partial x_2} & \frac{\partial^2 f}{\partial x_n \partial x_3} & \cdots & \frac{\partial^2 f}{\partial x_n^2} \end{bmatrix}$$

As a special case, in the context of logit:

Suppose that the log likelihood function is  $\ell(\theta)$ .  $\theta$  is a  $m$  dimension vector.

$$\theta = \begin{bmatrix} \theta_1 \\ \theta_2 \\ \theta_3 \\ \theta_4 \\ \cdots \\ \theta_m \end{bmatrix}$$

$$Hessian = \nabla^2(\ell) = \begin{bmatrix} \frac{\partial^2 \ell}{\partial \theta_1^2} & \frac{\partial^2 \ell}{\partial \theta_1 \partial \theta_2} & \frac{\partial^2 \ell}{\partial \theta_1 \partial \theta_3} & \cdots & \frac{\partial^2 \ell}{\partial \theta_1 \partial \theta_m} \\ \frac{\partial^2 \ell}{\partial \theta_2 \partial \theta_1} & \frac{\partial^2 \ell}{\partial \theta_2^2} & \frac{\partial^2 \ell}{\partial \theta_2 \partial \theta_3} & \cdots & \frac{\partial^2 \ell}{\partial \theta_2 \partial \theta_m} \\ \frac{\partial^2 \ell}{\partial \theta_3 \partial \theta_1} & \frac{\partial^2 \ell}{\partial \theta_3 \partial \theta_2} & \frac{\partial^2 \ell}{\partial \theta_3^2} & \cdots & \frac{\partial^2 \ell}{\partial \theta_3 \partial \theta_m} \\ \cdots & \cdots & \cdots & \cdots & \cdots \\ \frac{\partial^2 \ell}{\partial \theta_m \partial \theta_1} & \frac{\partial^2 \ell}{\partial \theta_m \partial \theta_2} & \frac{\partial^2 \ell}{\partial \theta_m \partial \theta_3} & \cdots & \frac{\partial^2 \ell}{\partial \theta_m \partial \theta_m} \end{bmatrix}$$

“In statistics, the observed information, or observed Fisher information, is the negative of the second derivative (the Hessian matrix) of the “log-likelihood” (the logarithm of the likelihood function). It is a sample-based version of the Fisher information.” (Direct quote from Wikipedia.)

Thus, the observed information matrix:

$$-Hessian = -\nabla^2(\ell)$$

Expected (Fisher) information matrix:

$$E[-\nabla^2(\ell)]$$

## Chapter 9

# Canonical link function

Inspired by a Stack Exchange post, I created the following figure:

$$\frac{\text{Paramter}}{\theta} \longrightarrow \gamma'(\theta) = \mu \longrightarrow \frac{\text{Mean}}{\mu} \longrightarrow g(\mu) = \eta \longrightarrow \frac{\text{Linearpredictor}}{\eta}$$

For the case of  $n$  time Bernoulli (i.e., Binomial), its canonical link function is logit. Specifically,

$$\frac{\text{Paramter}}{\theta = \beta^T x_i} \longrightarrow \gamma'(\theta) = \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}} \longrightarrow \frac{\text{Mean}}{\mu = \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}}} \longrightarrow g(\mu) = \log \frac{\frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}}}{1 - \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}}} \longrightarrow \frac{\text{Linearpredictor}}{\eta = \beta^T x_i}$$

Thus, we can see that,

$$\theta \equiv \eta$$

The link function  $g(\mu)$  relates the linear predictor  $\eta = \beta^T x_i$  to the mean  $\mu$ .

**Remarks:**

- (1) Parameter is  $\theta = \beta^T x_i$  (Not  $\mu$ !).
- (2)  $\mu = p(y = 1) = \frac{e^{\beta^T x_i}}{1 + e^{\beta^T x_i}}$  (Not logit!).
- (3) Link function (i.e.,  $g(\mu)$ ) = logit = logarithm of odds =  $\log \frac{\text{Event-Happened}}{\text{Event-Not-Happened}}$ .
- (4)  $g(\mu) = \log \frac{\mu}{1-\mu} = \beta^T x_i$ . Thus, link function = linear predictor = log odds!

- (5) Quote from the Stack Exchange post “Newton Method and Fisher scoring for finding the ML estimator coincide, these links simplify the derivation of the MLE.”

(Recall, we know that  $\mu$  or  $p(y = 1)$  is the mean function. Recall that,  $n$  trials of coin flips, and get  $p$  heads. Thus  $\mu = \frac{p}{n}$ .)

## Chapter 10

# Ordinary Least Squares (OLS)

Suppose we have  $n$  observation, and  $m$  variables.

$$\begin{bmatrix} x_{11} & x_{12} & x_{13} & \dots & x_{1m} \\ x_{21} & x_{22} & x_{23} & \dots & x_{2m} \\ \dots & & & & \\ x_{n1} & x_{n2} & x_{n3} & \dots & x_{nm} \end{bmatrix}$$

Thus, we can write it as the following  $n$  equations.

$$y_1 = \beta_0 + \beta_1 x_{11} + \beta_2 x_{12} + \dots + \beta_m x_{1m}$$

$$y_2 = \beta_0 + \beta_1 x_{21} + \beta_2 x_{22} + \dots + \beta_m x_{2m}$$

$$y_3 = \beta_0 + \beta_1 x_{31} + \beta_2 x_{32} + \dots + \beta_m x_{3m}$$

...

$$y_n = \beta_0 + \beta_1 x_{n1} + \beta_2 x_{n2} + \dots + \beta_m x_{nm}$$

We can combine all the  $n$  equations as the following one:

$$y_i = \beta_0 + \beta_1 x_{i1} + \beta_2 x_{i2} + \dots + \beta_m x_{im} (i \in [1, n])$$

We can further rewrite it as a matrix format as follows.

$$y = X\beta$$

Where,

$$y = \begin{bmatrix} y_1 \\ y_2 \\ y_3 \\ y_4 \\ \dots \\ y_n \end{bmatrix}$$

$$X = \begin{bmatrix} 1 & x_{11} & x_{12} & x_{13} & \dots & x_{1m} \\ 1 & x_{21} & x_{22} & x_{23} & \dots & x_{2m} \\ \dots & & & & & \\ 1 & x_{n1} & x_{n2} & x_{n3} & \dots & x_{nm} \end{bmatrix}$$

$$\beta = \begin{bmatrix} \beta_0 \\ \beta_1 \\ \beta_2 \\ \beta_3 \\ \dots \\ \beta_m \end{bmatrix}$$

Since later we need the inverse of  $X$ , we need to make it into a square matrix.

$$X^T y = X^T X \hat{\beta} \Rightarrow \hat{\beta} = (X^T X)^{-1} X^T y$$

We can use R to implement this calculation. As we can see, there is no need to do any iterations at all, but rather just pure matrix calculation.

```
X<-matrix(rnorm(1000),ncol=2) # we define a 2 column matrix, with 500 rows
X<-cbind(1,X) # add a 1 constant
beta_true<-c(2,1,2) # True regression coefficients
beta_true<-as.matrix(beta_true)
y=X%%beta_true+rnorm(500)

transposed_X<-t(X)
beta_hat<-solve(transposed_X%%X)%%transposed_X%%y
beta_hat

##           [,1]
## [1,] 1.929709
## [2,] 1.007278
## [3,] 1.952445
```

**Side Notes** The function of `as.matrix` will automatically make `c(2,1,2)` become the dimension of  $3 \times 1$ , you do not need to transpose the  $\beta$ .



## 10.1 Taylor series

$$\begin{aligned} f(x)|_a &= f(a) + \frac{f'(a)}{1!}(x-a) + \frac{f'(a)}{2!}(x-a)^2 + \frac{f''(a)}{3!}(x-a)^3 + \dots \\ &= \sum_{n=0}^{\infty} \frac{f^n(a)}{n!}(x-a)^n \end{aligned}$$

For example:

$$\begin{aligned} e^x|_{a=0} &= e^a + \frac{e^a}{1!}(x-a) + \frac{e^a}{2!}(x-a)^2 + \dots + \frac{e^a}{n!}(x-a)^n \\ &= 1 + \frac{1}{1!}x + \frac{1}{2!}x^2 + \dots + \frac{1}{n!}x^n \end{aligned}$$

if  $x = 2$

$$e^2 = 7.389056$$

$$e^2 \approx 1 + \frac{1}{1!}x = 1 + \frac{1}{1!}2 = 3$$

$$e^2 \approx 1 + \frac{1}{1!}x + \frac{1}{2!}x^2 = 1 + \frac{1}{1!}2 + \frac{1}{2!}2 = 5 \dots$$

$$e^2 \approx 1 + \frac{1}{1!}x + \frac{1}{2!}x^2 + \frac{1}{3!}x^2 + \frac{1}{4!}x^2 + \frac{1}{5!}x^2 = 7.2666\dots$$

## 10.2 References

1. Steffen Lauritzen's slides:

<http://www.stats.ox.ac.uk/~steffen/teaching/bs2HT9/scoring.pdf>

2. The Stack Exchange post:

<https://stats.stackexchange.com/questions/40876/what-is-the-difference-between-a-link-function-and-a-canonical-link-function>

3. Wikipedia for OLS

[https://en.wikipedia.org/wiki/Ordinary\\_least\\_squares](https://en.wikipedia.org/wiki/Ordinary_least_squares)

4. Gradient and Jacobian

<https://math.stackexchange.com/questions/1519367/difference-between-gradient-and-jacobian>

[https://www.youtube.com/watch?v=3xVMVT-2\\_t4](https://www.youtube.com/watch?v=3xVMVT-2_t4)

<https://math.stackexchange.com/questions/661195/what-is-the-difference-between-the-gradient-and-the-directional-derivative>

#### 5. Hessian

[https://en.wikipedia.org/wiki/Hessian\\_matrix](https://en.wikipedia.org/wiki/Hessian_matrix)

#### 6. Observed information

[https://en.wikipedia.org/wiki/Observed\\_information](https://en.wikipedia.org/wiki/Observed_information)

#### 7. Fisher information

[https://people.missouristate.edu/songfengzheng/Teaching/MTH541/Lecture%20notes/Fisher\\_info.pdf](https://people.missouristate.edu/songfengzheng/Teaching/MTH541/Lecture%20notes/Fisher_info.pdf)

#### 8. Link function

[https://en.wikipedia.org/wiki/Generalized\\_linear\\_model#Link\\_function](https://en.wikipedia.org/wiki/Generalized_linear_model#Link_function)

<https://stats.stackexchange.com/questions/40876/what-is-the-difference-between-a-link-function-and-a-canonical-link-function>

## Chapter 11

# Cholesky decomposition

### 11.1 Example 1

Use Cholesky decomposition to generate 1,000 trivariate normal deviates  $X_1, \dots, x_{1000}$  with mean  $\mu = (-2, 4, 3)$  and covariance matrix

$$X = \begin{bmatrix} 2 & -1 & 0.5 \\ -1 & 4 & 1 \\ 0.5 & 1 & 5 \end{bmatrix}$$

```
Nsim = 10
means = c(-2,4,3)
N_columns = 3

# Generating random standard normal distribution numbers
Generated_numbers = matrix(rnorm(N_columns * Nsim), nrow = N_columns)

# The provided covariance matrix
cov_matrix = rbind(c(2, -1, 0.5), c(-1, 4, 1), c(0.5, 1, 5))

# Cholesky decomposition
Cholesky_decom_results = chol(cov_matrix)

# Data is transformed using the Cholesky decomposition
adjusted_data = t(Generated_numbers) %*% Cholesky_decom_results

Final_data = t(t(adjusted_data) + means)
```

```

# calculating column means
colMeans(Final_data)

## [1] -2.022820  3.182124  2.178836

# calculating column variances
apply(Final_data,2,var)

## [1] 1.005080 2.099021 2.239064

# calculating covariance matrix
cov(Final_data)

##           [,1]      [,2]      [,3]
## [1,] 1.00508039 -0.13006437  0.06780402
## [2,] -0.13006437  2.09902097 -0.09608461
## [3,]  0.06780402 -0.09608461  2.23906434

```

## 11.2 Example 2

AR(1) Covariance Matrix with Correlation Rho and Variance SigmaSq. Note that, there is only one individual or participant in this data simulation.

```

n = 10;
SigmaSq = 5;
Rho = 0.8;

V = matrix(rep(n*n,0),n,n);

for (i in 1:n)
{
  for (j in i:n)
  {
    V[i,j]=SigmaSq*Rho^(j-i)
    V[j,i]=V[i,j]
  }
}

set.seed(123)
random_normal<-rnorm(n,2,1)
#chol(V) %*% random_normal

```

```
#colSums (chol(V))
b2<-t(as.matrix(random_normal))%*%chol(V)

pi = exp(b2)/(1 + exp(b2));

y<-ifelse(pi>runif(1),1,0)

y
```

```
##      [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [,10]
## [1,]    1    1    1    1    1    1    1    1    1    1
```

```
# The code above basically completes the generating job!
# The code below is to check
```

```
b = b2[2:n]
c = b2[1:(n-1)]
cor(b,c)
```

```
## [1] 0.8967058
```

```
sd(as.vector(b2))
```

```
## [1] 3.535119
```

```
# note that, you can not use var, as the mean is not zero, but rather it is 2
var(as.vector(b2))
```

```
## [1] 12.49707
```

```
#Not sure why the means are not the same ?
mean(as.vector(b2))
```

```
## [1] 10.01925
```

```
mean(random_normal)
```

```
## [1] 2.074626
```

### 11.3 Example 3

The following code very similar to the code shown above. However, it had only one observation. To illustrate the situation where there are more than one individual (or, participant), I did the code below.

```
n =25;    #the number of time points
m= 15;    # the number of participants or individuals, whichever ways you would like to
SigmaSq = 5;
Rho = 0.8;

filling_numbers<-rep(n*n,0)
V = matrix(filling_numbers,n,n);

for (i in 1:n)
{
  for (j in i:n)
  {
    V[i,j]=SigmaSq*Rho^(j-i)
    V[j,i]=V[i,j]
  }
}

set.seed(2345)
random_normal<-matrix(rnorm(m*n),nrow = m)
#chol(V) %%% random_normal
#colSums (chol(V))
b2<-random_normal%%chol(V)

pi = exp(b2)/(1 + exp(b2));

random_unfirom<-matrix(runif(m*n),nrow = m)

y<-ifelse(pi>random_unfirom,1,0)
y
```

```
##      [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8] [,9] [,10] [,11] [,12] [,13]
## [1,]  0    0    0    0    0    0    1    1    1    1    1    0    1
## [2,]  0    1    1    1    1    1    0    1    0    0    0    0    0
## [3,]  1    0    0    0    0    1    0    1    0    1    1    0    0
## [4,]  1    0    0    0    0    0    0    0    1    0    0    0    0
## [5,]  0    0    0    0    1    0    0    0    0    0    0    0    1
## [6,]  0    0    1    1    0    0    0    1    0    0    0    0    0
## [7,]  0    0    0    0    0    1    0    0    1    0    0    0    1
## [8,]  1    1    1    1    0    0    0    0    0    1    1    0    0
```

```
## [9,] 1 1 1 1 0 1 1 1 0 1 0 0 0
## [10,] 1 0 1 1 1 1 1 1 1 1 0 0 0
## [11,] 1 1 1 1 1 1 1 1 1 0 1 0 0
## [12,] 1 1 1 1 1 1 0 1 1 1 1 1 0
## [13,] 0 1 0 0 0 0 0 0 1 1 1 1 0
## [14,] 1 1 0 1 1 1 1 1 1 0 0 0 0
## [15,] 1 1 0 1 1 1 1 0 0 1 0 0 1
##      [,14] [,15] [,16] [,17] [,18] [,19] [,20] [,21] [,22] [,23] [,24] [,25]
## [1,] 1 1 1 1 0 1 0 0 1 0 0 0
## [2,] 0 1 1 1 1 1 0 0 0 1 0 1
## [3,] 0 1 0 0 1 1 1 1 0 0 1 1
## [4,] 1 0 0 0 0 0 1 0 1 0 0 0
## [5,] 0 0 1 1 1 0 1 1 1 0 1 1
## [6,] 1 1 0 0 1 0 0 1 0 1 1 1
## [7,] 0 0 0 0 0 1 0 0 1 0 0 0
## [8,] 0 0 0 0 0 0 0 0 0 1 1 1
## [9,] 0 1 1 1 0 0 1 0 0 0 0 0
## [10,] 1 1 1 1 1 1 1 1 1 1 1 1
## [11,] 0 0 0 1 0 0 0 0 1 1 1 0
## [12,] 0 0 0 0 1 1 1 0 0 0 1 0
## [13,] 0 1 0 0 0 1 0 1 1 1 1 1
## [14,] 0 0 0 0 1 0 1 0 0 0 0 0
## [15,] 0 0 0 0 0 0 1 0 1 0 0 1
```

*# The code above basically completes the generating job! The code below is to check*

```
# The following calculates variance
# calculate variance of each column
mean(apply(b2, 2, var))
```

```
## [1] 4.330903
```

```
# calculate variance of each row
mean(apply(b2, 1, var))
```

```
## [1] 3.568107
```

```
# The whole table
var(as.vector(b2))
```

```
## [1] 4.299165
```

```

# The following code calculates the correlation
b = b2[,2:n]
c = b2[,1:(n-1)]

collected_cor<-rep(0,m-1) #creating an empty vector to collect correlation.
for (i in 1:(m-1))
{collected_cor[i]<-cor(b[i,],c[i,])}
collected_cor

## [1] 0.8473037 0.7065013 0.6376223 0.5481540 0.7851062 0.6576329 0.4844481
## [8] 0.6950847 0.6731673 0.6409116 0.7966547 0.7184030 0.8001861 0.7913736

mean(collected_cor)

## [1] 0.6987535

mean(y)

## [1] 0.456

log(mean(y)/(1-mean(y)))

## [1] -0.1764564

# It will always get a value close to zero, since we set the mean to be zero when simu

```