

ABSTRACT

Proposal for an Adjusted Numerical Approximation to the Integrated Likelihood Function

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The primary focus of this prospectus is to motivate and explain an adapted version of numerical likelihood integration as a method for eliminating nuisance parameters from a statistical model.

Table of Contents

ABSTRACT	1
Table of Contents	2
Chapter 1. Introduction and Motivation	4
Chapter 2. Experiments and Statistical Models	5
Chapter 3. The Likelihood Function	8
3.1. Definition	8
3.2. Transformations	9
3.3. Maximum Likelihood Estimation	12
3.4. Regularity Conditions	13
3.5. The Bartlett Identities	14
3.6. One-Index Asymptotics	16
Chapter 4. Pseudolikelihood Functions	20
4.1. Model Parameter Decomposition	20
4.2. Types of Pseudolikelihoods	22
4.3. The Bartlett Identities Revisited	23
Chapter 5. Approximating the Integrated Likelihood Function	24
5.1. The Zero-Score Expectation Parameter	24
5.2. Two-Index Asymptotics	26
Chapter 6. Applications	31

	3
6.1. Multinomial Distribution	31
6.2. Standardized Mean Difference	31
References	32
Appendix A. Chapter 3	34
A.1. Definitions and Notation	34
A.2. Taylor's Theorem	35
Appendix B. Chapter 5	36
B.1. Desirable Properties of the Integrated Likelihood	36
B.2. Laplace's Method	36

CHAPTER 1

Introduction and Motivation

CHAPTER 2

Experiments and Statistical Models

The acquisition of knowledge regarding a population of interest has long been the impetus for the field of statistical inference. In all but the most basic of circumstances, limiting constraints such as time, accessibility, and cost make perfect knowledge of a population essentially impossible to obtain. It therefore becomes necessary to infer characteristics of the population based on a random and representative sample of observations drawn from it. The procedures by which these samples might be procured are themselves far from trivial, and indeed an entire branch of statistics has been dedicated to their study. However, we are primarily concerned in this paper with what occurs after the sample has been taken, and so we will generally take it for granted that a suitably representative sample of the population already exists.

Suppose (x_1, \dots, x_n) is one such sample. What information can we then glean from this sample about the population from which it has been drawn? Where is its point of central tendency located? Are its values clustered tightly around this point, or are they more diffuse? Are they distributed symmetrically or skewed to one side or the other? As the questions increase in complexity, so do the techniques required to answer them. Unfortunately the natural chaos of the real world all but guarantees there will never be an instrument capable of completely capturing the intricacies of a population whose properties we wish to infer. Hence, some amount of idealization will always be required in order to proceed.

This idealization typically comes in the form of additional assumptions that we impose on the population with the goal of sacrificing what we hope is only a small amount of accuracy in exchange for a large reduction in complexity. These assumptions are essentially never “true” in the sense that they are not a flawless representation of reality, but they may nevertheless serve as convenient approximations that are capable of producing answers with degrees of accuracy high enough to be useful in their own right. Taken as a whole, they form the basis for a statistical model.

A statistical model is best understood through the lens of what is known as an *experiment*. From a probabilistic perspective, an experiment refers to any procedure that could theoretically be replicated an infinite number of times and in which each replication results in exactly one outcome taken. The set of possible outcomes of an experiment, known as the *sample space*, must be well-defined, meaning that

assuming that there exists an unknown probability distribution P over the population of interest that generates the data (x_1, \dots, x_n) we observe from it.

We can think of this observed data as being the realized outcomes (“realizations”) of some random variable X that is distributed according to P . We will restrict our attention in this paper to distributions that are either absolutely continuous or discrete, meaning they admit a probability density or probability mass function, respectively, over their support. For the sake of brevity, we will refer to all such functions as density functions, with the understanding that it is actually a probability mass function in the event that X is discrete. Let $p(x)$ denote the unknown density function associated with P and \mathcal{P} the set of functions that we are willing to consider as candidates for $p(x)$. Out of necessity, we will proceed as though our choice of \mathcal{P} always contains $p(x)$ though in reality there is nothing specifically requiring it.

We will also assume the set \mathcal{P} is *parameterized*. That is, we assume there exists a *parameter* θ which indexes \mathcal{P} , acting as a label that allows us to differentiate between the densities it contains. For a particular value of the parameter θ , say θ_1 , we can refer to its corresponding density in \mathcal{P} with the notation $p(\cdot; \theta_1)$, and therefore \mathcal{P} itself may be written as $\mathcal{P} = \{p(\cdot; \theta) | \theta \in \Theta\}$. Θ is called the *parameter space* and represents the set of all possible values θ can take on.

In general, a model’s parameterization is not unique, and for a given parameter θ , we are free to choose any one-to-one function of θ as a new parameter. Once we have made our choice of parameterization, we will assume that Θ does contain a singular true parameter value, which we will denote by θ_0 . The conventional interpretation of θ_0 is as a fixed but unknown constant that represents the value of the parameter corresponding to the true density function in \mathcal{P} . Conversely, θ represents an arbitrary parameter value (corresponding to an arbitrary density in \mathcal{P}) that is allowed to range over all possible elements of Θ , including θ_0 . In other words, θ acts like a tuning dial for the population - rotate the dial and certain behaviors of the population (e.g. its location, scale, or shape) will change. Making inferences

regarding θ_0 is like trying to figure out the particular value (θ_0) to which a population's dial (θ) has been set. Note that it is possible for the value of θ_0 itself to change over time as well, depending on the population. In such cases, any estimate of θ_0 based on a cross-sectional sample drawn from the population is best thought of as an estimate of the true parameter value during the particular time in which the sample was collected.

Crucially, it must always be possible to identify the parameter in our model on the basis of the data we observe. A model is considered *identifiable* if having perfect knowledge of the population would enable us to determine θ_0 with absolute certainty. This is equivalent to requiring that for some observed data x and any two parameters $\theta_1, \theta_2 \in \Theta$, if $p(x; \theta_1) = p(x; \theta_2)$, then it must follow that $\theta_1 = \theta_2$. A model that is not identifiable could potentially have two or more distinct parameter values that give rise to the same probability distribution. For example, suppose Y is distributed uniformly on the interval $(0, \alpha + \beta)$, where $\alpha, \beta > 0$. If we use $\theta = (\alpha, \beta)$ as a parameter for the distribution of Y , then θ is unidentifiable since, for instance, the case where $\theta_1 = (0, 1)$ and $\theta_2 = (1, 0)$ implies that $p(y; \theta_1) = p(y; \theta_2)$ despite the fact that $\theta_1 \neq \theta_2$. This is obviously an undesirable property for a model to possess, and so we will consider only identifiable models in this paper as a means of avoiding it.

Finally, we must make a choice regarding the dimension of the parameter space Θ when formulating our models. *Parametric* models are defined as having finite-dimensional parameter spaces. Any model that is not parametric is called *nonparametric*. We will restrict our attention in this paper solely to parametric models whose parameter spaces are subsets of \mathbb{R}^d , where $d \in \mathbb{Z}^+$.

CHAPTER 3

The Likelihood Function**3.1. Definition**

Upon choosing a statistical model that we think best characterizes our population of interest, the obvious next step is to identify the true distribution in \mathcal{P} or at the very least, the one that best approximates the truth. This is equivalent to making inferences about θ_0 in the case where the model is parametric and identifiable. That is, given the particular form(s) we have chosen for the distributions in \mathcal{P} , the only unknown remaining is the value of θ_0 itself. Since this value is ultimately what controls the mechanism generating any sample of data $\mathbf{x} = (x_1, \dots, x_n)$ that we might observe from the population, it stands to reason that information regarding θ_0 can be inferred from the specific values of x_1, \dots, x_n that we obtain. To make this notion more rigorous, we require some method of analyzing the joint probability of our sample as a function of our parameter θ .

Given some observed data $X = \mathbf{x}$, the *likelihood function* for θ is defined as

$$(3.1.1) \quad L(\theta) \equiv L(\theta; \mathbf{x}) = p(\mathbf{x}; \theta), \quad \theta \in \Theta.$$

In other words, the value of the likelihood function evaluated at a particular $\theta \in \Theta$ is simply equal to the output of the model's density function evaluated at the same inputs. However, while $p(\mathbf{x}; \theta)$ is viewed primarily as a function of \mathbf{x} for fixed θ , the reverse is actually true for $L(\theta; \mathbf{x})$. Indeed, we regard the likelihood as being a function of the parameter θ for fixed \mathbf{x} . The positioning of the arguments θ and x is a reflection of this difference in perspectives.

When X is discrete, we may interpret $L(\theta; x)$ as the probability that $X = \mathbf{x}$ given that θ is the true parameter value. Crucially, this is *not* equivalent to the inverse probability that θ is the true parameter value given $X = \mathbf{x}$. The likelihood does not directly tell us anything about the probability that θ assumes

any particular value at all. Though intuitively appealing, this interpretation constitutes a fundamental misunderstanding of what a likelihood function is, and great care must be taken to avoid it.

When X is continuous, the likelihood for θ may still be defined as it is in Equation 3.1.1. However, we must forfeit our previous interpretation of $L(\theta)$ as a probability since the probability that X takes on any particular value is now 0. We may however still think of the likelihood as being proportional to the probability that X takes on a value “close” to x , meaning that that X is within a tiny neighborhood of x . Specifically, for two different samples x_1 and x_2 , if $L(\theta; \mathbf{x}_1) = c \cdot L(\theta; \mathbf{x}_2)$, where $c > 1$, then under this model we may conclude X is c times more likely to assume a value closer to \mathbf{x}_1 than \mathbf{x}_2 given that θ is the true value of the parameter.

As in the discrete case, we must also be careful when X is continuous to avoid using $L(\theta; \mathbf{x})$ to make probabilistic assertions regarding θ . Despite our use of probability in its definition, the likelihood itself is *not* a probability density function for the parameter θ and is subject to neither the same rules nor interpretations as one.

3.2. Transformations

There are a few useful transformations of the likelihood function that we will define here for use in future sections. The first is the *log-likelihood function*, which is defined as the natural logarithm of the likelihood function:

$$(3.2.1) \quad \ell(\theta) \equiv \ell(\theta; \mathbf{x}) = \log L(\theta; \mathbf{x}), \quad \theta \in \Theta.$$

In practice, we will typically eschew direct analysis of the likelihood in favor of the log-likelihood due to the nice mathematical properties logarithms possess. Chief among these properties is the ability to turn products into sums (i.e. $\log(ab) = \log(a) + \log(b)$ for $a, b > 0$). Sums tend to be easier to differentiate than products, making this a particularly useful feature for likelihood functions, which are often expressed as the product of marginal density functions when the observations are independent.

The other key property of logarithms that makes the log-likelihood so useful is that they are strictly increasing functions of their arguments (i.e. $\log x > \log y$ for $x > y > 0$). This monotonicity ensures

that the locations of a function's extrema are preserved when the function is passed to the argument of a logarithm. For example, for a positive function f with a global maximum, $\operatorname{argmax}_x f(x) = \operatorname{argmax}_x \log f(x)$.

In the general case in which θ is a d -dimensional vector, where d is an integer greater than 1, it follows that the first derivative of the log-likelihood with respect to θ will also be a d -dimensional vector, the second derivative will be a $d \times d$ matrix, the third derivative will be a $d \times d \times d$ array, and so forth. Consequently, we will use gradient notation (i.e. ∇ , ∇^2 , etc.) to emphasize the vector nature of these results.

The gradient of ℓ with respect to θ appears frequently enough in the analysis of likelihood functions that it has earned its own name - the *score function*. Formally, it is defined as

$$(3.2.2) \quad \mathcal{S}(\theta) \equiv \mathcal{S}(\theta; \mathbf{x}) = \nabla \ell(\theta; \mathbf{x}) = \begin{pmatrix} \frac{\partial \ell}{\partial \theta_1} \\ \vdots \\ \frac{\partial \ell}{\partial \theta_d} \end{pmatrix}, \quad \theta \in \Theta.$$

Similarly, the negative Hessian matrix of the log-likelihood function with respect to θ is called the *observed information* or just the *information* and is denoted by

$$(3.2.3) \quad \mathcal{J}(\theta) \equiv \mathcal{J}(\theta; \mathbf{x}) = -\nabla^2 \ell(\theta; \mathbf{x}) = - \begin{pmatrix} \frac{\partial^2 \ell}{\partial \theta_1^2} & \frac{\partial^2 \ell}{\partial \theta_1 \partial \theta_2} & \cdots & \frac{\partial^2 \ell}{\partial \theta_1 \partial \theta_d} \\ \frac{\partial^2 \ell}{\partial \theta_2 \partial \theta_1} & \frac{\partial^2 \ell}{\partial \theta_2^2} & \cdots & \frac{\partial^2 \ell}{\partial \theta_2 \partial \theta_d} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial^2 \ell}{\partial \theta_d \partial \theta_1} & \frac{\partial^2 \ell}{\partial \theta_d \partial \theta_2} & \cdots & \frac{\partial^2 \ell}{\partial \theta_d^2} \end{pmatrix}, \quad \theta \in \Theta.$$

Note that this is also equal to the negative gradient of the score function. The use of the term “information” here derives from the fact that the second partial derivatives of ℓ with respect to the components of θ are all related to the curvature of ℓ near its maximum - the sharper the curve, the less uncertainty and therefore more information we have about θ .

Recall that $L(\theta; x_1, \dots, x_n)$ is defined as a function of θ for a fixed sample of observations x_1, \dots, x_n , where we think of each x_i as being a realization of a random variable X_i . We may therefore interpret $L(\theta; x_1, \dots, x_n)$ as a random variable in the following sense: for a given θ , the value of $L(\theta; x_1, \dots, x_n)$

depends entirely on the values of X_1, \dots, X_n that we happened to observe, and so $L(\theta; x_1, \dots, x_n)$ is itself a random variable with respect to the joint probability distribution of X_1, \dots, X_n . The same is also true for any function or estimate based on the likelihood, as they ultimately will all depend on the data through it as well.

The random nature of these likelihood-based quantities further implies that finding their expectations and variances respect to $p(x_1, \dots, x_n; \theta_0)$ is a well-defined, nontrivial task. The variance of the score function in particular will be useful to know, as it also relates to the amount of information pertaining to θ_0 that is contained within the log-likelihood function of our model. Properly known as the *Fisher information* or the *expected information*, it is defined as

$$(3.2.4) \quad \mathcal{J}(\theta) = \mathbb{V}[\mathcal{S}(\theta; \mathbf{x}); \theta_0], \quad \theta \in \Theta.$$

However, since we are working in the more general framework in which $\mathcal{S}(\theta)$ is a $d \times 1$ random vector, it would be more accurate to speak of the *Fisher information matrix*, which is equal to the variance-covariance matrix of $\mathcal{S}(\theta)$. Hence, we have

$$\begin{aligned} \mathcal{J}(\theta) &= \mathbb{V}[\mathcal{S}(\theta; \mathbf{x}); \theta_0] && \text{(by Eq. 2.2.4)} \\ &= \text{Cov} \left[\left(\frac{\partial \ell}{\partial \theta_1} \dots \frac{\partial \ell}{\partial \theta_d} \right)^T; \theta_0 \right] && \text{(by Eq. 2.2.2)} \\ &= \begin{pmatrix} \mathbb{V} \left(\frac{\partial \ell}{\partial \theta_1} \right) & \text{Cov} \left(\frac{\partial \ell}{\partial \theta_1}, \frac{\partial \ell}{\partial \theta_2} \right) & \dots & \text{Cov} \left(\frac{\partial \ell}{\partial \theta_1}, \frac{\partial \ell}{\partial \theta_d} \right) \\ \text{Cov} \left(\frac{\partial \ell}{\partial \theta_2}, \frac{\partial \ell}{\partial \theta_1} \right) & \mathbb{V} \left(\frac{\partial \ell}{\partial \theta_2} \right) & \dots & \text{Cov} \left(\frac{\partial \ell}{\partial \theta_2}, \frac{\partial \ell}{\partial \theta_d} \right) \\ \vdots & \vdots & \ddots & \vdots \\ \text{Cov} \left(\frac{\partial \ell}{\partial \theta_d}, \frac{\partial \ell}{\partial \theta_1} \right) & \text{Cov} \left(\frac{\partial \ell}{\partial \theta_d}, \frac{\partial \ell}{\partial \theta_2} \right) & \dots & \mathbb{V} \left(\frac{\partial \ell}{\partial \theta_d} \right) \end{pmatrix}. \end{aligned} \quad \text{(by definition of covariance of a random vector)}$$

We have omitted the θ_0 arguments inside each of the variance and covariance operators in the final line above for the sake of avoiding overly cluttered notation, but it should be understood that all of these quantities are really integrals of the form

3.3. Maximum Likelihood Estimation

Maximum likelihood estimation is one of the most powerful and widespread techniques for obtaining point estimates of model parameters. The original intuition behind the method derives from the observation that when faced with a choice between two possible values of a parameter, the sensible choice is the one which makes the data we actually did observe more probable to have been observed. We have already defined the likelihood function as a means of capturing this probability, which makes expressing this decision rule in terms of it very easy - we simply choose for our estimate the option that produces the higher value of the likelihood function. That is, if $L(\theta_1; \mathbf{x}) > L(\theta_2; \mathbf{x})$, then under the preceding logic, θ_1 is the better estimate of the true parameter value.

This can be extended to include as many candidate parameter values as we would like. For n potential estimates of θ_0 , the best is the one that corresponds to the highest value of the likelihood function. If we follow this line of reasoning to its natural conclusion, we arrive at the notion of a *maximum likelihood estimate* of the parameter θ , which is defined as any value of θ that maximizes the output of the likelihood function among all possible choices of θ in Θ for fixed data \mathbf{x} . We will denote this using the notation $\hat{\theta}$ (pronounced “theta hat”). Formally, a maximum likelihood estimate of θ is defined as a function $\hat{\theta} = \hat{\theta}(\mathbf{x}) \in \Theta$ for which

$$(3.3.1) \quad L(\hat{\theta}; \mathbf{x}) = \sup_{\theta \in \Theta} L(\theta; \mathbf{x}),$$

or equivalently

$$(3.3.2) \quad \ell(\hat{\theta}; \mathbf{x}) = \sup_{\theta \in \Theta} \ell(\theta; \mathbf{x}).$$

In general, there is no guarantee that a maximum likelihood estimate for the parameter of a statistical model will exist, and if it does, it will not necessarily be unique.

3.4. Regularity Conditions

As a consequence of the random variable interpretation of likelihood-based quantities such as the score function and the MLE for θ_0 , a natural line of inquiry to investigate is how the behavior of these random variables changes as the sample size n increases. Of particular interest is the distribution to which the maximum likelihood estimate converges as n tends toward infinity, if indeed one exists. To that end, it will be useful to establish some *regularity conditions* for our models. We can think of these conditions as being assumptions similar to those we discussed in the introduction to this paper that, when satisfied, endow our models with certain properties that enable us to determine the aforementioned distribution.

For our purposes, we will call a model *regular* if it satisfies the following conditions:

- RC1) Any observations x_1, \dots, x_n belonging to a sample that has been drawn from the model's sample space are independent and identically distributed (i.i.d.) realizations of a random variable X with density function $p(x; \theta)$.
- RC2) For any two parameters $\theta_1, \theta_2 \in \Theta$, $\theta_1 \neq \theta_2 \implies p(x; \theta_1) \neq p(x; \theta_2)$.
- RC3) The parameter space Θ is a convex subset of \mathbb{R}^d .
- RC4) $\mathbb{E} \left([\mathcal{S}(\theta; x)] [\mathcal{S}(\theta; x)]^T; \theta \right)$ is positive definite for all $\theta \in \Theta$.
- RC5) There exists an open set $\Theta^* \subseteq \Theta$ of which θ_0 is an interior point.
- RC6) $p(x; \theta)$ is twice continuously differentiable with respect to θ for all $\theta \in \Theta^*$.
- RC7) The set $\{x \in \mathbb{R} : p(x; \theta) > 0\}$ does not depend on θ .
- RC8) The integral $\int_{-\infty}^{\infty} p(x; \theta) dx$ can be differentiated twice under the integral sign with respect to θ .
- RC9) For all $\theta \in \Theta^*$, there exists a radius $r > 0$ and a random function $M(\mathbf{x})$ which is allowed to depend on the data \mathbf{x} but not θ such that $\mathbb{E}[M(X); \theta_0] < \infty$ and $\sup_{\theta \in \Theta^*} |\nabla^2 \mathcal{S}(\theta; x)_{ijk}| \leq M(\mathbf{x})$ for all integers $1 \leq i, j, k \leq d$.

RC1) is the identifiability assumption of our model that we discussed in Section 1.1. We repeat it here is included merely to simplify calculations and frame our discussion in the context of a standard case in which likelihood theory holds. RC2-3) guarantee that it is always possible to solve the likelihood equation

for $\hat{\theta}$. RC4-6) ensure that there is no issue with defining derivatives of the log-likelihood function. Finally, by Leibniz's integral rule RC7) implies that , and that the remainder term in the first-order Taylor series expansion of the score function centered at θ_0 is negligible.

3.5. The Bartlett Identities

The Bartlett identities are a set of equations relating to the expectations of functions of derivatives of a log-likelihood function.

Consider the case in which a random variable X has a probability density f that depends on a scalar parameter θ . Denote the log-likelihood function for θ by $\ell(\theta; x) = \log f(x; \theta)$ and its first derivative with respect to θ by $\ell_\theta(\theta; x) = \frac{\partial}{\partial \theta} \ell(\theta; x)$. Taking the expectation of $\ell_\theta(\theta; x)$ gives

$$\begin{aligned}
 \mathbb{E}[\ell_\theta(\theta; x); \theta] &= \mathbb{E}\left[\frac{\partial}{\partial \theta} \ell(\theta; x); \theta\right] \\
 &= \int_{\mathbb{R}} \left[\frac{\partial}{\partial \theta} \ell(\theta; x)\right] f(x; \theta) dx \\
 &= \int_{\mathbb{R}} \left[\frac{\partial}{\partial \theta} \log f(x; \theta)\right] f(x; \theta) dx \\
 &= \int_{\mathbb{R}} \frac{\frac{\partial}{\partial \theta} f(x; \theta)}{f(x; \theta)} f(x; \theta) dx \\
 &= \int_{\mathbb{R}} \frac{\partial}{\partial \theta} f(x; \theta) dx \\
 &= \frac{d}{d\theta} \int_{\mathbb{R}} f(x; \theta) dx \quad (\text{by regularity of } f) \\
 &= \frac{d}{d\theta} 1 \\
 &= 0.
 \end{aligned}$$

Therefore,

$$(3.5.1) \quad \mathbb{E}[\ell_\theta(\theta; x); \theta] = 0 \text{ for all } \theta.$$

Equation 3.5.1 is called the first Bartlett identity. In words, it states that the expectation of the first derivative of the log-likelihood function of a statistical model with respect to the model parameter will

always be 0. Another name for ℓ_θ is the *score function*, and any pseudolikelihood that also satisfies the first Bartlett identity is said to be *score-unbiased*.

If we now consider the second derivative of $\ell(\theta; x)$, we have

$$\begin{aligned}
 \ell_{\theta\theta}(\theta; x) &= \frac{\partial^2}{\partial\theta^2} \ell(\theta; x) \\
 &= \frac{\partial}{\partial\theta} \left[\frac{\partial}{\partial\theta} \ell(\theta; x) \right] \\
 &= \frac{\partial}{\partial\theta} \left[\frac{\partial}{\partial\theta} \log f(x; \theta) \right] \\
 &= \frac{\partial}{\partial\theta} \left[\frac{\frac{\partial}{\partial\theta} f(x; \theta)}{f(x; \theta)} \right] \\
 &= \frac{\left[\frac{\partial^2}{\partial\theta^2} f(x; \theta) \right] f(x; \theta) - \left[\frac{\partial}{\partial\theta} f(x; \theta) \right] \left[\frac{\partial}{\partial\theta} f(x; \theta) \right]}{[f(x; \theta)]^2} \\
 &= \frac{\frac{\partial^2}{\partial\theta^2} f(x; \theta)}{f(x; \theta)} - \left[\frac{\frac{\partial}{\partial\theta} f(x; \theta)}{f(x; \theta)} \right]^2 \\
 &= \frac{\frac{\partial^2}{\partial\theta^2} f(x; \theta)}{f(x; \theta)} - \left[\frac{\partial}{\partial\theta} \log f(x; \theta) \right]^2 \\
 &= \frac{\frac{\partial^2}{\partial\theta^2} f(x; \theta)}{f(x; \theta)} - [\ell_\theta(\theta; x)]^2.
 \end{aligned}$$

Rearranging terms and taking expectations yields

$$\begin{aligned}
 \mathbb{E}[\ell_{\theta\theta}(\theta; x); \theta] + \mathbb{E}[(\ell_\theta(\theta; x))^2; \theta] &= \mathbb{E} \left[\frac{\frac{\partial^2}{\partial\theta^2} f(x; \theta)}{f(x; \theta)}; \theta \right] \\
 &= \int_{\mathbb{R}} \left[\frac{\frac{\partial^2}{\partial\theta^2} f(x; \theta)}{f(x; \theta)} \right] f(x; \theta) dx \\
 &= \int_{\mathbb{R}} \frac{\partial^2}{\partial\theta^2} f(x; \theta) dx \\
 &= \frac{d^2}{d\theta^2} \int_{\mathbb{R}} f(x; \theta) dx \quad (\text{by regularity of } f) \\
 &= \frac{d^2}{d\theta^2} 1 \\
 &= 0.
 \end{aligned}$$

Therefore,

$$(3.5.2) \quad \mathbb{E}[\ell_{\theta\theta}(\theta; x); \theta] + \mathbb{E}\left[(\ell_{\theta}(\theta; x))^2; \theta\right] = 0 \text{ for all } \theta.$$

Equation 3.5.2 is called the second Bartlett identity. The second term on the left-hand side can be further rewritten as

$$\begin{aligned} \mathbb{E}\left[(\ell_{\theta}(\theta; x))^2; \theta\right] &= \mathbb{V}[\ell_{\theta}(\theta; x); \theta] + \left(\mathbb{E}[\ell_{\theta}(\theta; x); \theta]\right)^2 \\ &= \mathbb{V}[\ell_{\theta}(\theta; x); \theta]. \end{aligned} \quad (\text{by the first Bartlett identity})$$

Another name for this quantity is the *expected information*. It follows from the second Bartlett identity that

$$(3.5.3) \quad \mathbb{E}[-\ell_{\theta\theta}(\theta; x); \theta] = \mathbb{V}[\ell_{\theta}(\theta; x); \theta].$$

The quantity $-\ell_{\theta\theta}(\theta; x)$ is called the *observed information*. Any pseudolikelihood that satisfies the second Bartlett identity is said to be *information-unbiased*.

It is possible to derive further Bartlett identities by continuing in this manner for an arbitrary number of derivatives of the log-likelihood function, provided that they exist. However, the first two are sufficient for our purposes of evaluating the validity of pseudolikelihoods as approximations to a genuine likelihood so we will not go further here. While the above derivations were performed under the assumption that θ is a scalar, the Bartlett identities also hold in the case where θ is a multi-dimensional vector.

3.6. One-Index Asymptotics

The one-index asymptotics framework describes the behavior of likelihood-based statistics as the sample size (n) grows to infinity while the dimension of the nuisance parameter (q) remains fixed. The aim of this section is to present a basic overview of the theory's results so that we will have a readily available baseline against which to compare the results of the following section discussing the two-index asymptotics framework, in which q is allowed to increase with n .

Assume the regularity conditions of the previous section apply to our model and let $\hat{\theta}$ denote the MLE for θ_0 . The traditional method for analyzing the asymptotic behavior of $\hat{\theta}$ is to use a first-order Taylor series expansion of the score function.¹

To do this, first note that the likelihood function for θ based on $\mathbf{x} = (x_1, \dots, x_n)$ as

$$L(\theta; \mathbf{x}) \equiv p(\mathbf{x}; \theta) = \prod_{i=1}^n p(x_i; \theta).$$

We may therefore write the log-likelihood function as

$$\begin{aligned} \ell(\theta; \mathbf{x}) &= \log L(\theta; \mathbf{x}) \\ &= \log p(\mathbf{x}; \theta) \\ &= \log \left[\prod_{i=1}^n p(x_i; \theta) \right] \\ &= \sum_{i=1}^n \log p(x_i; \theta) \\ &= \sum_{i=1}^n \ell(\theta; x_i). \end{aligned}$$

It follows that the score function can be written as

$$\begin{aligned} \mathcal{S}(\theta; \mathbf{x}) &= \frac{\partial}{\partial \theta} \ell(\theta; \mathbf{x}) \\ &= \frac{\partial}{\partial \theta} \sum_{i=1}^n \ell(\theta; x_i) \\ (3.6.1) \quad &= \sum_{i=1}^n \frac{\partial}{\partial \theta} \ell(\theta; x_i) \\ &= \sum_{i=1}^n \mathcal{S}(\theta; x_i). \end{aligned}$$

In other words, the score function for the parameter θ based on data x_1, \dots, x_n can be written as the sum of individual contributions $\mathcal{S}(\theta; x_i)$, $i = 1, \dots, n$, where each $\mathcal{S}(\theta; x_i)$ can be thought of as the score

¹See Appendix A for a review of Taylor's theorem and the conditions under which it is satisfied.

function for θ had we drawn only one observation, x_i . Note that these individual contributions are all independent from one another, a consequence of the independence assumption in RC1).

Equation 3.6.1 implies that a Taylor series expansion of $\mathcal{S}(\theta; \mathbf{x})$ will be equal to the sum of the Taylor series expansions of its individual contributions, plus a remainder term that grows with n . We have assumed each x_i is identically distributed, so it suffices to consider the expansion for an arbitrary contribution, $\mathcal{S}(\theta; x_i)$.

Since θ is a $d \times 1$ vector, the k -th derivative of the log-likelihood with respect to θ will be a k -dimensional array having d entries along each of its k indices. In particular, the score function (i.e. the first derivative of the log-likelihood) will also be a $d \times 1$ vector,

$$\mathcal{S}(\theta; x_i) = \begin{pmatrix} \mathcal{S}_1(\theta; x_i) \\ \vdots \\ \mathcal{S}_d(\theta; x_i) \end{pmatrix},$$

where each component is a function $\mathcal{S}_j : \Theta \rightarrow \mathbb{R}$. Similarly, the first and second derivatives of the score function will be a $d \times d$ matrix and a three-dimensional $d \times d \times d$ array, respectively. To simplify notation, we will perform a component-wise first-order Taylor series expansion of the score function around the point $\theta = \theta_0$ wherein each component $\mathcal{S}_j(\theta; x_i)$ of $\mathcal{S}(\theta; x_i)$ is expanded separately.

For any $\theta \in N_r(\theta_0)$, there exists $\bar{\theta}_j$ on the line segment connecting θ and θ_0 such that

$$\begin{aligned} \mathcal{S}_j(\theta; x_i) &= \mathcal{S}_j(\theta_0; x_i) + \nabla \mathcal{S}_j(\theta_0; x_i)^T (\theta - \theta_0) + \frac{1}{2} (\theta - \theta_0)^T \nabla^2 \mathcal{S}_j(\bar{\theta}_j; x_i) (\theta - \theta_0) \\ &= \mathcal{S}_j(\theta_0; x_i) + [\nabla \mathcal{S}_j(\theta_0; x_i) + \frac{1}{2} \nabla^2 \mathcal{S}_j(\bar{\theta}_j; x_i) (\theta - \theta_0)]^T (\theta - \theta_0) \\ &= \mathcal{S}_j(\theta_0; x_i) + [\nabla \mathcal{S}_j(\theta_0; x_i) + M(x_i) O(\|\theta - \theta_0\|)]^T (\theta - \theta_0), \end{aligned}$$

where the last equality follows as a result of RC7). Summing over each of the individual contributions to the score function yields

$$\begin{aligned}
\mathcal{S}_j(\theta; \mathbf{x}) &= \sum_{i=1}^n \mathcal{S}_j(\theta; x_i) \\
&= \sum_{i=1}^n [\mathcal{S}_j(\theta_0; x_i) + [\nabla \mathcal{S}_j(\theta_0; x_i) + M(x_i)O(\|\theta - \theta_0\|)]^T (\theta - \theta_0)] \\
&= \mathcal{S}_j(\theta_0; \mathbf{x}) + \left[\nabla \mathcal{S}_j(\theta_0; \mathbf{x}) + \left\{ \sum_{i=1}^n M(x_i) \right\} O(\|\theta - \theta_0\|) \right]^T (\theta - \theta_0) \\
&= \mathcal{S}_j(\theta_0; \mathbf{x}) + \left[\frac{1}{n} \nabla \mathcal{S}_j(\theta_0; \mathbf{x}) + \left\{ \frac{1}{n} \sum_{i=1}^n M(x_i) \right\} O(\|\theta - \theta_0\|) \right]^T n(\theta - \theta_0).
\end{aligned}$$

If we divide through by \sqrt{n} , we arrive at

$$\frac{1}{\sqrt{n}} \mathcal{S}_j(\theta; \mathbf{x}) = \frac{1}{\sqrt{n}} \mathcal{S}_j(\theta_0; \mathbf{x}) + \left[\frac{1}{n} \nabla \mathcal{S}_j(\theta_0; \mathbf{x}) + \left\{ \frac{1}{n} \sum_{i=1}^n M(x_i) \right\} O(\|\theta - \theta_0\|) \right]^T \sqrt{n}(\theta - \theta_0).$$

CHAPTER 4

Pseudolikelihood Functions

4.1. Model Parameter Decomposition

It is often the case that we are not interested in estimating the full parameter $\theta \in \Theta \subseteq \mathbb{R}^d$, but rather a different parameter ψ taking values in a set $\Psi \subseteq \mathbb{R}^p$, where $p < d$. In such an event, we refer to ψ as the *parameter of interest*.

Since ψ is of lower dimension than θ , it necessarily follows that there is another parameter λ , taking values in a set $\Lambda \subseteq \mathbb{R}^q$, where $p + q = d$, that is made up of whatever is “left over” from the full parameter θ . We refer to λ as the *nuisance parameter*, so named for its ability to complicate inference regarding the parameter of interest. Despite not being the object of study themselves, nuisance parameters are nevertheless capable of modifying the distributions of our observations and therefore must be accounted for when conducting inference or estimation regarding the parameter of interest.¹ The process by which this is accomplished is often nontrivial and can constitute a significant barrier that must be overcome.

While not required, we will assume the parameter of interest ψ is always one-dimensional. That is, $\Psi \subseteq \mathbb{R}$ and consequently $\Lambda \subseteq \mathbb{R}^{d-1}$. This restriction reflects the common habit of researchers to focus on scalar-valued summaries of vector quantities. For example, suppose we observe data $Y = (y_1, \dots, y_n)$, where each y_i is the outcome of some random variable $Y_i \sim N(\mu_i, \sigma_i^2)$, and we are interested in estimating the average of the population means, $\frac{1}{n} \sum_{i=1}^n \mu_i$. Rather than defining $\psi = (\mu_1, \dots, \mu_n)$, we can instead define $\psi = \frac{1}{n} \sum_{i=1}^n \mu_i$ directly, bypassing the need to estimate each μ_i individually before taking their average. This does carry the trade-off of increasing the dimension of the nuisance parameter, which must be dealt with before conducting inference or estimation on ψ . We will examine some of the issues posed by high-dimensional nuisance parameters in greater detail in the next chapter.

¹Note that nuisance parameters are not always uniquely defined. Depending on the choice of parameter of interest, there may be multiple or even infinite ways to define a nuisance parameter.

4.1.1. Explicit Parameters

Parameters of interest and nuisance parameters can be broadly classified into two categories, explicit or implicit. For a given statistical model, both types of parameter must occupy the same category - it is not possible for ψ to be explicit and λ to be implicit, or vice versa.

Let us first consider the case in which ψ and λ are *explicit* parameters. This means that ψ is a sub-vector of θ , so that all the components of ψ are also components of θ . Then there exists a set $I = \{I_1, \dots, I_p\} \subsetneq \{1, \dots, d\}$ such that

$$(4.1.1) \quad \psi = (\theta_{I_1}, \dots, \theta_{I_p}).$$

It immediately follows that λ is the sub-vector of all components of θ that are not part of ψ . More precisely, if we let $J = \{J_1, \dots, J_q\} \subsetneq \{1, \dots, q\}$ such that $I \cup J = \{1, \dots, d\}$ and $I \cap J = \emptyset$, then

$$(4.1.2) \quad \lambda = (\theta_{J_1}, \dots, \theta_{J_q}).$$

θ can therefore be decomposed as $\theta = (\psi, \lambda)$ when ψ and λ are explicit, provided we shuffle the indices appropriately.

4.1.2. Implicit Parameters

Now let us consider the case in which ψ and λ are *implicit* parameters. This means there exists some function $\varphi : \Theta \rightarrow \Psi$ for which the parameter of interest can be written as

$$(4.1.3) \quad \psi = \varphi(\theta).$$

As before, Ψ is still assumed to be a subset of \mathbb{R}^p where p is less than d . This reduction in dimension again implies the existence of a nuisance parameter $\lambda \in \Lambda \subseteq \mathbb{R}^{k-m}$. However, unlike in the explicit case, a closed form expression for λ in terms of the original components of θ need not exist. For this reason, implicit nuisance parameters are in general more difficult to eliminate compared to their explicit counterparts.

Note that when the parameter of interest and nuisance parameter are explicit, it is always possible to define a function φ such that

$$(4.1.4) \quad \varphi(\theta) = (\theta_{I_1}, \dots, \theta_{I_p}) \equiv \psi,$$

where $\{I_1, \dots, I_p\}$ is defined as above. Hence, the first case is really just a special example of this more general one in which $\psi = \varphi(\theta)$. With this understanding in mind, we will use the notation $\psi = \varphi(\theta)$ to refer to the parameter of interest in general, only making the distinction between implicitness and explicitness when the difference is relevant to the situation.

4.2. Types of Pseudolikelihoods

The natural solution to the hindrance nuisance parameters pose to making inferences on the parameter of interest is to find a method for eliminating them from the model altogether. Since one way of uniquely specifying a model is through its likelihood function, this is equivalent to eliminating the nuisance parameters from the likelihood function itself. The result of this elimination is what is known as a pseudolikelihood function.

In general, a *pseudolikelihood function* for ψ is a function of the data and ψ only, having properties resembling that of a genuine likelihood function. Suppose $\psi = \varphi(\theta)$ for some function φ and parameter $\theta \in \Theta$. If we let $\Theta(\psi) = \{\theta \in \Theta : \varphi(\theta) = \psi\}$, then associated with each $\psi \in \Psi$ is the set of likelihoods $\mathcal{L}_\psi = \{L(\theta) : \theta \in \Theta(\psi)\}$.

Any summary of the values in \mathcal{L}_ψ that does not depend on λ theoretically constitutes a pseudolikelihood function for ψ . There exist a variety of methods to obtain this summary but among the most popular are profiling (maximization), conditioning, and integration, each with respect to the nuisance parameter. We will explore each of these methods in more detail in the following sections.

4.2.1. The Profile Likelihood

The profile likelihood is the most straightforward method for eliminating a nuisance parameter from a likelihood function.

For example, suppose we are interested in estimating the mean of a random variable Y , where $Y \sim N(\mu, \sigma^2)$. The full model parameter is $\theta = (\mu, \sigma^2)$ but since we are only interested in estimating the mean, the parameter of interest is $\psi = \mu$ and the nuisance parameter is $\lambda = \sigma^2$.

4.2.2. The Conditional Likelihood

4.2.3. The Marginal Likelihood

4.2.4. The Integrated Likelihood

4.3. The Bartlett Identities Revisited

The Bartlett identities offer an alternative way of characterizing the difference between likelihood and pseudolikelihood functions. A genuine likelihood function of θ can be characterized as any nonnegative random function of θ for which all of the Bartlett identities hold. Similarly, we can think of a pseudolikelihood function of θ as being any nonnegative random function of θ for which at least one of the Bartlett identities does not hold. Hence, the identities act as a litmus test of sorts for determining the validity of a pseudolikelihood as an approximation to the genuine likelihood from which it originated - the more identities it does satisfy, the better the approximation.

CHAPTER 5

Approximating the Integrated Likelihood Function

5.1. The Zero-Score Expectation Parameter

Let $\psi = \varphi(\theta)$ and λ denote the parameter of interest and nuisance parameter, respectively, for some statistical model $(\mathcal{S}, \mathcal{P}_\theta)$. Then the general expression to obtain an integrated likelihood for ψ may be written as

$$(5.1.1) \quad \bar{L}(\psi) = \int_{\Lambda} L(\psi, \lambda) \pi(\lambda|\psi) d\lambda,$$

where $\pi(\lambda|\psi)$ is a conditional prior density for λ given ψ .

Severini (2007) considered the problem of selecting $\pi(\lambda|\psi)$ such that when the likelihood function is integrated with respect to this density, the result is useful for non-Bayesian inference. To do this, he outlined four properties (see Appendix B) that an integrated likelihood function must satisfy if it is to be of any use. He went on to prove that an integrated likelihood satisfying these properties could be obtained by first constructing a new nuisance parameter $\phi \in \Phi$ that is unrelated to the parameter of interest (in the sense that its maximum likelihood estimator remains roughly constant for all values of ψ) and then choosing a prior density $\pi(\phi)$ that is independent of ψ . Once chosen, the desired integrated likelihood function for ψ is given by

$$(5.1.2) \quad \bar{L}(\psi) = \int_{\Phi} \tilde{L}(\psi, \phi) \pi(\phi) d\phi,$$

where $\tilde{L}(\psi, \phi)$ is the likelihood function for the model after it has been reparameterized in terms of ϕ . It is important to note that the exact choice of prior density for ϕ is not particularly important; the only restriction we place upon it is that it must not depend on ψ .

Suppose that we have an explicit parameter of interest and nuisance parameter, so that $\theta = (\psi, \lambda)$. Then Severini (2007) defines this new nuisance parameter ϕ as the solution to the equation

$$(5.1.3) \quad \mathbb{E}(\ell_\lambda(\psi, \lambda); \psi_0, \lambda_0) \Big|_{(\psi_0, \lambda_0) = (\hat{\psi}, \phi)} = 0,$$

where $\ell_\lambda(\psi, \lambda) = \frac{\partial \ell(\psi, \lambda)}{\partial \lambda}$, ψ_0 and λ_0 denote the true values of ψ and λ , and $\hat{\psi}$ is the MLE for ψ_0 . In other words, for a particular value of $(\psi, \lambda, \hat{\psi})$, we can find the corresponding value of ϕ by solving for it in Equation 5.1.3. ϕ is called the *zero-score expectation* (ZSE) parameter because it is defined as the value that makes the expectation of the score function (where the derivative is taken with respect to λ) evaluated at the point $(\hat{\psi}, \phi)$ equal to zero.

Note that ϕ is a function of the data through $\hat{\psi}$. Normally we avoid creating such dependencies in our parameters as it renders them useless for the purpose of parameterizing a statistical model. However, from the perspective of the likelihood function, once the data have been collected they are considered fixed in place and there is no issue with using a quantity such as ϕ that depends on the data to parameterize it.

For a given value of $(\psi, \phi, \hat{\psi})$, it is also possible to solve Equation 5.1.3 for λ . This allows us to write Equation 5.1.2 in terms of $L(\psi, \lambda)$:

$$(5.1.4) \quad \bar{L}(\psi) = \int_{\Phi} L(\psi, \lambda(\psi, \phi)) \pi(\phi) d\phi.$$

Severini (2018) proved that reparameterizing the nuisance parameter in terms of the ZSE parameter yields the same desirable properties in the subsequent integrated likelihood when ψ and λ are implicit. Suppose $\psi = \varphi(\theta)$, for some function $\varphi : \Theta \rightarrow \Psi$, and consider the set of all values of θ satisfying $\varphi(\theta) = \hat{\psi}$. Call this set $\Omega_{\hat{\psi}}$ so that

$$(5.1.5) \quad \Omega_{\hat{\psi}} = \{\omega \in \Theta : \varphi(\omega) = \hat{\psi}\}.$$

Elements of $\Omega_{\hat{\psi}}$ take the form $(\hat{\psi}, \phi)$, where $\phi \in \Lambda$.

5.1.1. Weight Functions

5.2. Two-Index Asymptotics

Earlier we discussed the one-index asymptotics setting, in which the sample size (n) of the model diverged to infinity while the dimension of the nuisance parameter (q) remained fixed. Now we turn our attention to the two-index asymptotics setting which describes the behavior of likelihood and pseudolikelihood functions as n and q both tend to infinity, with q growing at least as fast as n . Under such a framework, De Bin, Sartori, and Severini (2015) showed that estimates for ψ based on a suitably constructed integrated likelihood function will outperform those coming from more traditional pseudolikelihoods, such as the profile likelihood. Such findings provide the motivation for our ensuing examination of two-index asymptotics theory, insofar as it relates to the performance of the integrated likelihood function as a method of inference regarding a parameter of interest.

To mirror the strategy used by Sartori (2003) and De Bin, Sartori, and Severini (2015), we will frame our discussion in terms of a stratified sample of data in which each stratum contributes one component to the overall nuisance parameter. Consider a model with parameter $\theta = (\psi, \lambda)$ where ψ is the parameter of interest and $\lambda = (\lambda_1, \dots, \lambda_q)$ is a q -dimensional nuisance parameter. For the sake of reducing complexity in our notation, we will only consider the case in which ψ and the individual components of λ are scalar parameters, though the results of this section should hold in the case where they are all vectors as well. Suppose that we have divided the model's population into q strata and collected a sample of size m_i from each stratum such that observation j from stratum i may be modeled as

$$(5.2.1) \quad X_{ij} \sim p_{ij}(x_{ij}; \psi, \lambda_i),$$

where $i = 1, \dots, q$ and $j = 1, \dots, m_i$, making the total sample size $n = \sum_{i=1}^q m_i$.¹ Hence, there is a one-to-one correspondence between the strata and the components of λ ; assume that each $\lambda_i \in \Lambda$, where the space Λ is the same for all i , and they all have the same interpretation within their respective strata.

¹It will be convenient to work under the restriction that the stratum sample sizes are all identical, meaning there exists some positive integer m such that $m_i = m$ for all i . However, as both Sartori (2003) and De Bin, Sartori, and Severini (2015) note, we could also assume a looser condition in which $m_i = K_i m$ where $0 < K_i < \infty$ without compromising our results.

Assume that all the regularity conditions set forth in Section 3.4 apply, except possibly for RC1) - it is not necessary to assume that the observations are i.i.d. here, and in fact it is perfectly acceptable for the p_{ij} 's in Equation 5.2.1 to differ from one another. We will also allow for the possibility of dependence among observations within a stratum, though not between them.

Let $\mathbf{x}_i = (x_{i1}, \dots, x_{im})$ denote the sample of observations from stratum i , so that their joint density may be written as $p_i(\mathbf{x}_i; \psi, \lambda_i)$. Therefore, the likelihood and log-likelihood for the i th stratum are

$$(5.2.2) \quad L^{(i)}(\psi, \lambda_i) = p_i(\mathbf{x}_i; \psi, \lambda_i),$$

and

$$(5.2.3) \quad \ell^{(i)}(\psi, \lambda_i) = \log L^{(i)}(\psi, \lambda_i),$$

respectively. For a particular choice of weight function $g(\lambda_i; \psi)$, the integrated likelihood for ψ in stratum i is given by

$$(5.2.4) \quad \bar{L}^{(i)}(\psi) = \int_{\Lambda} L^{(i)}(\psi, \lambda_i) g(\lambda_i; \psi) d\lambda_i.$$

From here, we proceed by using Laplace's method as described by Tierney and Kadane (1986) (see Appendix B for a brief review) to obtain an analytic approximation to $\bar{L}^{(i)}(\psi)$. Setting $h(\lambda_i) = -\frac{1}{m}\ell^{(i)}(\psi, \lambda_i)$ and $f(\lambda_i) = g(\lambda_i; \psi)$, we may rewrite the integral in Equation 5.2.4 as

$$\bar{L}^{(i)}(\psi) = \int_{\Lambda} f(\lambda_i) \exp[-mh(\lambda_i)] d\lambda_i.$$

One consequence of the regularity conditions we have assumed is that $L^{(i)}(\psi, \lambda_i)$ is that an MLE for θ_0 , $\hat{\theta}$, exists and is unique. This further implies the existence and uniqueness of a conditional MLE for the true value of each stratum-specific nuisance parameter given ψ - denote this value for the i th stratum by $\hat{\lambda}_{i\psi}$. By definition this value maximizes $\ell^{(i)}(\psi, \lambda_i)$ as a function of λ_i , and so it also maximizes $-h(\lambda_i)$ since the two functions differ only by a multiplicative constant $\frac{1}{m}$.

The Laplace approximation to $\bar{L}^{(i)}(\psi)$ is then given by

$$(5.2.5) \quad \hat{\hat{L}}^{(i)}(\psi) = f(\hat{\lambda}_{i\psi}) \sqrt{\frac{2\pi}{m}} \sigma \exp[-mh(\hat{\lambda}_{i\psi})],$$

where

$$\begin{aligned} \sigma &= \left[\frac{\partial^2 h}{\partial \lambda_i^2} \Big|_{\lambda_i = \hat{\lambda}_{i\psi}} \right]^{-1/2} \\ &= \left[-\frac{1}{m} \frac{\partial^2 \ell^{(i)}(\psi, \lambda_i)}{\partial \lambda_i^2} \Big|_{\lambda_i = \hat{\lambda}_{i\psi}} \right]^{-1/2} \\ &= \left[\frac{1}{m} \mathcal{J}(\hat{\lambda}_{i\psi}) \right]^{-1/2}. \end{aligned}$$

Here, $\mathcal{J}(\hat{\lambda}_{i\psi})$ denotes the observed information function for λ_i only (i.e. the negative second partial derivative of the log-likelihood with respect to λ_i) evaluated at $\hat{\lambda}_{i\psi}$. Plugging in the appropriate quantities for f , h , and σ into Equation 5.2.5, we arrive at

$$\begin{aligned} (5.2.6) \quad \hat{\hat{L}}^{(i)}(\psi) &= f(\hat{\lambda}_{i\psi}) \sqrt{\frac{2\pi}{m}} \sigma \exp[-mh(\hat{\lambda}_{i\psi})] \\ &= g(\hat{\lambda}_{i\psi}; \psi) \sqrt{\frac{2\pi}{m}} \left[\frac{1}{m} \mathcal{J}(\hat{\lambda}_{i\psi}) \right]^{-1/2} \exp \left\{ -m \cdot -\frac{1}{m} \ell^{(i)}(\psi, \hat{\lambda}_{i\psi}) \right\} \\ &= \frac{\sqrt{2\pi}}{m} L^{(i)}(\psi, \hat{\lambda}_{i\psi}) g(\hat{\lambda}_{i\psi}; \psi) [\mathcal{J}(\hat{\lambda}_{i\psi})]^{-1/2} \\ &= \frac{\sqrt{2\pi}}{m} L_P^{(i)}(\psi) g(\hat{\lambda}_{i\psi}; \psi) [\mathcal{J}(\hat{\lambda}_{i\psi})]^{-1/2}, \end{aligned}$$

where $L_P^{(i)}(\psi) = L^{(i)}(\psi, \hat{\lambda}_{i\psi})$ is the profile likelihood for ψ . The error in this approximation is

$$(5.2.7) \quad \bar{L}^{(i)}(\psi) = \hat{\hat{L}}^{(i)}(\psi) \left\{ 1 + O\left(\frac{1}{m}\right) \right\}.$$

Let

$$(5.2.8) \quad \bar{\ell}^{(i)}(\psi) = \log \bar{L}^{(i)}(\psi)$$

denote the integrated log-likelihood for ψ . Putting the results in Equation 5.2.6, Equation 5.2.7, and Equation 5.2.8 together, we have

$$\begin{aligned}
\bar{\ell}^{(i)}(\psi) &= \log \bar{L}^{(i)}(\psi) && \text{(by Equation 5.2.8)} \\
&= \log \left(\hat{\bar{L}}^{(i)}(\psi) \left\{ 1 + O\left(\frac{1}{m}\right) \right\} \right) && \text{(by Equation 5.2.7)} \\
&= \log \hat{\bar{L}}^{(i)}(\psi) + \log \left\{ 1 + O\left(\frac{1}{m}\right) \right\} && \text{(by Equation 5.2.6)} \\
&= \log \left\{ \frac{\sqrt{2\pi}}{m} L_P^{(i)}(\psi) g(\hat{\lambda}_{i\psi}; \psi) [\mathcal{J}(\hat{\lambda}_{i\psi})]^{-1/2} \right\} + O\left(\frac{1}{m}\right) \\
&= \log L_P^{(i)}(\psi) + \log g(\hat{\lambda}_{i\psi}; \psi) - \frac{1}{2} \log \mathcal{J}(\hat{\lambda}_{i\psi}) + \frac{1}{2} \log(2\pi) - \log(m) + O\left(\frac{1}{m}\right) \\
&= \ell_P^{(i)}(\psi) + \log g(\hat{\lambda}_{i\psi}; \psi) - \frac{1}{2} \log \mathcal{J}(\hat{\lambda}_{i\psi}) + \frac{1}{2} \log(2\pi) - \log(m) + O\left(\frac{1}{m}\right),
\end{aligned}$$

where $\ell_P^{(i)}(\psi) = \ell^{(i)}(\psi, \hat{\lambda}_{i\psi})$ is the profile log-likelihood for ψ . Since log-likelihoods are equivalent up to additive constants, we can discard the $\frac{1}{2} \log(2\pi)$ term in the final line above to arrive at our final approximation for the integrated log-likelihood in stratum i :

$$(5.2.9) \quad \hat{\bar{\ell}}^{(i)}(\psi) = \ell_P^{(i)}(\psi) + \log g(\hat{\lambda}_{i\psi}; \psi) - \frac{1}{2} \log \mathcal{J}(\hat{\lambda}_{i\psi}).$$

The error in this approximation is given by

$$(5.2.10) \quad \bar{\ell}^{(i)}(\psi) = \hat{\bar{\ell}}^{(i)}(\psi) - \log(m) + O\left(\frac{1}{m}\right).$$

Since the observations between the strata are independent, we may write the likelihood and log-likelihood functions for the entire model as

$$(5.2.11) \quad L(\psi, \lambda) = \prod_{i=1}^q L^{(i)}(\psi, \lambda_i)$$

and

$$(5.2.12) \quad \ell(\psi, \lambda) = \sum_{i=1}^q \ell^{(i)}(\psi, \lambda_i),$$

respectively. If we define the weight function

$$(5.2.13) \quad G(\lambda; \psi) \equiv \prod_{i=1}^q g(\lambda_i; \psi)$$

then the integrated likelihood function for ψ becomes separable. That is,

$$(5.2.14) \quad \begin{aligned} \bar{L}(\psi) &= \int_{\Lambda^q} L(\psi, \lambda) G(\lambda; \psi) d\lambda \\ &= \int_{\Lambda} \cdots \int_{\Lambda} \left[\prod_{i=1}^q L^{(i)}(\psi, \lambda_i) \right] \left[\prod_{i=1}^q g(\lambda_i; \psi) \right] d\lambda_1 \cdots d\lambda_q \\ &= \prod_{i=1}^q \int_{\Lambda} L^{(i)}(\psi, \lambda_i) g(\lambda_i; \psi) d\lambda_i \\ &= \prod_{i=1}^q \bar{L}^{(i)}(\psi). \end{aligned}$$

Let $\bar{\ell}(\psi) = \log \bar{L}(\psi)$ denote the integrated log-likelihood function for ψ . Taking the logarithm of both sides in Equation 5.2.14, we have

$$(5.2.15) \quad \bar{\ell}(\psi) = \sum_{i=1}^q \bar{\ell}^{(i)}(\psi).$$

Plugging in our approximation to $\bar{\ell}^{(i)}(\psi)$ and its error term in Equation 5.2.9 and Equation 5.2.10, respectively, yields

$$(5.2.16) \quad \begin{aligned} \bar{\ell}(\psi) &= \sum_{i=1}^q \left[\ell_P^{(i)}(\psi) + \log g(\hat{\lambda}_{i\psi}; \psi) - \log \mathcal{J}(\hat{\lambda}_{i\psi}) - \log(m) + O\left(\frac{1}{m}\right) \right] \\ &= \ell_P(\psi) + \sum_{i=1}^q \log g(\hat{\lambda}_{i\psi}; \psi) - \frac{1}{2} \sum_{i=1}^q \log \mathcal{J}(\hat{\lambda}_{i\psi}) - q \log(m) + O\left(\frac{q}{m}\right). \end{aligned}$$

CHAPTER 6

Applications**6.1. Multinomial Distribution****6.2. Standardized Mean Difference**

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APPENDIX A

Chapter 3

A.1. Definitions and Notation

A.1.1. Neighborhoods

A *neighborhood* of a point $\mathbf{p} \in \mathbb{R}^d$ is the set of all points $\mathbf{x} \in \mathbb{R}^d$ such that the Euclidean distance between \mathbf{p} and \mathbf{x} is less than some radius $r > 0$. We use the following notation to refer to such neighborhoods:

$$N_r(\mathbf{p}) = \{\mathbf{x} \in \mathbb{R}^d : \|\mathbf{x} - \mathbf{p}\| < r\}.$$

A.1.2. Line Segments

For two points $\mathbf{x}_1, \mathbf{x}_2 \in \mathbb{R}^d$, a third point $\bar{\mathbf{x}}$ is said to be on the *line segment* connecting \mathbf{x}_1 and \mathbf{x}_2 if there exists $\omega \in [0, 1]$ such that $\bar{\mathbf{x}} = \omega\mathbf{x}_1 + (1 - \omega)\mathbf{x}_2$. We use the following notation to refer to such line segments:

$$LS(\mathbf{x}_1, \mathbf{x}_2) = \{\omega\mathbf{x}_1 + (1 - \omega)\mathbf{x}_2 : \omega \in [0, 1]\}.$$

A.1.3. Higher Order Partial Derivatives

For $\boldsymbol{\alpha} \in \mathbb{N}^d$ and $\mathbf{x} \in \mathbb{R}^d$, we define the following multi-index notation:

$$|\boldsymbol{\alpha}| = \alpha_1 + \dots + \alpha_d, \quad \boldsymbol{\alpha}! = \alpha_1! \dots \alpha_d!, \quad \mathbf{x}^{\boldsymbol{\alpha}} = x_1^{\alpha_1} \dots x_d^{\alpha_d}.$$

For a function $f : \mathbb{R}^d \rightarrow \mathbb{R}$ that is k times differentiable at the point $\mathbf{a} \in \mathbb{R}^d$, define the following notation to indicate the mixed partial derivatives of f at \mathbf{a} up to and including order k :

$$D^{\boldsymbol{\alpha}} f = \frac{\partial^{|\boldsymbol{\alpha}|} f}{\partial x_1^{\alpha_1} \dots \partial x_d^{\alpha_d}}, \quad |\boldsymbol{\alpha}| \leq k.$$

A.2. Taylor's Theorem

Let $f : \mathbb{R}^d \rightarrow \mathbb{R}$ be a function that is $(k+1)$ -times continuously differentiable in a neighborhood $N_r(\mathbf{x}_0)$ of some point $\mathbf{x}_0 \in \mathbb{R}^d$ and suppose there exists M satisfying $|D^\alpha f| \leq M$ for all $x \in N_r(x_0)$ and all α such that $|\alpha| = k+1$. Then

$$f(\mathbf{x}) = \sum_{0 \leq |\alpha| \leq k} \frac{D^\alpha f(\mathbf{x}_0)}{\alpha!} (\mathbf{x} - \mathbf{x}_0)^\alpha + R_k(\mathbf{x}),$$

where the remainder term $R_k(\mathbf{x})$ satisfies

$$|R_k(\mathbf{x})| \leq \frac{M}{\alpha!} |\mathbf{x} - \mathbf{x}_0|^\alpha$$

for all α such that $|\alpha| = k+1$

APPENDIX B

Chapter 5

B.1. Desirable Properties of the Integrated Likelihood**B.1.1. Property 1**

Suppose the likelihood function for a parameter θ can be decomposed as the product $L(\theta) = L_1(\psi)L_2(\lambda)$. Then the integrated likelihood for ψ should satisfy

$$\bar{L}(\psi) = L_1(\psi).$$

B.1.2. Property 2**B.1.3. Property 3****B.1.4. Property 4****B.2. Laplace's Method**

Let θ be a scalar parameter taking values in \mathbb{R} and consider an integral of the form

$$I = \int_{-\infty}^{\infty} f(\theta) \exp[-nh(\theta)] d\theta.$$

Suppose the function $-h(\theta)$ is smooth, bounded, and unimodal so that it attains a maximum at a point $\hat{\theta}$. Then Laplace's method states that an approximation for I is given by

$$\hat{I} = f(\hat{\theta}) \sqrt{\frac{2\pi}{n}} \sigma \exp[-nh(\hat{\theta})],$$

where

$$\sigma = \left[\frac{\partial^2 h}{\partial \theta^2} \bigg|_{\theta=\hat{\theta}} \right]^{-1/2}.$$

It can further be shown that

$$I = \hat{I} \left\{ 1 + O\left(\frac{1}{n}\right) \right\},$$

where the term n may be interpreted as the sample size.