# **ADVANCED** BAYESIAN MODELING

Logistic Regression Modeling

## Data Model

$$X_i = (x_{i1}, \dots, x_{ik}), \quad y_i \qquad \qquad i = 1, \dots, n$$

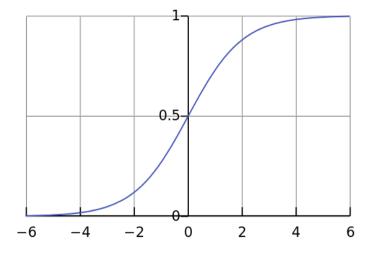
X has k explanatory variables (usually including an intercept)

$$y_i \mid \beta, X_i \quad \sim \quad \text{indep. } \mathrm{Bin}(n_i, p_i) \qquad \qquad n_i \text{ fixed and known}$$

$$logit(p_i) = X_i\beta$$
  $p_i = logit^{-1}(X_i\beta) = \frac{e^{X_i\beta}}{1 + e^{X_i\beta}}$ 

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The inverse logit:



From: Logistic regression. (2017, July 8). In *Wikipedia*, *The Free Encyclopedia*. Retrieved August 4, 2017 from https://en.wikipedia.org/w/index.php?title=Logistic\_regression&oldid=789649742

### **Priors**

In general logistic regression, conjugate priors are not very useful.

Flat priors are often a poor choice (later).

Consider instead proper priors from a standard (non-conjugate) distribution, but reasonably vague ...

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### Suggestion for weakly informative priors (see BDA3, Sec. 16.3):

- 1. Center/standardize explanatory variables:
  - ► Center indicator variables that don't define the intercept.
  - ► Center and scale quantitative variables to 0.5 sample standard deviation.
  - ► Use these centered/scaled variables to form any higher-order terms (e.g., interactions).
- 2. Choose independent scaled- $t_1$  priors for coefficients in  $\beta$ :
  - $t_1(0,2.5^2)$  for variables not defining the intercept
  - $t_1(0,10^2)$  for variables defining the intercept

The  $t_1(\mu, \sigma^2)$  distribution, also called the *Cauchy*, is a generalization of the t-distribution with one degree of freedom.

Like the normal, it is symmetric about  $\mu$  and unimodal, but has heavier tails, which may be useful if a few coefficients turn out to be large.

In JAGS:

dt(mu, sigmasqinv, 1)

Note: Parameterization uses *inverse* of  $\sigma^2$ .

Why not use flat priors?

Because, in certain cases, they would lead to improper posteriors.

Occurs when there is *separation*: A linear combination exists that predicts y perfectly (except possibly along some boundary).

Can occur even when the columns of X are linearly independent.

Details: BDA3, Sec. 16.3

# Overdispersion

Mean and variance of binomial are closely linked:

$$E(y_i \mid \beta, X_i) = n_i p_i \qquad var(y_i \mid \beta, X_i) = n_i p_i (1 - p_i)$$

Common problem: Data exhibit variances that are incompatible with the means – typically larger than they should be: **overdispersion**.

### Possible causes:

- Missing explanatory variables
- Dependent (underlying) Bernoulli variables
- ▶ Different probabilities in (underlying) Bernoulli distributions

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The **chi-square discrepancy** measures overdispersion:

$$T(y,\theta,X) = \sum_{i=1}^{n} \frac{(y_i - n_i p_i)^2}{n_i p_i (1-p_i)}$$

(When the model is correct and the  $n_i$ s are large enough, it has an approximate  $\chi_n^2$  distribution.)

When there is overdispersion, it tends to be larger.

Can be used to compute a posterior predictive p-value for overdispersion:

$$\Pr(T(y^{\mathsf{rep}}, \theta, X) \geq T(y, \theta, X) \mid y)$$

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One possible way to account for overdispersion:

Add a separate random effect to the linear combination of each observation.

$$logit(p_i) = X_i\beta + \epsilon_i \qquad \epsilon_i \sim iid N(0, \sigma_{\epsilon}^2)$$

Then add a reasonably noninformative prior on  $\sigma_{\epsilon}^2$  (e.g., flat on  $\sigma_{\epsilon}$ ).

Remark:

Overdispersion is generally a detectable problem only if the  $n_i$ s are not too small.

In particular, it is irrelevant to a purely binary response  $(n_i = 1)$ , in which case the model should be checked in other ways (e.g., trying additional explanatory variables).