

Statistical Geophysics

Chapter 2

Inferential Statistics 1



Parameter estimation

- We will be studying problems of **statistical inference**.
- Many problems of inference have been dichotomized into two areas: **estimation of parameters** and **tests of hypotheses**.
- Parameter estimation: Let X be a random variable, whose density is $f_X(x; \theta)$, where the form of the density is assumed known except that it contains an unknown parameter θ .
- The problem is then to use the observed values x_1, \dots, x_n of a random sample X_1, \dots, X_n to **estimate** the value of θ or the value of some function of θ , say $\tau(\theta)$.

Estimator and estimate

- Any statistic $T = g(X_1, \dots, X_n)$ whose values are used to estimate θ is defined to be an **estimator** of θ .
- That is, T is a known function of observable random variables that is itself a random variable.
- An **estimate** is the realized value $t = g(x_1, \dots, x_n)$ of an estimator, which is a function of the realized values x_1, \dots, x_n .
- Example:** $\bar{X}_n = \frac{1}{n} \sum_{i=1}^n X_i$ is an estimator of a mean μ and \bar{x}_n is an estimate of μ . Here, T is \bar{X}_n , t is \bar{x}_n and $g(\cdot)$ is the function defined by summing the arguments and then dividing by n .

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Estimation by Maximum Likelihood

Background

- In 1921, **R. A. Fisher** pointed out an attractive rationale, called **maximum likelihood** (ML), for estimating parameters.
- This procedure says one should examine the **likelihood function** of the sample values and take as the estimates of the unknown parameters those values that **maximize** this likelihood function.
- ML is unifying concept to cover a broad range of problems.
- It is generally accepted as the best rationale to apply in estimating parameters, when one is willing to assume the form of the population probability law is known.

Likelihood function

- If X_1, \dots, X_n are an i.i.d. sample from a population with pdf or pmf $f(x|\theta)$, the **likelihood function** is defined by

$$L(\theta) = L(\theta|x_1, \dots, x_n) = \prod_{i=1}^n f(x_i|\theta) .$$

- **Maximum likelihood principle:** Given x_1, \dots, x_n take as the estimate of θ the value $\hat{\theta}$ that maximizes the likelihood, that is,

$$L(\hat{\theta}) = \max_{\theta} L(\theta) .$$

- The value $\hat{\theta}$ that maximizes the likelihood is called the maximum likelihood estimate (MLE) for θ .

Log-likelihood and score function

- It is often more convenient to work with the logarithm of the likelihood function, called the **log-likelihood**:

$$l(\theta) = \ln L(\theta) = \sum_{i=1}^n \ln f(x_i|\theta) .$$

- If the log-likelihood is differentiable (in θ), possible candidates for the MLE are the values that solve

$$s(\theta) = \frac{\partial}{\partial \theta} l(\theta) = 0 .$$

- The first derivative of the log-likelihood is called the **score function**.

Example

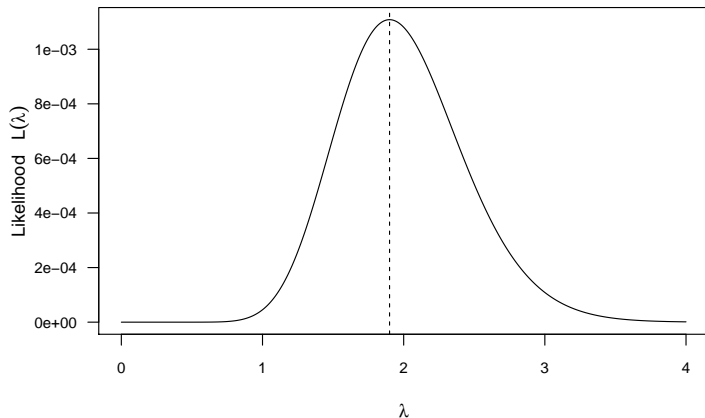
- Let x_1, \dots, x_n be realizations from $X_i \stackrel{i.i.d.}{\sim} \mathcal{P}(\lambda)$ ($i = 1, \dots, n$) with unknown parameter λ .
- The aim is to estimate λ by maximum likelihood.
- Likelihood function:

$$\begin{aligned} L(\lambda) &= f(x_1, \dots, x_n | \lambda) \\ &= f(x_1 | \lambda) \cdots f(x_n | \lambda) \\ &= \prod_{i=1}^n f(x_i | \lambda) \\ &= \prod_{i=1}^n \left(\frac{\lambda^{x_i}}{x_i!} \exp(-\lambda) \right) . \end{aligned}$$

Example

Likelihood for i.i.d. sample of $n=10$ from $X \sim \text{Pois}(\lambda=2)$

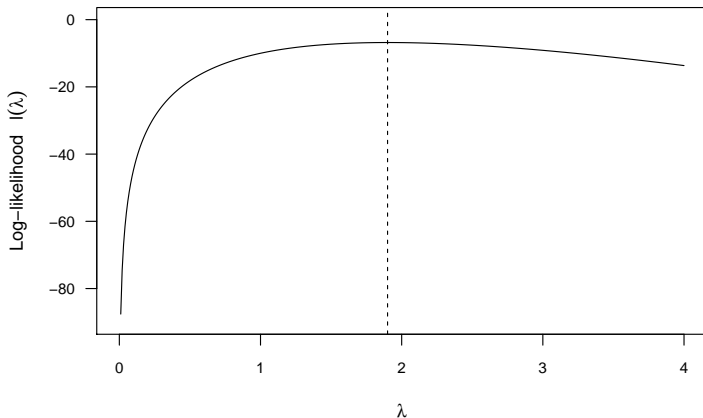
$x_1 = 1$ $x_2 = 1$ $x_3 = 0$ $x_4 = 3$ $x_5 = 1$ $x_6 = 4$ $x_7 = 1$ $x_8 = 1$ $x_9 = 2$ $x_{10} = 5$



Example

Log-likelihood for i.i.d. sample of $n=10$ from $X \sim \text{Pois}(\lambda=2)$

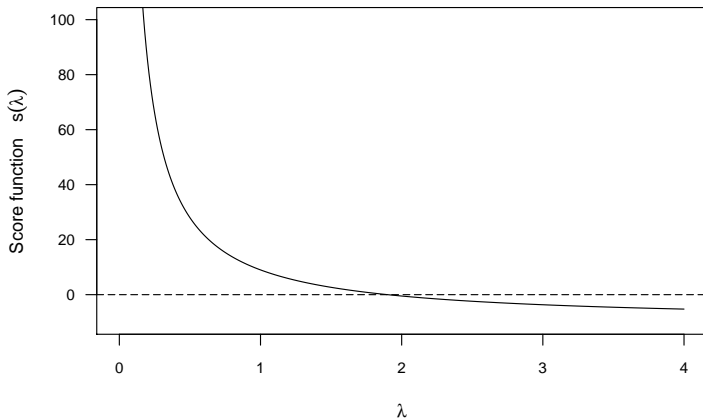
$x_1 = 1$ $x_2 = 1$ $x_3 = 0$ $x_4 = 3$ $x_5 = 1$ $x_6 = 4$ $x_7 = 1$ $x_8 = 1$ $x_9 = 2$ $x_{10} = 5$



Example

Score function for i.i.d. sample of $n=10$ from $X \sim \text{Pois}(\lambda=2)$

$x_1 = 1$ $x_2 = 1$ $x_3 = 0$ $x_4 = 3$ $x_5 = 1$ $x_6 = 4$ $x_7 = 1$ $x_8 = 1$ $x_9 = 2$ $x_{10} = 5$



Numerical optimization

Newton-Raphson method

- Suppose that we want to approximate the solution to $s(\theta) = 0$.
- Let us also suppose that we have somehow found an initial approximation to this solution, say $\theta^{(0)}$.
- If $\theta^{(k)}$ is an approximation to $s(\theta) = 0$ and if $s'(\theta^{(k)}) \neq 0$, the next approximation is given by

$$\theta^{(k+1)} = \theta^{(k)} - \frac{1}{s'(\theta^{(k)})} \cdot s(\theta^{(k)}) .$$

- This iterative scheme continues until a prespecified convergence criterion is met.

Other estimation methods

- The **method of moments** uses sample moments to estimate the parameters of an assumed probability law.
- **Least squares estimation** minimizes the sum of the squares of the deviations of the observed values and the fitted values.
- **Bayesian estimation** is based on combining the evidence contained in the data with prior knowledge, based on **subjective probabilities**, of the values of unknown parameters.

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Properties of Estimators

Evaluating estimators

- We have outlined reasonable techniques for finding out estimators of parameters.
- Are some of many possible estimators **better** in some sense, than others?
- When we are faced with the choice of two or more estimators for the same parameter, it becomes important to develop **criteria** for comparing them.
- We will now define certain **properties**, which an estimator may or may not possess, that will help us in deciding whether one estimator is better than another.

Unbiasedness

Definition:

- An estimator $T = g(X_1, \dots, X_n)$ is defined to be an **unbiased** estimator of an unknown parameter θ if and only if

$$E(T) = \theta \text{ for all values of } \theta.$$

- The difference $E(T) - \theta$ is called the **bias** of T and can be either positive, negative, or zero.
- An estimator T of θ is said to be **asymptotically unbiased** if

$$\lim_{n \rightarrow \infty} E(T) = \theta .$$

Precision of estimation

- For observations x_1, \dots, x_n an estimator T yields an estimate $t = g(x_1, \dots, x_n)$.
- In general, the estimate will not be equal to θ .
- For unbiased estimators the **precision** of the estimation method is captured by the variance of the estimator, $\text{Var}(T)$.
- The square root of $\text{Var}(T)$ (the standard deviation of T) is called the **standard error**, which in general has to be estimated itself.

Lower bound for variance

- Let X be a random variable with density $f(x, \theta)$. Under certain **regularity conditions**:

$$\text{Var}(T) \geq \frac{1}{n\mathbb{E} \left[\left(\frac{\partial}{\partial \theta} \ln f(x, \theta) \right)^2 \right]} ,$$

where T is an unbiased estimator of θ .

- The equation above is called the **Cramér-Rao inequality**, and the right-hand side is called the **Cramér-Rao lower bound** for the variance of unbiased estimators of θ .

Mean-squared error

Definition:

- The **mean-squared error** (MSE) of $T = g(X_1, \dots, X_n)$ (as an estimator for θ) is

$$\text{MSE}(T) = \text{E}[(T - \theta)^2] = \text{Var}(T) + (\text{E}(T) - \theta)^2 .$$

- Suppose T is an unbiased estimator of θ , then $\text{MSE}(T) = \text{Var}(T)$.

Consistency

Definition:

- Let $T = g(X_1, \dots, X_n)$ be an estimator for θ . Then, T is a **consistent** estimator for θ if

$$\lim_{n \rightarrow \infty} P(|T - \theta| \geq \epsilon) = 0 \text{ for any } \epsilon > 0 .$$

- From the **Chebyshev inequality** we know that

$$\begin{aligned} P(|T - \theta| \geq \epsilon) &\leq \frac{1}{\epsilon^2} E[(T - \theta)^2] \\ &= \frac{1}{\epsilon^2} \text{MSE}(T) . \end{aligned}$$

- It follows that if $\text{MSE}(T) \rightarrow 0$ as $n \rightarrow \infty$, then T is consistent.

Efficiency

Definition:

- If T_1 and T_2 are two estimators of θ , then T_1 is **more efficient** than T_2 if

$$\text{MSE}(T_1) \leq \text{MSE}(T_2) \text{ for any value of } \theta$$

with strict inequality holding somewhere.

- For two **unbiased** estimators T_1 and T_2 of θ , T_1 is **more efficient** than T_2 if

$$\text{Var}(T_1) \leq \text{Var}(T_2) \text{ for any value of } \theta$$

with strict inequality holding somewhere.

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Confidence Intervals

Interval estimation

- So far, we have dealt with the **point estimation** of a parameter.
- It seems desirable that a point estimate should be accompanied by some measure of the possible error of the estimate.
- We might make the inference of estimating that the true value of the parameter is contained in some interval.
- **Interval estimation**: Define two statistics $T_1 = g_1(X_1, \dots, X_n)$ and $T_2 = g_2(X_1, \dots, X_n)$, where $T_1 \leq T_2$, so that $[T_1, T_2]$ constitutes an interval for which the probability can be determined that it contains the unknown θ .

Confidence interval

Definition:

- Given a random sample X_1, \dots, X_n let $T_1 = g_1(X_1, \dots, X_n)$ and $T_2 = g_2(X_1, \dots, X_n)$ be two statistics satisfying $T_1 \leq T_2$ for which

$$P(T_1 \leq \theta \leq T_2) = 1 - \alpha .$$

- Then the random interval $[T_1, T_2]$ is called a $(1 - \alpha)$ -confidence interval for θ .
- $1 - \alpha$ is called the confidence coefficient and T_1 and T_2 are called the lower and upper confidence limits, respectively.
- A value $[t_1, t_2]$, where $t_j = g_j(x_1, \dots, x_n)$ ($j = 1, 2$) is an observed $(1 - \alpha)$ -confidence interval for θ .

One-sided confidence interval

Definition:

- Let $T_1 = -\infty$ and $T_2 = g_2(X_1, \dots, X_n)$ be a statistic for which

$$P(\theta \leq T_2) = 1 - \alpha .$$

- Then T_2 is called a **one-sided upper confidence limit** for θ .
- Similarly, let $T_2 = \infty$ and $T_1 = g_1(X_1, \dots, X_n)$ be a statistic for which

$$P(T_1 \leq \theta) = 1 - \alpha .$$

- Then T_1 is called a **one-sided lower confidence limit** for θ .

Confidence intervals for the mean (σ^2 known)

100(1 - α) %-confidence interval for μ (scenario σ^2 known)

- For a normally distributed random variable X :

$$\left[\bar{X} - z_{1-\alpha/2} \frac{\sigma}{\sqrt{n}}, \bar{X} + z_{1-\alpha/2} \frac{\sigma}{\sqrt{n}} \right] .$$

- For an arbitrarily distributed random variable X and $n > 30$,

$$\left[\bar{X} - z_{1-\alpha/2} \frac{\sigma}{\sqrt{n}}, \bar{X} + z_{1-\alpha/2} \frac{\sigma}{\sqrt{n}} \right]$$

is an **approximate** confidence interval for μ .

- For $0 < p < 1$, z_p is the p -quantile of the standard normal distribution, that is, it is the value for which $F(z_p) = \Phi(z_p) = p$. Hence, $z_p = \Phi^{-1}(p)$.

Confidence intervals for the mean (σ^2 unknown)

100(1 - α) %-confidence interval for μ (scenario σ^2 unknown)

- For a normally distributed random variable X :

$$\left[\bar{X} - t_{1-\alpha/2}(n-1) \frac{S}{\sqrt{n}}, \bar{X} + t_{1-\alpha/2}(n-1) \frac{S}{\sqrt{n}} \right],$$

where $S = \sqrt{\frac{1}{n-1} \sum_{i=1}^n (X_i - \bar{X})^2}$ and $t_{1-\alpha/2}(n-1)$ being the $(1 - \alpha/2)$ -quantile of the **t-distribution** with $n - 1$ degrees of freedom.

- For an arbitrarily distributed random variable X and $n > 30$,

$$\left[\bar{X} - z_{1-\alpha/2} \frac{S}{\sqrt{n}}, \bar{X} + z_{1-\alpha/2} \frac{S}{\sqrt{n}} \right]$$

is an **approximate** confidence interval for μ .

Confidence intervals for the variance

100(1 - α) %-confidence interval for σ^2

- For a normally distributed random variable X :

$$\left[\frac{(n-1)S^2}{\chi_{1-\alpha/2}^2(n-1)}, \frac{(n-1)S^2}{\chi_{\alpha/2}^2(n-1)} \right],$$

where $\chi_{1-\alpha/2}^2(n-1)$ and $\chi_{\alpha/2}^2(n-1)$ denote the $(1 - \alpha/2)$ -quantile and $(\alpha/2)$ -quantile, respectively, of the **chi-square distribution** with $n - 1$ degrees of freedom.

Confidence interval for a proportion

100(1 - α) %-confidence interval for π

- In **dichotomous** populations and for $n > 30$, an **approximate** confidence interval for $\pi = P(X = 1)$ is given by

$$\left[\hat{\pi} - z_{1-\alpha/2} \sqrt{\frac{\hat{\pi}(1 - \hat{\pi})}{n}}, \hat{\pi} + z_{1-\alpha/2} \sqrt{\frac{\hat{\pi}(1 - \hat{\pi})}{n}} \right],$$

where $\hat{\pi} = \bar{X}$ denotes the relative frequency.