# R Programming and Advanced Data Analysis Bayesian Computing

#### Introduction

- 1. Approximating integrals
  - a. Monte Carlo
  - b. Importance sampling
- 2. Markov Chain Monte Carlo
  - a. Gibbs sampler
  - b. Metropolis-Hastings sampler

() Bayesian Computing Introduction 2/40

## **Monte Carlo**

- Suppose  $\mathbf{X} = (X_1, \dots, X_n)$  is an iid sample from Beta(a, b).
- We would like to compute the moment generating function

$$\theta = \int_0^1 e^{tx} \frac{x^{a-1}(1-x)^{b-1}}{B(a,b)} dx.$$

- There are no unknown parameters involved, but exact calculation involves a difficult integration.
- However, we can use a Monte Carlo (MC) method to approximate  $\theta$ , to any desired degree of accuracy.
- **Problem**: How many samples are required to construct a confidence interval for  $\theta$  of width 0.1?

- Let  $g(x) = e^{tx}$  be the integrand.
- An MC estimate for  $\theta$  is

$$\widehat{\theta} = \frac{1}{R} \sum_{r=1}^{R} g(X_r).$$

By the Law of Large Numbers,  $\widehat{\theta} \overset{\text{a.s.}}{\to} \theta$  as  $R \to \infty$ .

• We have  $Var(\widehat{\theta}) = \frac{1}{R}Var[g(X)]$ . Approximate  $\sigma^2 = Var[g(X)]$  with

$$\hat{\sigma}^2 = \frac{1}{R} \sum_{r=1}^{R} \left( g(X_r) - \widehat{\theta} \right)^2.$$

• By the Central Limit Theorem

$$\frac{\sqrt{R}\left[\widehat{\theta} - E(\widehat{\theta})\right]}{\widehat{\Phi}} \xrightarrow{\mathcal{L}} \mathsf{N}(0,1).$$

ullet Using this, we can obtain an approximate (1-lpha) level CI for heta as

$$\widehat{\theta} \pm z_{\alpha/2} \frac{\widehat{\sigma}}{\sqrt{R}}$$
.

()

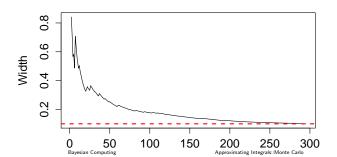
#### Algorithm

- 1. Sample  $X_i$  from Beta(a, b)
- 2. Until width of CI  $= 2 \cdot z_{\alpha/2} \, {\hat \sigma \over \sqrt R}$  is less than  $\delta = 0.1$

```
t <- 1; a <- 2; b <- 1/2
g <- function(x) { exp(t*x) }
alpha <- 0.05
delta <- 0.1
z <- rbeta(1, a, b)
widths <- c(Inf)
while(tail(widths, 1) >= delta) {
    z \leftarrow c(z, rbeta(1, a, b))
    R <- length(z)
    est.mc <- mean(g(z))
    var.mc < -1/R * (R-1)/R * var(g(z))
    widths <- c(widths, 2 * qnorm(1 - alpha/2) * sqrt(var.mc))
ci \leftarrow est.mc + c(-1,1) * qnorm(1 - alpha/2) * sqrt(var.mc)
```

```
> cat("Estimate is", est.mc, "\n")
> cat("CI = [", ci, "], width =", ci[2] - ci[1], "\n")
> cat("Number of samples needed:", length(z), "\n")
> plot(widths, type="l", xlab="Iteration", ylab="Width")
> abline(h = delta, lwd = 2, lty = 2, col = "red")
```

```
Estimate is 2.240291
CI = [ 2.190351 2.290232 ], width = 0.09988176
Number of samples needed: 295
```



- Importance Sampling is a technique to reduce the variance in Monte Carlo approximation (Owen, 2013; Jones et al., 2009).
- Let supp $(f) = \{x : f(x) \neq 0\}$  for a function f.
- Suppose we want to compute

$$\theta = \int_{\mathsf{supp}(p)} g(x) p(x) dx$$

where p is a density ("the nominal distribution") and g is a given function.

• Suppose we have another density q ("the importance distribution") with  $supp(g \cdot p) \subseteq supp(q)$ . We may write

$$\begin{split} \theta &= \int_{\text{supp}(\rho)} g(x) p(x) dx \\ &= \int_{\text{supp}(\rho)} \frac{g(x) p(x)}{q(x)} q(x) dx \\ &= \int_{\text{supp}(q)} \frac{g(x) p(x)}{q(x)} q(x) dx - \underbrace{\int_{\text{supp}(q) \setminus \text{supp}(\rho)} \frac{g(x) p(x)}{q(x)} q(x) dx}_{=0} \\ &= \mathsf{E} \left[ \frac{g(X) p(X)}{q(X)} \right], \quad \text{where } X \sim q. \end{split}$$

• Another Monte Carlo estimator is then

$$\hat{\theta}_q = \frac{1}{R} \sum_{r=1}^{R} \frac{g(X_r)p(X_r)}{g(X_r)}, \quad X_1, \dots, X_R \stackrel{\text{iid}}{\sim} q.$$

• By the Law of Large Numbers,  $\hat{\theta}_a \stackrel{a.s.}{\to} \theta$  as  $R \to \infty$ .

• Can also show that  $\mathsf{E}(\hat{ heta}_q) = heta$  and  $\mathsf{Var}(\hat{ heta}_q) = \sigma_q^2/R$ , where

$$\sigma_q^2 = \int_{\text{supp}(p)} \frac{[g(x)p(x)]^2}{q(x)} dx - \theta^2$$

- To get a variance reduction, choose q so that it gives higher weight to the region of supp(p) where |g(x)|p(x) is large.
- Also, q must give enough weight to regions where |g(x)|p(x) is nonzero. I.e. tails of q cannot be too light.
- The optimal variance is achieved at

$$q^*(x) = \frac{|g(x)|p(x)}{\int_{\mathsf{SUPP}(p)} |g(z)|p(z)dz},$$

but this result is usually not useful, since computing the denominator is as hard as the original integration problem.

- The variance  $\sigma_q^2$  can be estimated by  $\hat{\sigma}_q^2$ , the sample variance of  $g(X_r)p(X_r)/q(X_r)$  for  $r=1,\ldots,R$ .
- The Central Limit Theorem gives that

$$\sqrt{R} rac{(\hat{ heta}_q - heta)}{\hat{\sigma}_q} \stackrel{\mathcal{L}}{ o} \mathsf{N}(0,1) \quad ext{as } R o \infty.$$

• Using this, we can obtain an approximate  $(1-\alpha)$  level confidence interval for  $\theta$  as

$$\hat{\theta}_q \pm z_{\alpha/2} \frac{\hat{\sigma}_q}{\sqrt{R}}.$$

- The following example is from Jones et al. (2009, Section 20.2.3).
- Suppose we would like to compute

$$\theta = \int_0^1 e^{-x^2/2} dx = \sqrt{2\pi} \, \mathsf{P}(0 < Z < 1), \quad \text{for } Z \sim \mathsf{N}(0,1).$$

• A simple/obvious Monte Carlo estimator is

$$\hat{\theta} = \frac{\sqrt{2\pi}}{R} \sum_{r=1}^{R} I(0 < Z_r < 1), \quad \text{for } Z_1, \dots, Z_R \stackrel{\text{iid}}{\sim} \mathsf{N}(0,1)$$

• We have that

$$\sum_{r=1}^{R} I(0 < Z_r < 1) \sim \mathsf{Binomial}(R, \theta)$$

$$\implies \mathsf{Var}(\hat{\theta}) = \frac{2\pi}{R} \theta (1 - \theta) \approx \frac{1.413}{R}$$

It turns out that we can do much better...

Recall our objective

$$\theta = \int_0^1 e^{-x^2/2} dx = \sqrt{2\pi} \, \mathsf{P}(0 < Z < 1), \quad \text{for } Z \sim \mathsf{N}(0,1).$$

• A 2nd order Taylor series approximation around x = 0 yields

$$e^{x^2/2} \approx 1 + x^2/2 \implies e^{-x^2/2} \approx \frac{1}{1 + x^2/2}$$

• The expression  $1/(1+x^2/2)$  can be turned into a density on  $x \in (0,1)$ ;

$$q(x) = \frac{1/(1+x^2/2)}{\int_0^1 1/(1+z^2/2)dz} = \frac{1/(1+x^2/2)}{\sqrt{2}\arctan(1/\sqrt{2})}.$$

Using q as the importance density, our estimator becomes

$$\hat{\theta}_q = \frac{1}{R} \sum_{r=1}^{R} \frac{g(X_r)p(X_r)}{g(X_r)} = \frac{1}{R} \sum_{r=1}^{R} e^{-X_r^2/2} \sqrt{2} \arctan(1/\sqrt{2})(1+X_r^2/2),$$

where  $X_1, \ldots, X_R \stackrel{\text{iid}}{\sim} q$ .

- q is a nonstandard density, but we can sample from it using the inverse CDF method.
- The inverse CDF method can be used with continuous univariate distributions where the cumulative distribution function (CDF) is available.
- The CDF of q is

$$Q(x) = \int_0^x q(x) dx = \int_0^x \frac{1/(1 + x^2/2)}{\sqrt{2} \arctan(1/\sqrt{2})} dx = \frac{\arctan(x/\sqrt{2})}{\arctan(1/\sqrt{2})}$$

• The inverse of Q(x) is

$$Q^{-1}(u) = \sqrt{2} \tan \left( \operatorname{arctan}(1/\sqrt{2}) \right)$$

• Taking  $X = Q^{-1}(U)$  with  $U \sim U(0,1)$ , it can be shown that  $X \sim q$ .

• The exact answer (computed numerically)

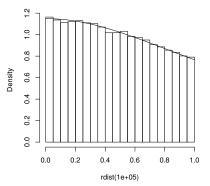
```
> sqrt(2*pi) * (pnorm(1) - pnorm(0))
[1] 0.8556244
```

• Simple Monte Carlo estimator. This time fix R.

```
g \leftarrow function(x) \{ sqrt(2*pi)*(x > 0 & x < 1) \}
alpha <- 0.05
delta <- 0.1
R <- 10000
widths.simple <- numeric(R)
z < - rnorm(R)
for (r in 1:R){
    w \leftarrow g(z[1:r])
    est.mc <- mean(w)
    var.mc <- 1/r * (r-1)/r * var(w)
    widths.simple[r] <- 2 * qnorm(1 - alpha/2) * sqrt(var.mc)
ci \leftarrow est.mc + c(-1,1) * qnorm(1 - alpha/2) * sqrt(var.mc)
cat("Estimate is", est.mc, "\n")
cat("CI = [", ci, "], width = ", ci[2] - ci[1], "\n")
```

```
Estimate is 0.8632828
CI = [ 0.839938 0.8866275 ], width = 0.04668948
```

```
# Density and drawing function for the truncated Cauchy-like
    distribution.
rdist <- function(n) {
    u <- runif(n)
    sqrt(2) * tan(u * atan(1/sqrt(2)))
}
ddist <- function(x) { 1 / ((1 + x^2/2) * sqrt(2) * atan(1 / sqrt(2))) }
# A quick test to make sure they are consistent with each other
hist(rdist(100000), freq = FALSE, main = NULL)
curve(ddist, add = TRUE)</pre>
```

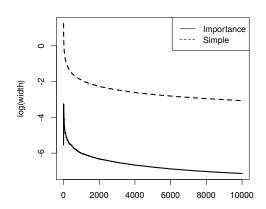


#### Importance sampling estimator

```
widths.imp <- numeric(R)
x <- rdist(R)
for (r in 1:R){
    w <- exp(-x[1:r]^2/2) * sqrt(2) * atan(1/sqrt(2)) * (1 + x[1:r]^2/2)
    est.mc <- mean(w)
    var.mc <- 1/r * (r-1)/r * var(w)
    widths.imp[r] <- 2 * qnorm(1 - alpha/2) * sqrt(var.mc)
}
ci <- est.mc + c(-1,1) * qnorm(1 - alpha/2) * sqrt(var.mc)
cat("Estimate is", est.mc, "\n")
cat("CI = [", ci, "], width =", ci[2] - ci[1], "\n")
cat("Number of samples used:", R, "\n")</pre>
```

```
> cat("Estimate is", est.mc, "\n")
Estimate is 0.8554805
> cat("CI = [", ci, "], width =", ci[2] - ci[1], "\n")
CI = [ 0.8550783 0.8558828 ], width = 0.0008045813
```

Compare accuracy of simple and Importance Sampling estimator



() Bayesian Computing Markov Chain Monte Carlo 20/40

- Markov Chain Monte Carlo (MCMC) describes a class of algorithms which produce approximate draws from a given density f(x).
- Often, the functional form of f is known except for a normalizing constant.
- Even when f is fully specified, may not be easy to draw directly from it.
- An MCMC algorithm produces a sequence X<sub>1</sub>, X<sub>2</sub>,... of dependent draws. The sequence is constructed in such a way that it "behaves like" draws from f more as we progress through the sequence.
- Technical justifications for MCMC are beyond the scope of the workshop, but are discussed in books Robert and Casella (2005) and Meyn and Tweedie (2009). A more applied book for practitioners is Gelman et al. (2013).

() Bayesian Computing Markov Chain Monte Carlo 21/40

- MCMC algorithms have found widespread use in the Bayesian paradigm of statistics.
- The fundamental assumption in Bayesian statistics is the framework

$$\mathbf{X} \sim f(\mathbf{x} \mid \mathbf{\theta})$$
  $\leftarrow$  likelihood,  $\mathbf{\theta} \sim f(\mathbf{\theta})$   $\leftarrow$  prior.

- The prior distribution represents our apriori belief about the parameter θ. Often, we have no idea, and the prior is chosen for mathematical convenience and/or to express our lack of knowledge.
- Using the likelihood, a prior, and observed data x, all inference about  $\theta$  should be based on the **posterior** distribution

$$f(\theta \mid \mathbf{x}) = \frac{f(\mathbf{x} \mid \theta)f(\theta)}{f(\mathbf{x})} \propto f(\mathbf{x} \mid \theta)f(\theta)$$

• The distribution  $f(\theta \mid x)$  can sometimes be recognizable. E.g., sometimes  $f(x \mid \theta)$  and  $f(\theta)$  may be chosen so that the posterior will be in the same family as the prior (i.e. *conjugate* prior).

() Bayesian Computing Markov Chain Monte Carlo 22/40

- In most settings,  $f(\theta \mid x)$  will not have a recognizable form.
- Typically, the denominator of the posterior density,

$$f(\mathbf{x}) = \int f(\mathbf{x} \mid \boldsymbol{\theta}) f(\boldsymbol{\theta}) d\theta,$$

will be too complicated to evaluate analytically.

- MCMC algorithms allow us to generate draws  $\theta^{(1)}, \theta^{(2)}, \ldots$  from the posterior without knowing the denominator.
- MCMC algorithms are also useful when  $f(\theta \mid x)$  is completely known, but does not have a familiar form.

() Bayesian Computing Markov Chain Monte Carlo 23/40

 Another important distribution for Bayesians is the posterior predictive distribution,

$$f(\tilde{\mathbf{x}} \mid \mathbf{x}) = \int f(\tilde{\mathbf{x}} \mid \boldsymbol{\theta}) f(\boldsymbol{\theta} \mid \mathbf{x}) d\theta,$$

where  $\tilde{x}$  represents a new observation.

- The posterior predictive distribution allows inference on new observations through the posterior distribution.
- $f(\tilde{x} \mid x)$  is usually too complicated to determine analytically. But we can generate draws from it by

$$\tilde{\mathbf{X}}^{(r)} \sim f(\tilde{\mathbf{x}} \mid \boldsymbol{\theta}^{(r)}),$$
  
 $\boldsymbol{\theta}^{(r)} \sim f(\boldsymbol{\theta} \mid \mathbf{x}).$ 

The draws  $\theta^{(r)}$ ,  $r=1,\ldots,R$ , from the posterior are obtained using the MCMC sampler.

() Bayesian Computing Markov Chain Monte Carlo 24/40

- Math Gym (http://mathgym.umbc.edu) is a program at UMBC which provides "workouts" to aid students in introductory math courses.
- Question: Is there evidence that it helps students pass the courses?
- Denote Y<sub>i</sub> as the indicator of whether the ith student passed. Let x<sub>i</sub> contain predictors computed from the course, section, FYI, # of days attending math gym, and Quiz Zero score.
- A Bayesian probit regression model is

$$Y_i \sim \mathsf{Ber}(p_i), \quad p_i = \Phi(\mathbf{x}_i^{\top} \boldsymbol{\beta}), \qquad \leftarrow \mathsf{likelihood}, \\ \boldsymbol{\beta} \sim \mathsf{N}(\boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}), \qquad \leftarrow \mathsf{prior}.$$

- Here,  $\Phi(\cdot)$  is the CDF of N(0,1);  $\mu_{\beta}$  and  $\Sigma_{\beta}$  are hyperparameters to describe our uncertainty of  $\beta$  before observing the data. We will take  $\mu_{\beta} = \mathbf{0}$  and  $\Sigma_{\beta} = \sigma_{\beta}^2 \mathbf{I}$  for some  $\sigma_{\beta}^2 >> 1$ .
- The joint distribution of  $\mathbf{Y} = (Y_1, \dots, Y_n)$  and  $\beta$  is

$$f(\boldsymbol{y}, \boldsymbol{\beta}) = \mathsf{N}(\boldsymbol{\beta} \mid \boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}) \cdot \prod_{i=1}^{n} \left[ \Phi(\boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}) \right]^{y_{i}} \left[ 1 - \Phi(\boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}) \right]^{1 - y_{i}}.$$

• We will make use of the basis function decomposition

$$f(x) = \sum_{j=1}^{J} \beta_j h_j(x),$$

where  $h_1(x), \ldots, h_J(x)$  are prespecified basis functions and  $\beta_1, \ldots, \beta_J$  are unknown coefficients. Chapter 5 of Hastie et al. (2013) discusses the basics.

• We will specify K knot points  $\xi_1 < \ldots < \xi_K$  so that the behavior of f(x) can be modeled differently in the regions

$$(-\infty, \xi_1), (\xi_1, \xi_2), \dots, (\xi_{K-1}, \xi_K), (\xi_K, \infty).$$

• To support a piecewise linear function, whose pieces are connected at the knots (i.e. continuous), an appropriate basis is

$$h_1(x) = 1, h_2(x) = x,$$
  
 $h_3(x) = (x - \xi_1) \cdot I(x > \xi_1), \dots, h_{K+2}(x) = (x - \xi_K) \cdot I(x > \xi_K).$ 

## **Basis Example**

... Demonstration ...

(See basis-example.Rmd)

- The receiver operating characteristic (ROC) curve is often used to evaluate classifiers when there are two classes (Fawcett, 2003).
- There are four possible outcomes.

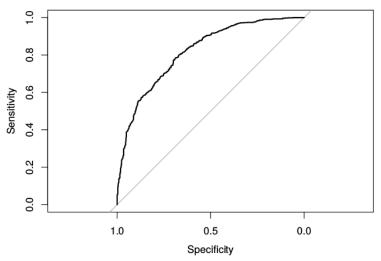
Classify	$True\ Class = 1$	True Class $= 0$
1	True Positive	False Positive
0	False Negative	${f T}$ rue ${f N}$ egative

- An ROC plot displays Sensitivity =  $\frac{TP}{TP+FN}$  vs. Specificity =  $\frac{FP}{FP+TN}$ .
- Classification methods such as probit regression produce an estimate  $\hat{p}(x) = \hat{P}(\text{Class} = 1 \mid x)$ . To obtain a predicted class, we often consider

$$\widehat{\mathsf{Class}}(\mathbf{x}) = \begin{cases} 1 & \text{if } \hat{p}(\mathbf{x}) > c, \\ 0 & \text{o.w.} \end{cases}$$

• We can produce an ROC curve by considering all the possible cutoffs

$$\Big\{(\mathsf{Sensitivity}(c),\mathsf{Specificity}(c)):c\in[0,1]\Big\}.$$



AUC is the Area Under the Curve from x = 1 to x = 0.

## **Metropolis-Hastings Sampler**

## **Metropolis-Hastings Sampler**

- Suppose f(x) is our target density, from which we want to draw. Let  $q(x^* \mid x)$  be a proposal density which should be easy to draw from.
- Starting with a given  $\boldsymbol{X}^{(0)}$ 
  - 1. Draw  $X^* \sim q(X^* \mid X^{(0)})$
  - 2. Draw  $U \sim \text{Uniform}(0,1)$  and let

$$\boldsymbol{\mathit{X}}^{(1)} = \left\{ \begin{aligned} \boldsymbol{\mathit{X}}^* & \text{if } \boldsymbol{\mathit{U}} < \min \left\{ 1, \frac{f(\boldsymbol{\mathit{X}}^*)}{f(\boldsymbol{\mathit{X}}^{(0)})} \frac{q(\boldsymbol{\mathit{X}}^{(0)} | \boldsymbol{\mathit{X}}^*)}{q(\boldsymbol{\mathit{X}}^* | \boldsymbol{\mathit{X}}^{(0)})} \right\}, \\ \boldsymbol{\mathit{X}}^{(0)} & \text{o.w.} \end{aligned} \right.$$

Repeat starting from  $\boldsymbol{X}^{(1)}$ , then  $\boldsymbol{X}^{(2)}$ , etc to generate a chain which eventually behaves like a draw from  $f(\boldsymbol{x})$ .

- A common choice of  $q(\mathbf{x}^* \mid \mathbf{x})$  is  $N(\mathbf{x}^* \mid \mathbf{x}, \mathbf{V})$ . The variance  $\mathbf{V}$  controls how candidate points are drawn and is a tuning parameter. This choice simplifies step 2, since  $q(\mathbf{x}^* \mid \mathbf{x}) = q(\mathbf{x} \mid \mathbf{x}^*)$ .
- For Bayesians: the normalizing constant of f is not needed, since

$$\frac{f(\boldsymbol{\theta}^* \mid \mathsf{Data})}{f(\boldsymbol{\theta} \mid \mathsf{Data})} = \frac{f(\mathsf{Data} \mid \boldsymbol{\theta}^*)f(\boldsymbol{\theta}^*)}{f(\mathsf{Data} \mid \boldsymbol{\theta})f(\boldsymbol{\theta})}, \quad \text{which is easy to compute.}$$

## **Metropolis-Hastings Example**

## ... Demonstration ...

(See sim.Rmd and mathgym.Rmd)

- Again suppose that f(x) is our target density, and we wish to draw  $X = (X_1, \dots, X_k)$  from f.
- Starting with a given  $\boldsymbol{X}^{(0)}$ , obtain  $\boldsymbol{X}^{(1)}$  by the following.
  - 1. Draw  $X_1^{(1)} \sim f(X_1 \mid X_2^{(0)}, X_2^{(0)}, \dots, X_k^{(0)})$ .
  - 2. Draw  $X_2^{(1)} \sim f(X_2 \mid X_1^{(1)}, X_3^{(0)}, \dots, X_k^{(0)})$ .

:

- k. Draw  $X_k^{(1)} \sim f(X_1 \mid X_1^{(1)}, X_2^{(1)}, \dots, X_{k-1}^{(1)}).$
- Repeat starting from  $\boldsymbol{X}^{(r)}$  to obtain  $\boldsymbol{X}^{(r+1)}$  for  $r=1,2,\ldots$
- Eventually the chain starts to behave like a draw from f(x).
- More generally, the Gibbs sampler can be applied when  $\mathbf{X} = (\mathbf{X}_1, \dots, \mathbf{X}_k)$  is decomposed into vectors of size  $\geq 1$ .
- For many problems, we can readily sample from the k conditional distributions.

• For probit regression, Albert and Chib (1993) suggest an augmented data model which leads to a convenient Gibbs sampler,

$$Y_i = I(Z_i > 0), \quad i = 1, ..., n$$
  
 $Z_i \sim \mathsf{N}(\mathbf{x}_i^{\top} \boldsymbol{\beta}, 1)$   
 $\boldsymbol{\beta} \sim \mathsf{N}(\boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}).$ 

• The joint density of all random variables is

$$f(\mathbf{y}, \mathbf{z}, \boldsymbol{\beta}) = \prod_{i=1}^{n} \left\{ \left[ I(z_i > 0) \right]^{y_i} \left[ I(z_i \leq 0) \right]^{1-y_i} \mathsf{N}(z_i \mid \mathbf{x}_i^{\top} \boldsymbol{\beta}, 1) \right\} \times \mathsf{N}(\boldsymbol{\beta} \mid \boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}).$$

• Integrating z out of the augmented model yields the original model.

• To derive the step for drawing z, we have

$$\begin{split} f(\boldsymbol{z} \mid \boldsymbol{\beta}, \boldsymbol{y}) &\propto \prod_{i=1}^{n} \left\{ \left[ \textit{I}(z_{i} > 0) \right]^{y_{i}} \left[ \textit{I}(z_{i} \leq 0) \right]^{1-y_{i}} \mathsf{N}(z_{i} \mid \boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}, 1) \right\} \times \\ & \mathsf{N}(\boldsymbol{\beta} \mid \boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}) \\ &\propto \prod_{i=1}^{n} \left\{ \left[ \textit{I}(z_{i} > 0) \right]^{y_{i}} \left[ \textit{I}(z_{i} \leq 0) \right]^{1-y_{i}} \mathsf{N}(z_{i} \mid \boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}, 1) \right\} \end{split}$$

• Therefore,

$$[\boldsymbol{z} \mid \boldsymbol{\beta}, \boldsymbol{y}] = [z_1 \mid \boldsymbol{\beta}, \boldsymbol{y}] \cdots [z_n \mid \boldsymbol{\beta}, \boldsymbol{y}]$$

where

$$[z_i \mid \boldsymbol{\beta}, \mathbf{y}] \sim \begin{cases} I(z_i > 0) \cdot \mathsf{N}(z_i \mid \mathbf{x}_i^{\top} \boldsymbol{\beta}, 1), & \text{if } y_i = 1, \\ I(z_i \leq 0) \cdot \mathsf{N}(z_i \mid \mathbf{x}_i^{\top} \boldsymbol{\beta}, 1), & \text{if } y_i = 0. \end{cases}$$

• To derive the step for drawing  $\beta$ , we have

$$\begin{split} f(\boldsymbol{\beta} \mid \boldsymbol{z}, \boldsymbol{y}) &\propto \prod_{i=1}^{n} \left\{ \left[ I(z_{i} > 0) \right]^{y_{i}} \left[ I(z_{i} \leq 0) \right]^{1-y_{i}} \mathsf{N}(z_{i} \mid \boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}, 1) \right\} \times \\ & \mathsf{N}(\boldsymbol{\beta} \mid \boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}) \\ &\propto \prod_{i=1}^{n} \left\{ \mathsf{N}(z_{i} \mid \boldsymbol{x}_{i}^{\top} \boldsymbol{\beta}, 1) \right\} \mathsf{N}(\boldsymbol{\beta} \mid \boldsymbol{\mu}_{\boldsymbol{\beta}}, \boldsymbol{\Sigma}_{\boldsymbol{\beta}}) \end{split}$$

• After some calculation, it can be shown that

$$[eta \mid oldsymbol{z}, oldsymbol{y}] = [eta \mid oldsymbol{z}] \sim \mathsf{N}( ilde{oldsymbol{\mu}}_eta, ilde{oldsymbol{\Sigma}}_eta),$$

where 
$$\tilde{\Sigma}_{\beta} = (\Sigma_{\beta}^{-1} + \pmb{X}^{ op} \pmb{X})^{-1}$$
 and  $\tilde{\mu}_{\beta} = \tilde{\Sigma}_{\beta} (\Sigma_{\beta}^{-1} \mu_{\beta} + \pmb{X}^{ op} \pmb{z})$ .

## **Gibbs Example**

... Demonstration ...

(See sim.Rmd and mathgym.Rmd)

#### References I

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