

Bayesian Regression

Bayesian Models for Ecologists

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Outline

- Be able to write proper Bayesian regression models for different types of data.
- Appreciate one-to-one relationship between math and JAGS code.
- Be able to interpret coefficients of general linear models.

A great follow-up

This book should be in your library:



The general Bayesian set-up

Recall that the posterior distribution of the unobserved quantities conditional on the observed ones is proportional to their joint distribution:

$$[\theta, \sigma^2 | \mathbf{y}] \propto [\theta, \sigma^2, \mathbf{y}].$$

The joint distribution can be factored into a likelihood and priors for simple Bayesian models:

$$[\theta, \sigma^2, \mathbf{y}] = [\mathbf{y} | \theta, \sigma^2] [\theta, \sigma^2] \stackrel{\text{ind.}}{=} [\mathbf{y} | \theta, \sigma^2] [\theta] [\sigma^2]$$

A deterministic model of an ecological or socioenvironmental process is embedded in the likelihood like this:

$$[\theta, \sigma^2, \mathbf{y}] \propto [\mathbf{y} | g(\theta, x), \sigma^2] [\theta] [\sigma^2]$$

Simple Bayesian regression models

We use likelihood to connect the underlying process to data:

$$\underbrace{[y_i \mid \mu_i, \sigma^2]}_{\text{stochastic model}}, \quad i = 1, \dots, n$$

We formulate the deterministic model:

$$\mu_i = \underbrace{g(\beta, \mathbf{x}_i)}_{\text{deterministic model}}, \quad i = 1, \dots, n$$

where β is a vector of regression coefficients and $\mathbf{x}_i$ is a vector of predictor variables.

Assuming conditional independence of the data,

$$[\beta, \sigma^2 \mid \mathbf{y}] \propto \prod_{i=1}^n [y_i \mid g(\beta, \mathbf{x}_i), \sigma^2] \times [\beta] [\sigma^2]$$

We choose appropriate deterministic functions (linear or non-linear) and appropriate probability distributions to compose specific models.

Identical notation

$$y_i = g(\beta, x_i) + \varepsilon_i$$
$$\varepsilon_i \sim \text{Normal}(0, \sigma^2)$$

is the same as:

$$y_i | \beta, \sigma^2 \sim \text{Normal}(g(\beta, x_i), \sigma^2),$$

but the second notation is much more flexible because it generalizes to distributions that do not have additive errors.

You don't have to be Normal!

Assuming you have one predictor x :

Data (y-values)	Distribution	"Mean" function	Link
continuous, real valued	Normal	$\mu = \beta_0 + \beta_1 x$	NA (i.e. identity)
discrete, strictly positive	Poisson	$\mu = e^{\beta_0 + \beta_1 x}$	$\log(\mu) = \beta_0 + \beta_1 x$
0 or 1	Bernoulli	$\mu = \frac{\exp(\beta_0 + \beta_1 x)}{\exp(\beta_0 + \beta_1 x) + 1}$	$\text{logit}(\mu) = \log\left(\frac{\mu}{1-\mu}\right) = \beta_0 + \beta_1 x$
[0, 1]	Beta	$\mu = \frac{\exp(\beta_0 + \beta_1 x)}{\exp(\beta_0 + \beta_1 x) + 1}$	$\text{logit}(\mu) = \log\left(\frac{\mu}{1-\mu}\right) = \beta_0 + \beta_1 x$
continuous, strictly positive, variance $\uparrow$ as a f(mean)	lognormal	$\mu = e^{\beta_0 + \beta_1 x}$	$\log(\mu) = \beta_0 + \beta_1 x$
continuous, strictly positive, constant variance	Gamma	$\mu = e^{\beta_0 + \beta_1 x}$	$\log(\mu) = \beta_0 + \beta_1 x$

Continuous and real valued data

Suppose you have collected some continuous data $\mathbf{y} = (-10.7, -4.3, \dots, 49)$ at n sites, along with a predictor x_i measured at each site i which you believe is likely to affect these measurements. Write a model regressing y on x as follows:

- 1 Choose a specific stochastic and deterministic model.
- 2 Specify (vague) priors for your parameters.
- 3 Write out the DAG and express posterior distribution as proportional to joint distribution for your model.
- 4 Write the JAGS code for the model.
- 5 Interpret the coefficients of your model.

Normal data, continuous and real valued

Stochastic model:

$$y_i | \beta_0, \beta_1, \sigma^2 \stackrel{\text{ind.}}{\sim} \text{Normal}(g(\beta_0, \beta_1, x_i), \sigma^2) \quad i = 1, \dots, n$$

Deterministic model:

$$\mu_i = g(\beta_0, \beta_1, x_i) = \beta_0 + \beta_1 x_i$$

Priors:

$$\beta_0 \sim ?$$

$$[\beta_0] = ?$$

$$\beta_1 \sim ?$$

$$[\beta_1] = ?$$

$$\sigma^2 \sim ?$$

$$[\sigma^2] = ?$$

Normal data, continuous and real valued

DAG:

Posterior distribution:

$$\begin{aligned} [\beta_0, \beta_1, \sigma \mid \mathbf{y}] &\propto [\beta_0, \beta_1, \sigma, \mathbf{y}] \\ &\propto [\mathbf{y} \mid \beta_0, \beta_1, \sigma][\beta_0][\beta_1][\sigma] \\ &\propto \prod_{i=1}^n \text{Normal}(y_i \mid g(\beta_0, \beta_1, x_i), \sigma^2) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \\ &\quad \times \text{uniform}(\sigma \mid 0, 100) \\ g(\beta_0, \beta_1, x_i) &= \beta_0 + \beta_1 x_i \end{aligned}$$

Normal data, continuous and real valued

JAGS code for the model:

```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
sigma ~ dunif(0, 100)
tau <- 1/sigma^2
for (i in 1:length(y)){
  mu[i] <- b0 + b1 * x[i]
  y[i] ~ dnorm(mu[i], tau)
}
```

Interpretation:

- β_0 : expected outcome when $x = 0$
- β_1 : average change in the outcome for a one-unit change in x
- σ : std deviation of the outcomes about their respective means

Counts, discrete and non-negative

You have collected some count data ($y = 12, 17, 1, 0, 31, \dots, 25$) at n sites, along with a covariate x_i at each location which you believe is likely to affect these counts. Write a model regressing y on x .

- 1 Choose a specific stochastic and deterministic model.
- 2 Specify (vague) priors for your parameters.
- 3 Write out the DAG and express posterior distribution as proportional to joint distribution for your model.
- 4 Write the JAGS code for the model.
- 5 Interpret the coefficients of your model.

Poisson, discrete and non-negative

Stochastic model:

$$y_i | \beta_0, \beta_1 \stackrel{\text{ind.}}{\sim} \text{Poisson}(g(\beta_0, \beta_1, x_i)) \quad i = 1, \dots, n$$

Deterministic model:

$$\mu_i = g(\beta_0, \beta_1, x_i) = e^{\beta_0 + \beta_1 x_i}$$

Priors:

$$\beta_0 \sim ?$$

$$[\beta_0] = ?$$

$$\beta_1 \sim ?$$

$$[\beta_1] = ?$$

Poisson, discrete and non-negative

DAG:

Posterior distribution:

$$\begin{aligned} [\beta_0, \beta_1 \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{Poisson}(y_i \mid g(\beta_0, \beta_1, x_i)) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \\ g(\beta_0, \beta_1, x_i) &= e^{\beta_0 + \beta_1 x_i} \end{aligned}$$

Poisson, discrete and non-negative

JAGS code:

```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
for(i in 1:length(y)){
  log(mu[i]) <- b0 + b1 * x[i]
  y[i] ~ dpois(mu[i])
}
```

or

```
mu[i] <- exp(b0 + b1 * x[i])
y[i] ~ dpois(mu[i])
```

Poisson, discrete and non-negative

$$\begin{aligned} [\beta_0, \beta_1 \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{Poisson}(y_i \mid g(\beta_0, \beta_1, x_i)) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \\ \mu_i = g(\beta_0, \beta_1, x_i) &= e^{\beta_0 + \beta_1 x_i} = e^{\beta_0} e^{\beta_1 x_i} \end{aligned}$$

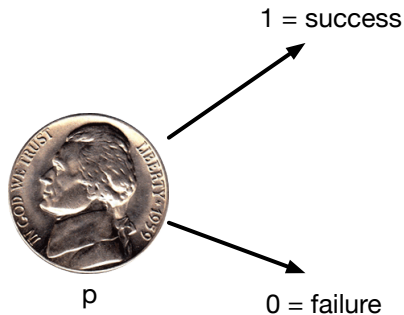
Interpretation: (Exponentiate coeff. and report multiplicative change in mean counts.)

- e^{β_0} : average count when $x = 0$
- e^{β_1} : multiplicative change in the mean count per one unit change in x

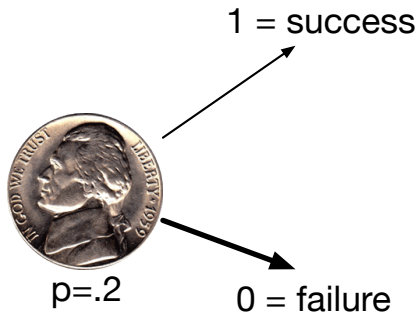
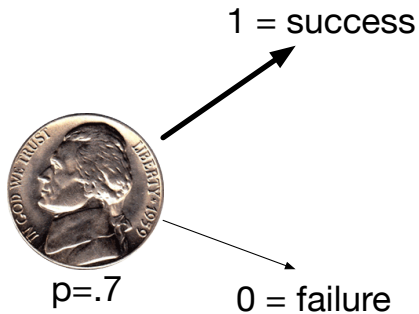
For example: “Mean western toad juvenile abundance is reduced by a factor of 5.1 (95% CI: 3.4, 10.8) per unit change in UV-B radiation.”

What was the estimate of β_1 ?

Binary data, 0 or 1 (aka logistic)



Binary data, 0 or 1 (aka logistic)



Binary data, 0 or 1 (aka logistic)

What happens when we want to relate p to a predictor x , where x can be any value on the real line? How do we connect x to $p \in [0, 1]$?

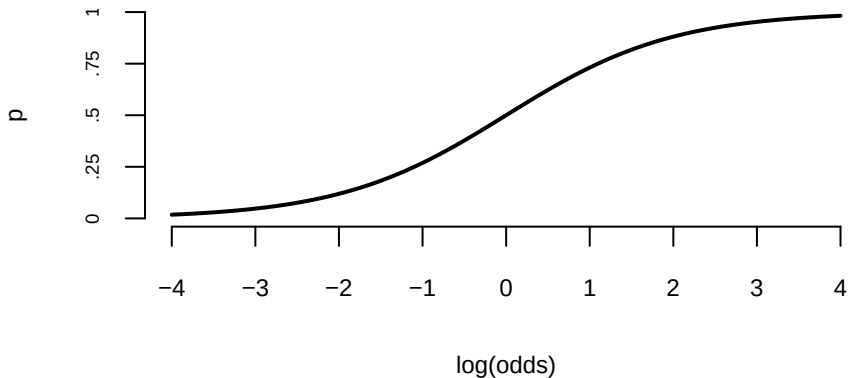
- odds: $\frac{p}{1-p} \in [0, \infty)$
- log odds: $\log(\text{odds}) = \log\left(\frac{p}{1-p}\right) \in (-\infty, \infty)$

Moving between probability and log odds and relating to x :

- $\text{logit}(p) = \log\left(\frac{p}{1-p}\right) = x$
 - ▶ input to $\text{logit}()$ is probability p , output is log odds x
- Inverting the above, we obtain $p = \frac{e^x}{e^x + 1} = \text{inverse logit}(x) = \text{expit}(x)$
 - ▶ input to inverse $\text{logit}()$ is $x = \text{log odds}$, output is probability p

Binary data, 0 or 1 (aka logistic)

Inverse logit mapping: input is log odds = $\log\left(\frac{p}{1-p}\right)$, output is probability



Binary data, 0 or 1 (aka logistic)

You have collected some binary data ($y = 1, 0, 0, 1, 1, 0, 1, \dots, 1$) at n sites, along with a covariate x_i measured at each location which you believe is likely to affect these counts. Write a model regressing y on x .

- 1 Choose a specific stochastic and deterministic model.
- 2 Specify (vague) priors for your parameters.
- 3 Write out the DAG and express posterior distribution as proportional to joint distribution for your model.
- 4 Write the JAGS code for the model.
- 5 Interpret the coefficients of your model.

Bernoulli, 0 or 1 (aka logistic)

Stochastic model:

$$y_i | \beta_0, \beta_1 \overset{\perp}{\text{ind.}} \text{Bernoulli}(g(\beta_0, \beta_1, x_i)) \quad i = 1, \dots, n$$

Deterministic model:

$$\mu_i = p_i = g(\beta_0, \beta_1, x_i) = \frac{e^{\beta_0 + \beta_1 x_i}}{e^{\beta_0 + \beta_1 x_i} + 1}$$

Priors:

$$\beta_0 \sim ?$$

$$[\beta_0] = ?$$

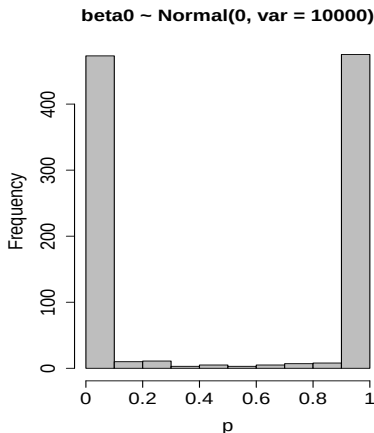
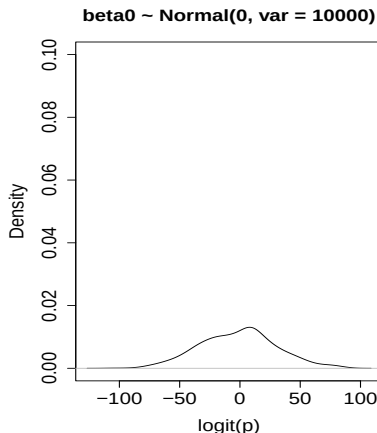
$$\beta_1 \sim ?$$

$$[\beta_1] = ?$$

Choosing reasonable flat priors on logit intercept

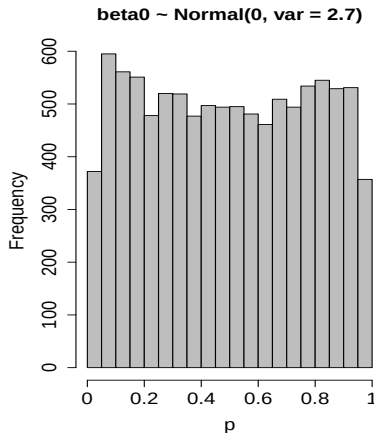
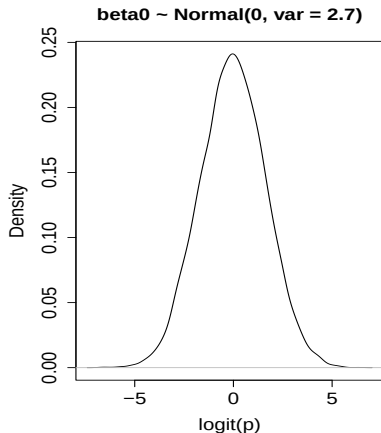
Imagine for now that we have no predictor, so $p_i = \mu_i = \frac{e^{\beta_0}}{e^{\beta_0} + 1}$.

If we use the same $\text{Normal}(0, \text{large variance})$ prior as before, what is the induced prior for the success probability p ?

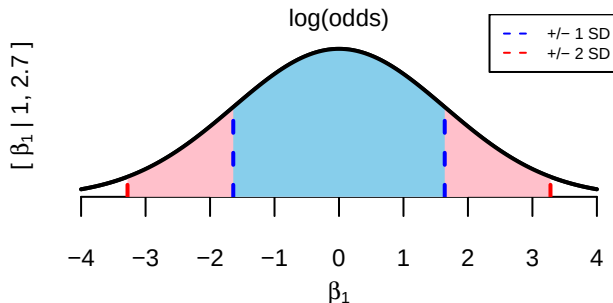
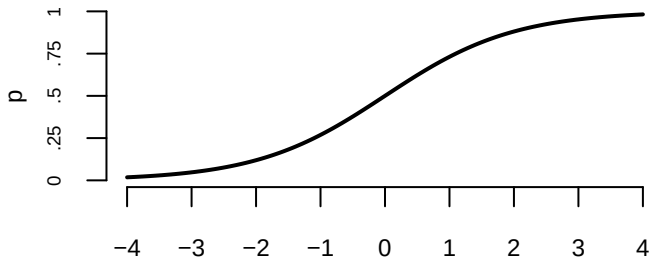


Choosing reasonable flat priors on logit intercept

Now instead consider the prior $\beta_0 \sim \text{Normal}(0, 2.7)$.



Choosing reasonable flat priors on logit effects



Bernoulli, 0 or 1 (aka logistic)

Returning to case with a single predictor, the posterior distribution is:

$$\begin{aligned} [\beta_0, \beta_1 \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{Bernoulli}(y_i \mid g(\beta_0, \beta_1, x_i)) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 2.7) \text{Normal}(\beta_1 \mid 0, 2.7) \\ g(\beta_0, \beta_1, x_i) &= \frac{e^{\beta_0 + \beta_1 x_i}}{e^{\beta_0 + \beta_1 x_i} + 1} \end{aligned}$$

Bernoulli, 0 or 1 (aka logistic)

JAGS code for the model:

```
b0 ~ dnorm(0, 1/2.7)
b1 ~ dnorm(0, 1/2.7)
for(i in 1:length(y)){
  logit(p[i]) <- b0 + b1 * x[i]
  y[i] ~ dbern(p[i])
}
```

or

```
p[i] <- inv.logit(b0 + b1 * x[i])
y[i] ~ dbern(p[i])
```

Bernoulli, 0 or 1 (aka logistic)

$$[\beta_0, \beta_1 \mid \mathbf{y}] \propto \prod_{i=1}^n \text{Bernoulli}(y_i \mid g(\beta_0, \beta_1, x_i)) \\ \times \text{Normal}(\beta_0 \mid 0, 2.7) \text{Normal}(\beta_1 \mid 0, 2.7)$$

$$p_i = g(\beta_0, \beta_1, x_i) = \frac{e^{\beta_0 + \beta_1 x_i}}{e^{\beta_0 + \beta_1 x_i} + 1} \iff \frac{p_i}{1 - p_i} = e^{\beta_0 + \beta_1 x_i} = e^{\beta_0} e^{\beta_1 x_i}$$

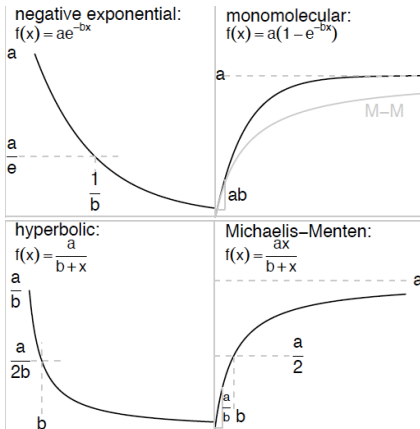
Interpretation: (Exponentiate coef. and report odds and odds ratios.)

- e^{β_0} : odds when $x = 0$
- e^{β_1} : multiplicative change in the odds for a one unit change in x

For example: “the odds of detecting weevils in upland willow stems were 3.2 (95% CI: 2.3, 4.8) times greater than detecting them in riparian willow stems.”

What was the estimate of β_1 ? (Might be helpful to identify/define x_i).

Nonlinear regression



{Figures c/o Bolker, B. 2008. *Ecological Models and Data in R*. Princeton University Press, Princeton, NJ. USA.}

Centering and standardizing

The remainder of the slides apply to all of the general linear models, but we will use a simple linear model for Normally distributed data as an example.

Centering predictor data

Rather than the usual linear model $y_i = \beta_0 + \beta_1 x_i + \varepsilon_i$ where $\varepsilon_i \stackrel{iid}{\sim} N(0, \sigma^2)$, consider the following linear model:

$$y_i = \beta_0 + \beta_1 (x_i - \bar{x}) + \varepsilon_i$$

where $\bar{x} = \sum_{i=1}^n x_i$ is the sample mean of the predictor.

Why complicate things?

- To reduce autocorrelation in MCMC chain and speed convergence.
- To make the intercept more easily interpretable.

Centering predictor data

$$\begin{aligned} [\beta_0, \beta_1, \sigma \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{Normal}(y_i \mid g(\beta_0, \beta_1, x_i), \sigma^2) \times \\ &\quad \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \times \\ &\quad \text{uniform}(\sigma \mid 0, 100) \\ g(\beta_0, \beta_1, x_i) &= \beta_0 + \beta_1(x_i - \bar{x}) \end{aligned}$$

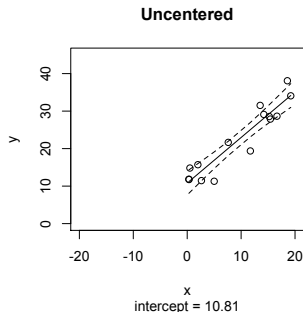
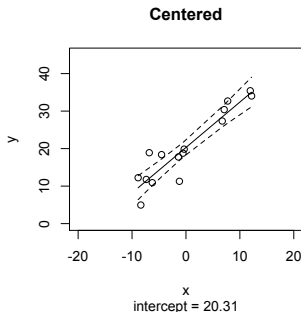
```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
sigma ~ dunif(0, 100)
tau <- 1/sigma^2
xBar <- mean(x)
for (i in 1:length(y)){
  mu[i] <- b0 + b1 * (x[i] - xBar)
  y[i] ~ dnorm(mu[i], tau)
}
b0_UC <- b0 - b1 * xBar
```


Exercise

What is the interpretation of β_0 in this model?

What is the interpretation of β_1 in this model?

Recovering uncentered parameters



$$\begin{aligned}\beta_0^* &= \beta_0 - \beta_1 \bar{x} \\ \beta_1^* &= \beta_1\end{aligned}$$

- For this to work properly, all the coefficients in the model must be *added*.
- Slopes will not be the same if there is an interaction term or quadratic. In these cases, back transforming is not simple.

Standardizing predictor data

$$y_i = \beta_0 + \beta_1 \left(\frac{x_i - \bar{x}}{s_x} \right)$$

where s_x is sample standard deviation of predictor

$$(s_x^2 = \frac{1}{n-1} \sum_{i=1}^n (x_i - \bar{x})^2)$$

Why complicate things?

- To reduce autocorrelation in MCMC chain and speed convergence.
- To make the intercept more easily interpretable.
- To make coefficients more easily comparable

Standardizing predictor data

$$\begin{aligned} [\beta_0, \beta_1, \sigma \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{Normal}(y_i \mid g(\beta_0, \beta_1, x_i), \sigma^2) \times \\ &\quad \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \times \\ &\quad \text{uniform}(\sigma \mid 0, 100) \\ g(\beta_0, \beta_1, x_i) &= \beta_0 + \beta_1 \left(\frac{x_i - \bar{x}}{s_x} \right) \end{aligned}$$

```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
sigma ~ dunif(0, 100)
tau <- 1/sigma^2
xBar <- mean(x)
xSD <- sd(x)
for (i in 1:length(y)){
  mu[i] <- b0 + b1 * ((x[i] - xBar)/xSD)
  y[i] ~ dnorm(mu[i], tau)
}
```

Recovering unstandardized parameters

$$\begin{aligned}y_i &= \beta_0 + \beta_1 \left(\frac{x_i - \bar{x}}{s_x} \right) \\&= \beta_0 + \frac{\beta_1}{s_x} x_i - \frac{\beta_1 \bar{x}}{s_x} \\&= \beta_0^* + \beta_1^* x_i \\ \beta_0^* &= \beta_0 - \frac{\beta_1 \bar{x}}{s_x} \\ \beta_1^* &= \frac{\beta_1}{s_x}\end{aligned}$$

- This only works if there are not squared values or interactions.
- It is fine to make predictions using $\hat{y}_i = \beta_0 + \beta_1 \frac{x_i - \bar{x}}{s_x}$ and plot $\hat{y}_i$ against x_i and the observed y_i

lognormal, data continuous and > 0 (log link)

$$\begin{aligned} [\beta_0, \beta_1, \sigma \mid \mathbf{y}] &\propto \prod_{i=1}^n \text{lognormal}(y_i \mid \log(g(\beta_0, \beta_1, x_i)), \sigma^2) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \\ &\quad \times \text{uniform}(\sigma \mid 0, 5) \\ g(\beta_0, \beta_1, x_i) &= e^{\beta_0 + \beta_1 x_i} \end{aligned}$$

Talk about the interpretation of σ .

```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
sigma ~ dunif(0, 5)
tau <- 1/sigma^2
for(i in 1:length(y)){
  mu[i] <- exp(b0 + b1 * x[i])
  y[i] ~ dlnorm(log(mu[i]), tau)
}
```

lognormal, data continuous and > 0 (not log link)

$$\begin{aligned} [\beta_0, \beta_1, \sigma \mid \mathbf{y}] &\propto \prod_{i=2}^n \text{lognormal}(y_i \mid \log(g(\beta_0, \beta_1, y_{i-1})), \sigma^2) \\ &\quad \times \text{Normal}(\beta_0 \mid 0, 1000) \text{Normal}(\beta_1 \mid 0, 1000) \\ &\quad \times \text{uniform}(\sigma \mid 0, 5) \text{uniform}(y_1 \mid 1, 1E6) \end{aligned}$$

$$g(\beta_0, \beta_1, y_{i-1}) = y_{i-1} e^{\beta_0 + \beta_1 y_{i-1}}$$

```
b0 ~ dnorm(0, .001)
b1 ~ dnorm(0, .001)
sigma ~ dunif(0, 5); tau <- 1/sigma^2
y[1] ~ dunif(1, 1E6)
for(i in 2:length(y)){
  mu[i] <- y[i-1] * exp(b0 + b1 * y[i-1])
  y[i] ~ dlnorm(log(mu[i]), tau)
}
```