Optimisation

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STAT 3150-Statistical Computing

Lecture Objectives

- Give examples of optimisation problems in statistics.
- Apply the Newton-Raphson algorithm to compute Maximum Likelihood estimates.

Motivation

- Optimisation is a very old and broad field of applied mathematics.
 - It is older than calculus: see Heron's problem.
- It plays an important role in statistics, but also computational biology, economics, mathematical finance, etc.
- Our discussion will necessarily be brief, but I will try to give a sample of the types of statistical problems that can be solved using optimisation.
 - For more details, I recommend *Optimization* (2013) by Kenneth Lange.

Problem statement

- In very general terms, the basic optimisation problem is as follows.
- Let A be a set (e.g. a subset of \mathbb{R}^n , a set of integers, etc.), and let f be a real-valued function on A.
- We are looking for a value $x_0 \in A$ such that:
 - Minimisation: $f(x_0) \le f(x)$ for all $x \in A$.
 - Maximisation: $f(x_0) \ge f(x)$ for all $x \in A$.
- The set A can be defined via a combination of equality and inequality constraints.
 - For a multinomial model, we want K probabilities p_k such that $0 \le p_k \le 1$ and $\sum_{k=1}^K p_k = 1$.

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Lasso regression

- In linear regression, we find an estimate $\hat{\beta}$ by minimising least squares:

$$\hat{\beta} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^p} \sum_{i=1}^n \left(Y_i - \mathbf{X}_i^T \beta \right)^2.$$

• In lasso regression, we add a constraint on the L1 norm of β :

$$\hat{\beta}_{Lasso} = \operatorname*{arg\,min}_{\beta \in \mathbb{R}^p} \sum_{i=1}^n \left(Y_i - \mathbf{X}_i^T \beta \right)^2, \quad \sum_{j=1}^p |\beta_j| \le \lambda.$$

• Why? By adding the L1 constraint, some of the components of the solution $\hat{\beta}_{Lasso}$ can be exactly zero, leading to *variable* selection.

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Dimension reduction

- Let \mathbf{Y} be a p-dimensional random vector.
- In dimension reduction, we are looking for a linear combination $w^T\mathbf{Y}$ that optimises a given criterion.
- · For example:
 - In Principal Component Analysis, we are looking to maximise ${\rm Var}\left(w^T{\bf Y}\right)$.
 - In Independent Component Analysis, we are looking to minimise mutual information.

Mixture models

· Recall from earlier this semester mixture models:

$$F(x) = \sum_{k=1}^{K} \pi_k F_k(x),$$

where F_k is a distribution function for all k and $\sum_{k=1}^K \pi_k = 1$.

- The likelihood can be difficult to maximise directly.
- Instead, we typically use the **Expectation-Maximization** algorithm.

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Machine learning

- A common goal in machine learning is prediction: given a vector of features ${\bf X}$, find a function f such that

$$L(Y, f(\mathbf{X}))$$

is minimised.

• Here $L(Y, \cdot)$ is a loss function. For example:

$$L(Y, f(\mathbf{X})) = (Y - f(\mathbf{X}))^2$$

$$L(Y, f(\mathbf{X})) = |Y - f(\mathbf{X})|$$

Newton-Raphson method i

- The Newton-Raphson method is the only optimisation technique we will discuss.
 - It is a very important technique and serves as motivation for more advanced methods.
- · Suppose we have a twice differentiable function $f:A\to\mathbb{R}$, where $A\subseteq\mathbb{R}^n$. We want to find the maximum of f on A.
- $\boldsymbol{\cdot}$ We can approximate f using a Taylor expansion:

$$f(\mathbf{x} + \mathbf{h}) \approx f(\mathbf{x}) + \nabla f(\mathbf{x})^T \mathbf{h} + \frac{1}{2} \mathbf{h}^T D^2 f(\mathbf{x}) \mathbf{h},$$

where $\nabla f(\mathbf{x})$ is the vector of first-order derivatives and $D^2f(\mathbf{x})$ is the matrix of second-order derivatives.

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Newton-Raphson method ii

- · They are called the gradient and the Hessian, respectively.
- In other words, we are approximating f in a neighbourhood of \mathbf{x} using a quadratic function of \mathbf{h} .
- We know how to find the maximum of a quadratic function: take the derivative, set it equal to zero, and solve for h.
 - Why? The derivative of a quadratic is a linear function. Easy to solve!
- \cdot The derivative of the Taylor approximation with respect to ${f h}$ is

$$\frac{d}{d\mathbf{h}}\left[f(\mathbf{x}) + \nabla f(\mathbf{x})^T\mathbf{h} + \frac{1}{2}\mathbf{h}^TD^2f(\mathbf{x})\mathbf{h}\right] = \nabla f(\mathbf{x}) + D^2f(\mathbf{x})\mathbf{h}.$$

Newton-Raphson method iii

- Setting the derivative equal to zero and solving for \boldsymbol{h} gives

$$\mathbf{h} = -[D^2 f(\mathbf{x})]^{-1} \nabla f(\mathbf{x}).$$

- Therefore, the approximate maximum of our function f is a neighbourhood of ${\bf x}$ occurs at

$$\mathbf{x} + \mathbf{h} = \mathbf{x} - [D^2 f(\mathbf{x})]^{-1} \nabla f(\mathbf{x}).$$

· In one-dimension, this simplifies to

$$x + h = x - \frac{f'(x)}{f''(x)}.$$

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Newton-Raphson method iv

Algorithm

Let \mathbf{x}_0 be an initial value. Then construct a sequence via

$$\mathbf{x}_{n+1} = \mathbf{x}_n - [D^2 f(\mathbf{x}_n)]^{-1} \nabla f(\mathbf{x}_n).$$

If the initial value \mathbf{x}_0 is "sufficiently close" to a maximum (or minimum) of f, then the sequence \mathbf{x}_n will converge to that maximum (or minimum).

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Example 1 i

- Consider the function $f(x) = -x^2 + \sin(x)$.
- · The derivatives are

$$f'(x) = -2x + \cos(x), \quad f''(x) = -2 - \sin(x).$$

Example 1 ii

```
# Create functions
gradient <- function(x) -2*x + cos(x)
hessian <- function(x) -2 - sin(x)
# Set-up parameters
x current <- 1
x next <- x current -
  gradient(x_current)/hessian(x current)
tol <- 10<sup>^</sup>-10
iter <-1
```

Example 1 iii

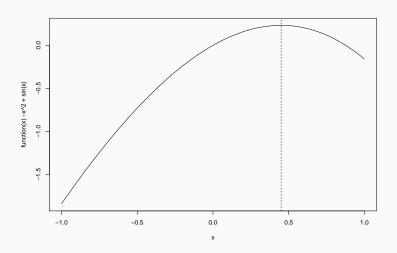
```
while (abs(x_current - x_next) > tol &
       iter < 100) {
  iter <- iter + 1
    x_current <- x_next
    step <- gradient(x_current)/hessian(x_current)</pre>
    x_next <- x_current - step</pre>
x next
```

[1] 0.4501836

Example 1 iv

```
plot(function(x) -x^2 + \sin(x), from = -1, to = 1)
abline(v = x_next, lty = 2)
```

Example 1 v



When does it go wrong?

- · Newton-Raphson usually goes wrong in one of two ways:
 - $\cdot \nabla f(\mathbf{x}_0) = 0$ and you're stuck at \mathbf{x}_0 .
 - The sequence \mathbf{x}_n does not converge.
- That's why we kept track of the number of iterations, so that we aren't stuck in an infinite loop.
- · When you reach the maximum number of iterations, you can:
 - Try a different initial value.
 - Look at a graph and see if there is indeed a maximum/minimum.
 - · Consider using a different approach (e.g. gradient descent).

Example (not convergent)

- Consider $f(x) = 0.25x^4 x^2 + 2x$.
- The derivatives are

$$f'(x) = x^3 - 2x + 2$$
, $f''(x) = 3x^2 - 2$.

· If we start at $x_0 = 0$, then at the next iteration we get

$$x_1 = x_0 - \frac{f'(x_0)}{f''(x_0)} = 0 - \frac{2}{-2} = 1.$$

Next we get

$$x_2 = x_1 - \frac{f'(x_1)}{f''(x_1)} = 1 - \frac{1}{1} = 0.$$

- · Therefore, we are stuck in an infinite loop.
 - · Solution: Use a different starting value.

Exercise

Find the minimum for the function

$$f(x) = 5x + 5x^2 + 5\log(1 + e^{-x}).$$

Solution i

We need to compute the first and second derivatives:

$$f'(x) = 5 + 10x - \frac{5e^{-x}}{1 + e^{-x}} = 5 + 10x - \frac{5}{1 + e^{x}},$$
$$f'' = 10 + \frac{5e^{x}}{(1 + e^{x})^{2}}.$$

Solution ii

```
gradient <- function(x) 5 + 10*x - 5/(1+exp(x))
hessian <- function(x) 10 + 5*exp(x)/(1+exp(x))^2

x_current <- 1
x_next <- x_current -
    gradient(x_current)/hessian(x_current)

tol <- 10^-10
iter <- 1</pre>
```

Solution iii

```
while (abs(x_current - x_next) > tol &
       iter < 100) {
  iter <- iter + 1
    x current <- x next
    step <- gradient(x current)/hessian(x current)</pre>
    x next <- x current - step
x next
```

```
## [1] -0.2223235
```

Poisson regression i

- Poisson regression is a generalization of linear regression where the outcome variable follows a Poisson distribution.
- Let (Y_i, X_i) , $i = 1, \ldots, n$, be pairs of outcome and covariate variables. We assume the following relationship:

$$Y_i \sim Pois(\mu(X_i)), \quad \log(\mu(X_i)) = \beta_0 + \beta_1 X_i.$$

- · We will use the Newton-Raphson method to find the Maximum Likelihood estimate $\hat{\theta}=(\hat{\beta}_0,\hat{\beta}_1)$.
- · First, let's write down the likelihood:

$$L(\theta) = \prod_{i=1}^{n} \frac{\mu(X_i)^{Y_i} \exp(-\mu(X_i))}{Y_i!}.$$

Poisson regression ii

· Take the logarithm:

$$\ell(\theta) = \sum_{i=1}^{n} Y_i \log(\mu(X_i)) - \mu(X_i) - \log(Y_i!)$$

$$= \sum_{i=1}^{n} Y_i (\beta_0 + \beta_1 X_i) - \exp(\beta_0 + \beta_1 X_i) - \log(Y_i!)$$

$$\propto \sum_{i=1}^{n} Y_i (\beta_0 + \beta_1 X_i) - \exp(\beta_0 + \beta_1 X_i).$$

Poisson regression iii

· The first-order derivatives are

$$\frac{\partial \ell(\theta)}{\partial \beta_0} = \sum_{i=1}^n Y_i - \exp(\beta_0 + \beta_1 X_i),$$

$$\frac{\partial \ell(\theta)}{\partial \beta_1} = \sum_{i=1}^n Y_i X_i - X_i \exp(\beta_0 + \beta_1 X_i).$$

Poisson regression iv

· The second-order derivatives are

$$\frac{\partial^2 \ell(\theta)}{\partial \beta_0^2} = -\sum_{i=1}^n \exp(\beta_0 + \beta_1 X_i),$$

$$\frac{\partial^2 \ell(\theta)}{\partial \beta_1^2} = -\sum_{i=1}^n X_i^2 \exp(\beta_0 + \beta_1 X_i),$$

$$\frac{\partial^2 \ell(\theta)}{\partial \beta_0 \partial \beta_1} = -\sum_{i=1}^n X_i \exp(\beta_0 + \beta_1 X_i).$$

Poisson regression v

- We will fit a Poisson regression model to the "famous" horseshoe crab dataset.
 - · The main covariate is the weight of female crabs.
 - The outcome is the number of satellite males for that female crab.
 - The dataset can be found in the **asbio** package.

Poisson regression vi

```
library(asbio)
data(crabs)
y_vec <- crabs$satell</pre>
x vec <- crabs$weight
gradient <- function(theta) {</pre>
  mu x <- exp(theta[1] + theta[2]*x vec)
  ell b0 < -sum(y vec - mu x)
  ell_b1 <- sum(y_vec*x_vec - x_vec*mu_x)
  return(c(ell_b0, ell_b1))
```

Poisson regression vii

```
hessian <- function(theta) {</pre>
 mu_x <- exp(theta[1] + theta[2]*x_vec)</pre>
  ell b0b0 < - -sum(mu x)
  ell b1b1 <- -sum(x vec^2*mu x)
  ell b0b1 < - -sum(x vec*mu x)
  hess <- matrix(c(ell b0b0, ell b0b1,
                    ell b0b1, ell b1b1),
                  ncol = 2)
  return(hess)
```

Poisson regression viii

Poisson regression ix

```
while(sum((x current - x next)^2) > tol &
      iter < 100) {
  iter <- iter + 1
  x_current <- x_next
  step <- solve(hessian(x_current),</pre>
                 gradient(x current))
  x next <- x current - step
}
x next
```

```
## [1] -0.4284053 0.5893041
```

Poisson regression x

• We can compare this to the built-in function glm in R:

Poisson regression xi

```
library(tidyverse)
log lik <- function(row) {</pre>
    mu_vec <- exp(row$beta0 + row$beta1*x vec)</pre>
    sum(dpois(y_vec, lambda = mu_vec, log = TRUE))
}
data plot \leftarrow expand.grid(beta0 = seq(-2, 0, 0.1),
                           beta1 = seq(0, 1, 0.1)) \%>\%
  purrrlyr::by_row(log_lik, .collate = "cols",
                     .to = "loglik")
```

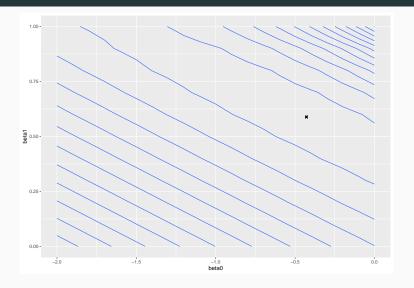
Poisson regression xii

head(data_plot)

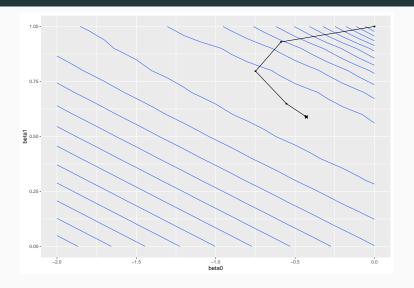
```
## # A tibble: 6 x 3
     beta0 beta1 loglik
##
##
     <dbl> <dbl> <dbl>
## 1 -2
              0 -1563.
## 2 -1.9
              0 -1515.
## 3 -1.8
              0 -1468.
## 4 -1.7
              0 - 1420.
## 5 -1.6
              0 -1373.
## 6 -1.5
              0 - 1326.
```

Poisson regression xiii

Poisson regression xiv



Poisson regression xv



Some vocabulary

- In the context of Maximum Likelihood estimation, the gradient and Hessian are given different names.
- The gradient $\nabla \ell(\theta \mid \mathbf{x})$ is called the score function.
 - Exercise: The expected value of the score function is 0.
- The Hessian $D^2\ell(\theta \mid \mathbf{x})$ is related to the Fisher information matrix $\mathcal{I}(\theta)$:

$$\mathcal{I}(\theta) = -E\left(D^2\ell(\theta \mid \mathbf{x})\right).$$

Final remarks

- · Newton-Raphson is an important optimisation method.
 - In particular, it is commonly used with generalized linear models.
- There exists optimisation methods that are gradient-free.
 - · See Nelder-Mead and the function optim.
- · Where to go from here:
 - · Gradient descent
 - · Regularized regression (e.g. lasso)
 - · MM algorithms