Logistic Regression

Andrew Zieffler

Preparing the Data for Analysis

```
# Read in the data
grad <- read.csv(file = "http://www.tc.umn.edu/~zief0002/Data/grad.csv")

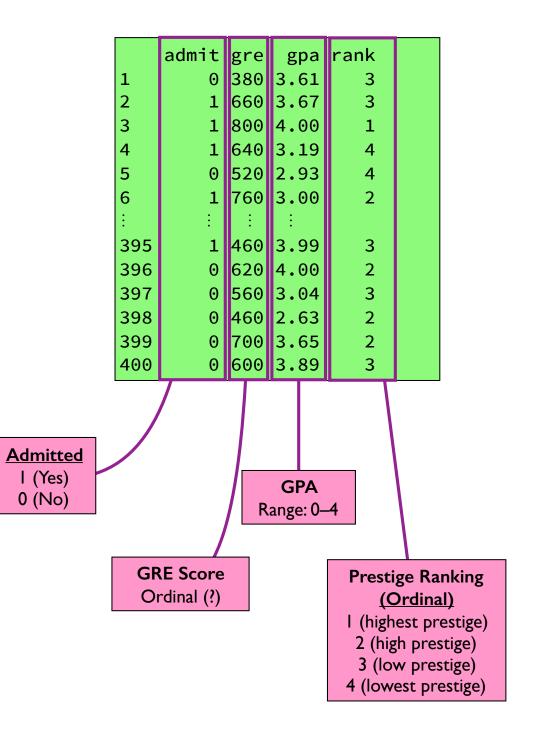
# Create new admission variable that is a factor
grad$admit2 <- as.factor(grad$admit)
levels(grad$admit2) <- c("Not Admitted", "Admitted")

# Create new rank variable that is a factor
grad$rank2 <- as.factor(grad$rank, ordered = TRUE)
levels(grad$rank2) <- c("Highest", "High", "Low", "Lowest")</pre>
```

Research Question

How do variables, such as Graduate Record Exam scores, grade point average, and prestige of the undergraduate institution, effect admission into graduate school?





> summary(grad)

admit Min. :0.0000 1st Qu.:0.0000 Median :0.0000 Mean : 0.3175

3rd Qu.:1.0000

Max. :1.0000 gre

Min. :220.0 1st Qu.:520.0

Median :580.0 :587.7 Mean

3rd Qu.:660.0

:800.0 Max.

gpa

Min. :2.260 1st Qu.:3.130

Min. :1.000 1st Qu.:2.000

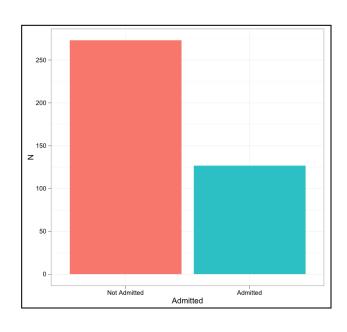
rank

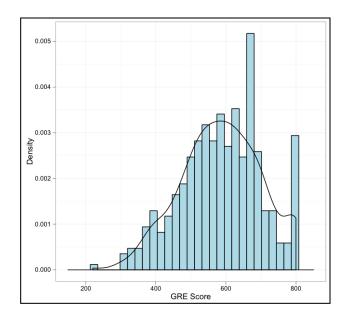
Median :3.395 Median :2.000

Mean :3.390 Mean :2.485

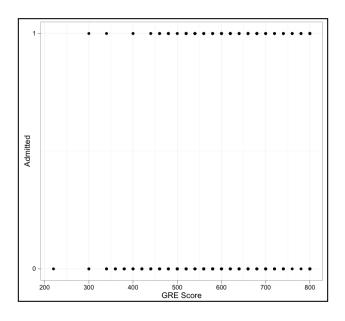
3rd Qu.:3.670 3rd Qu.:3.000

:4.000 Max. :4.000 Max.



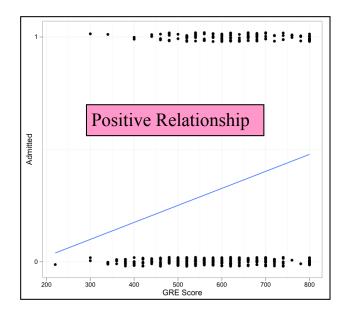


Examine relationship between GRE score and admission.



Problem: Hard to see relationships because of overplotting.

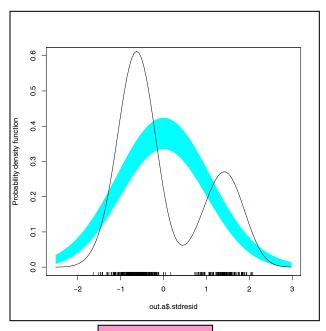
Solution: Add a small amount of random noise to the admit variable



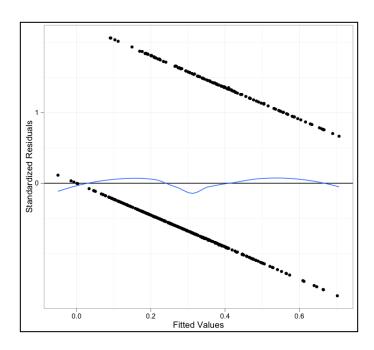
A correlation between a categorical, dichotomous variable and a quantitative variable is called a **point-biserial** correlation. It is a special case of Pearson's *r* when one of the variables is dummy coded.

• Students with higher GRE scores are more likely, on average, to be admitted to graduate school.

A correlation between a categorical, dichotomous variable and a ordinal variable is called a **rank-biserial** correlation. This can be obtained by adding the argument method="spearman" in the cor() function. (In this case they are not that different.)



The errors are not normally distributed.



Two parallel lines

$$\epsilon_i = 1 - \pi_i$$

$$\epsilon_i = -\pi_i$$

Linear model fitted to dichotomous outcomes is called *linear-probability model*

$$E(Y_i) = \beta_0 + \beta_1(X_i) + \epsilon_i$$

Let
$$\pi_i \equiv P(Y = 1 | X = x_i)$$
 then $E(Y_i) = \pi_i(1) + (1 - \pi_i)(0) = \pi_i$

$$\pi_i = eta_0 + eta_1(X_i) + \epsilon_i$$
 Re-expression of the model

$$\epsilon_i = Y_i - \hat{Y} = Y_i - \pi_i$$

$$\underline{\text{If } Y_i = 1} \qquad \epsilon_i = 1 - \pi_i$$

$$\begin{array}{c|c}
\hline
\mathbf{If Y_i = 0} \\
\end{array} \qquad \epsilon_i = -\pi_i$$

Errors from the linear-probability model are dichotomous...cannot be normally distributed

May (or may not) be a problem depending on sample size...remember the C.L.T.

$$Var(\epsilon_i) = E(\epsilon_i^2) - (E(\epsilon_i))^2$$

$$E(\epsilon_i) = 0$$
 Assumption of the model

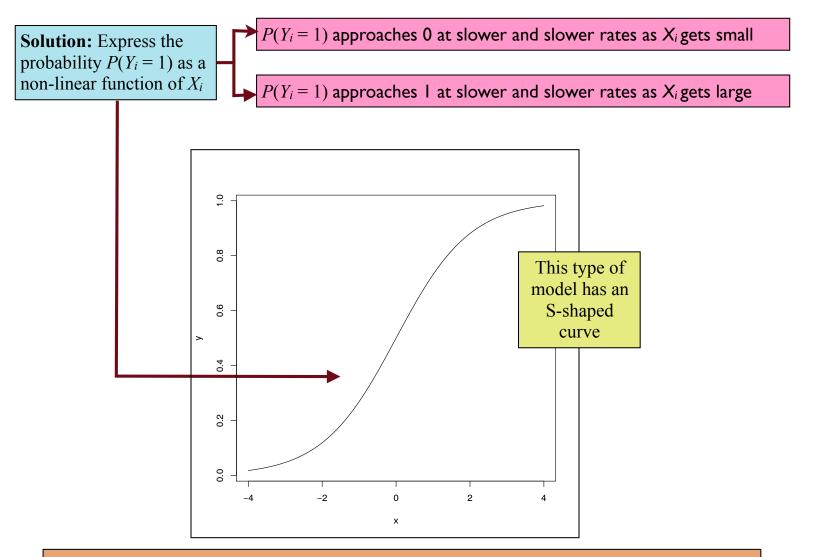
$$Var(\epsilon_i) = E(\epsilon_i^2)$$

$$= \pi_i (\epsilon_{i|Y_i=1})^2 + (1 - \pi_i) (\epsilon_{i|Y_i=0})^2$$

$$= \pi_i (1 - \pi_i)^2 + (1 - \pi_i) (-\pi_i)^2$$

Unless π_i is the same across all x_i , the errors will not have the same variance at each x_i

Furthermore, there is no guarantee that the predicted outcomes will be constrained to the [0, 1] range...this is a problem with linearity (constant rate of change)



Several key features

- The probability of Y_i never go lower than 0 or above 1
- A person with an *X*-value of 0 has less probability of *Y* (e.g., being admitted) than a person with an *X*-value of 1 (monotonic)
- For really small (or really large) values of *X*, the probability of *Y* does not change much (definitely not a constant rate of change)

Transformation of the Probabilities Probit Transformation

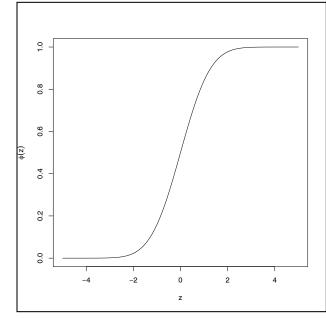
To map the probability P(Y) to the [0, 1] space, we apply a transformation

$$\pi_i = \phi(\beta_0 + \beta_1 X_i)$$

The transformation used is any function that can fit the criteria we had before (monotonic, nonlinear, maps to [0, 1] space)

Common nonlinear model that meets these specifications is the cumulative density function (CDF)—although any CDF would work, here we use the unit normal distribution's CDF

$$\phi(z) = \frac{1}{2\pi} \int_{-\inf}^{z} exp\left(-\frac{1}{2}z^{2}\right) dz$$



$$\phi(\beta_0 + \beta_1 X_i) = \frac{1}{2\pi} \int_{-\inf}^{\beta_0 + \beta_1 X_i} exp\left(-\frac{1}{2}z^2\right) dz$$

This is called a **probit transformation**, and the model is the **linear probit model**.

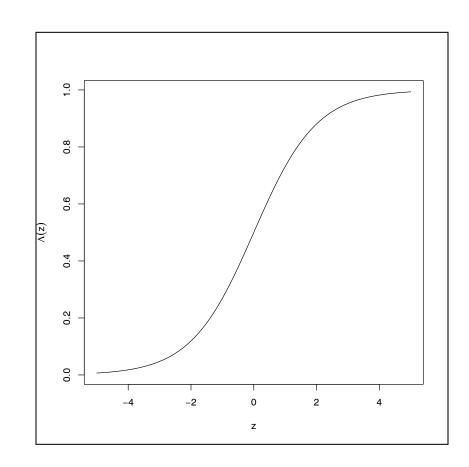
Logistic Transformation

Another common nonlinear model that meets these specifications is the **logistic** function

$$\Lambda(z) = \frac{1}{1 + e^{-z}}$$

We map the probability onto the logistic function

$$\pi_i = \Lambda(\beta_0 + \beta_1 X_i)$$



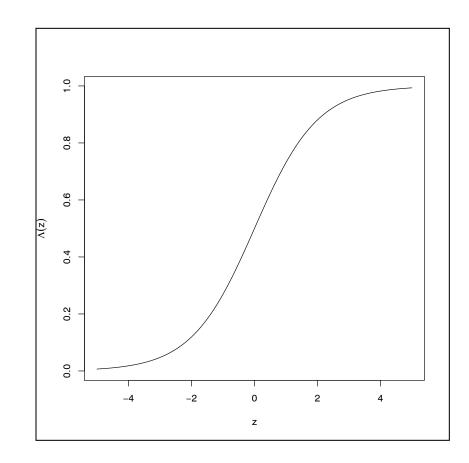
$$\Lambda(\beta_o + \beta_1 X_i) = \frac{1}{1 + e^{-(\beta_o + \beta_1 X_i)}}$$

This is called a **logistic transformation**, and the model is the **linear logistic model**.

In practice (once the variances are equalized) there is no difference between the probit and logit transformations

Two advantages to using the logit transformation

- I. Equation for logit transformation is **simpler** than for probit transformation
- 2. When we back transform, the inverse of the logit transformation is directly **interpretable**. (The inverse of the probit transformation is not interpretable directly.)



Re-expressing the linear logistic regression

$$\pi_i = \frac{1}{1 + e^{-(\beta_o + \beta_1 X_i)}}$$

► Taking the natural logarithm of both sides of the equation

$$\frac{\pi_i}{1 - \pi_i} = \exp(\beta_o + \beta_1 X_i)$$

$$ln\left(\frac{\pi_1}{1-\pi_i}\right) = \beta_o + \beta_1 X_i$$

This ratio,
$$\frac{\pi_i}{1-\pi_i}$$

The inverse transformation of the probability is called the **logit**.

is called the **odds**. Odds refers to the relative chances of an occurrence, Y_i

Logits

Logits are log-odds. It is the log transformed odds that $Y_i = 1$ rather than 0

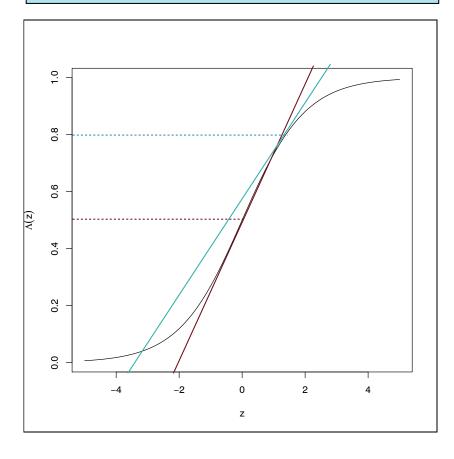
Another way to think about logistic regression, is to fit a linear model on the log-odds (logit) of the outcome variable *Y*

$$\frac{\pi_1}{1 - \pi_i} = exp(\beta_o + \beta_1 X_i)$$
$$= exp(\beta_o) \cdot exp(\beta_1 X_i)$$
$$= exp(\beta_o) \cdot exp(\beta_1)^{X_i}$$

Increasing X_i by one unit changes the logit, on average by β_1 and changes the odds by a factor of e^{β_1}

	Logit $ln\left(\frac{\pi_i}{-}\right)$	Odds π_i	Probability
	$ln\left(\frac{\pi_i}{1-\pi_i}\right)$	$\frac{1}{1-\pi_i}$	π_i
+1	-4	0.01831564	0.01798621
	-3	0.04978707	0.04742587
	÷	÷	÷
	0	I	0.5
+1	, 4	2.718282	0.7310586
	:	:	:
	3	20.08554	0.9525741
+1	4	54.59815	0.9820138

Consider the slope of the tangent line of the logistic function at a single probability (Y-value). This will help us understand the relationship between π_i and X_i



$$\frac{\partial \pi_i}{\partial \beta_1} = \beta_1 \cdot \pi_i \cdot (1 - \pi_i)$$

The tangent line at $\pi_i = 0.8$ is less steep (lower linear rate of change) than the tangent line at $\pi_i = 0.5$.

- This suggests the rate of change in probability of *Y* is different depending on the level of *X*
- The slope of these tangent lines can be computed by determining the partial derivative of the logistic function

π_i	$\beta_1 \pi_i (1 - \pi_i)$
0.01	$\beta_1 \times 0.0099$
0.05	$\beta_1 \times 0.0475$
0.1	$\beta_1 \times 0.09$
0.2	$\beta_1 \times 0.16$
0.5	$\beta_1 \times 0.25$
0.8	$\beta_1 \times 0.16$
0.9	$\beta_1 \times 0.09$
0.99	$\beta_1 \times 0.0099$

To fit the logistic model, we use the glm() function.

- The family= argument specifies the error distribution. For binary outcomes this is a **binomial distribution**.
- Within the binomial() function, we then specify the specific transformation via the link= argument.

Increasing X_i by one unit changes the logit, on average by β_1 and changes the odds by a factor of e^{β_1}

```
> glm.a <- glm(admit ~ gre, data = grad, family = binomial(link = "logit"))
> summary(glm.a)
Call:
glm(formula = admit ~ gre, family = binomial(link = "logit"),
   data = grad)
```

Deviance Residuals:

```
Min
           10 Median
                          30
                                 Max
-1.1623 -0.9052 -0.7547 1.3486 1.9879
```

Coefficients:

```
Estimate Std. Error z value Pr(>|z|)
(Intercept) -2.901344  0.606038  -4.787  1.69e-06 ***
gre
             0.003582
                       0.000986
                                   3.633 0.00028 ***
```

Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1

(Dispersion parameter for binomial family taken to be 1)

Null deviance: 499.98 on 399 degrees of freedom Residual deviance: 486.06 on 398 degrees of freedom AIC: 490.06

Number of Fisher Scoring iterations: 4

Parameter estimates from the logistic regression (logits/log-odds)

$$ln\left(\frac{\pi_1}{1-\pi_i}\right) = \beta_o + \beta_1 X_i$$

- The intercept, -2.90, is the predicted log-odds for being admitted to graduate school when a applicant has a GRE score of 0
- The slope, 0.004, is the predicted change in the log-odds in admission for a one-point change in GRE score.

Parameter estimates from the logistic regression (odds)

$$\frac{\pi_i}{1 - \pi_i} = exp(\beta_o) \cdot exp(\beta_1)^{X_i}$$

- The intercept, $e^{-2.90} = 0.055$, is the predicted odds for being admitted to graduate school when a applicant has a GRE score of 0.
 - ✓ Recall odds values below 1 correspond to odds with higher numbers in the denominator (higher probability of not being admitted) than in the numerator
 - ✓ The reciprocal, 1/0.055 = 18.199, gives the odds of not being admitted for an applicant with a GRE score of 0.
- The slope, $e^{0.004} = 1.004$, is the predicted change in the odds in admission for a one-point change in GRE score.
 - ✓ A one-point change in GRE score is associated with an increase in the odds of being admitted to graduate school by a factor 1.004

Parameter estimates from the logistic regression (probabilities)

- The intercept, $e^{-2.90}/(1 + e^{-2.90}) = 0.052$, is the predicted probability of being admitted to graduate school when a applicant has a GRE score of 0.
 - ✓ The probability of not being admitted for the same GRE score is 0.948.
- The slope is not directly interpretable in terms of the predicted change in probability
 - ✓ Recall the change in probability is a function of the probability (non-linear change), namely $\beta_1\pi_i(1-\pi_i)$
 - ✓ Thus, the predicted change in probability (effect of GRE) is dependent on where the applicant is on the GRE spectrum

Re-expressing the Joint Probability

The joint probability, assuming independence can be expressed as

$$p(y_1, y_2, \dots, y_n) = \prod_{i=1}^{n} p(y_i)$$

Substituting the probability for a binary variable we get,

$$X^{1-c} = \frac{X^1}{X^c}$$

Using this, we can re-write the product as...

$$= \prod_{i=1}^{n} \pi_i^{y_i} (1 - \pi_i)^{1 - y_i}$$

$$= \prod_{i=1}^{n} \pi_i^{y_i} \frac{(1-\pi_i)^1}{(1-\pi_i)^{y_i}}$$

$$= \prod_{i=1}^{n} \frac{\pi_i^{y_i}}{(1-\pi_i)^{y_i}} (1-\pi_i)$$
Re-organize
$$= \prod_{i=1}^{n} \left[\frac{\pi_i}{(1-\pi_i)} \right]^{y_i} (1-\pi_i)$$
Simplify

From the model, recall

$$\frac{\pi_i}{(1-\pi_i)} = e^{\beta_0 + \beta_1(X_{i1})}$$

We can use this equation and solve for both π_i and $(1 - \pi_i)$

 $\pi_i = e^{\beta_0 + \beta_1(X_{i1})} (1 - \pi_i)$ $= e^{\beta_0 + \beta_1(X_{i1})} - \pi_i e^{\beta_0 + \beta_1(X_{i1})}$ $\pi_i + \pi_i e^{\beta_0 + \beta_1(X_{i1})} = e^{\beta_0 + \beta_1(X_{i1})}$ $\pi_i \left(1 + e^{\beta_0 + \beta_1(X_{i1})} \right) = e^{\beta_0 + \beta_1(X_{i1})}$ $\pi_i = \frac{e^{\beta_0 + \beta_1(X_{i1})}}{1 + e^{\beta_0 + \beta_1(X_{i1})}}$ Solving for π_i $1 - \pi_i = \frac{1}{e^{\beta_0 + \beta_1(X_{i1})}}$ Solving for $(1 - \pi_i)$ $1 - \pi_i = \frac{e^{\beta_0 + \beta_1(X_i)}}{1 + e^{\beta_0 + \beta_1(X_{i1})}} \frac{1}{e^{\beta_0 + \beta_1(X_{i1})}}$ $= \frac{1}{1 + e^{\beta_0 + \beta_1(X_{i1})}}$ Substitute

Simplify

We now substitute the value of π_i into this equation

Maximum Likelihood Estimation for the Logistic Regression Model

Thus the joint probability can be expressed as...

$$p(y_1, y_2, \dots, y_n) = \prod_{i=1}^n \left[e^{\beta_0 + \beta_1(X_{i1})} \right]^{y_i} \left[\frac{1}{1 + e^{\beta_0 + \beta_1(X_i)}} \right]$$

If we treat the observed binary response vector

$$y_i = (y_1, y_2, ..., y_n)$$

as fixed (after all it is what we observed), then this becomes a function of the parameters β_0 and β_1 referred to as the **likelihood function** which is written as...

$$\mathcal{L}(\beta_0, \beta_1, \dots, \beta_k) = \prod_{i=1}^n \left[e^{\beta_0 + \beta_1(X_{i1}) + \dots + \beta_k(X_{ik})} \right]^{y_i} \left[\frac{1}{1 + e^{\beta_0 + \beta_1(X_{i1} + \dots + \beta_k(X_{ik}))}} \right]$$

Goal: Find the values

$$\hat{\beta}_0, \hat{\beta}_1, \dots, \hat{\beta}_k$$

that maximize the likelihood.

Inference for the Logistic Regression Model Model-Level Inference

(Dispersion parameter for binomial family taken to be 1)

Null deviance: 499.98 on 399 degrees of freedom

Residual deviance: 486.06 on 398 degrees of freedom

AIC: 490.06

Number of Fisher Scoring iterations: 4

The residual deviance associated with the intercept only logistic model

The residual deviance associated with the fitted logistic model

Deviance

-2(LL_{Model})

H₀: The reduced model and the full model fit equally well.

Model fit is measured through the deviance (smaller = better fit). If the two models fit equally well, there will be no difference in their deviances.

$$-2\ln(Lik_{\rm Reduced}) - (-2\ln(Lik_{\rm Full})) = -2\left[\ln(Lik_{\rm Reduced}) - \ln(Lik_{\rm Full})\right]$$

Difference in deviances (Likelihood ratio)

$$= -2 \ln \left(\frac{Lik_{\text{Reduced}}}{Lik_{\text{Full}}} \right)$$

499.98 - 486.06 = 13.92 Sample evidence against H_0

If *n* is sufficiently large, the likelihood ratio is $\sim \chi^2 (df_{\text{Full}} - df_{\text{Reduced}})$

Inference for the Logistic Regression Model

Parameter-Level Inference

Coefficients:

Null hypothesis associated with each parameter

$$H_0: \beta_j = 0$$

To examine the degree of statistical evidence against the null hypothesis, we carry out a **Wald test**.

$$Z = rac{\hat{eta}_j}{SE_{\hat{eta}_j}}$$
 $Z_j \sim \mathcal{N}(0, 1)$

$$3.633^2 = 13.19 \approx 13.92$$

When there is only one parameter that is different between a full and the reduced model, for large n the likelihood ratio (LR) = Z^2 _{Wald}

Confidence intervals can also be computed based on the Wald statistic

$$\hat{\beta}_j \pm 2 \times SE_{\hat{\beta}_j}$$

 $0.003582 \pm 2 \times 0.000986$

[0.00161, 0.005554]

Confidence limits for the effect on the logit

Most statisticians prefer to give confidence limits for the odds ratio rather than the logit.

$$[e^{0.00161}, e^{0.005554}]$$

[1.001611, 1.005569]

Confidence limits for the effect on the odds ratio

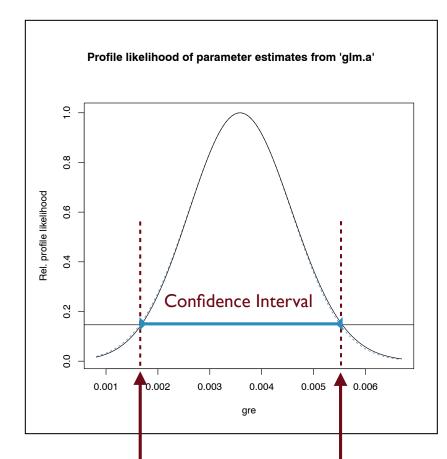
Better Confidence Intervals Profile Likelihood Cls

Confidence intervals based on the Wald statistic are formed by **inverting** the Wald tests

- Inversion determines the parameter values for which the null hypothesis is not rejected
- For small sample sizes the Wald test can be terribly wrong, which means the CI based on the inversion will be wrong as well

Better CIs can be produced by inverting the LR test. This is done by **profiling the likelihood**.

- A set of potential parameter values are chosen
- For this set of values, the likelihood is maximized over the other parameters included in the model
- The result is plotted in a profile plot



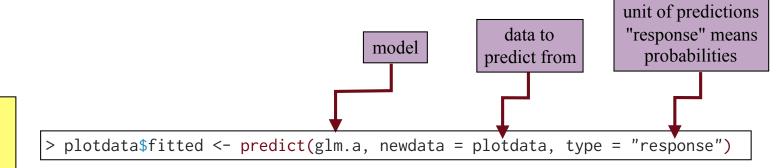
```
> confint(glm.a)
Waiting for profiling to be done...
2.5 % 97.5 %
(Intercept) -4.119988259 -1.739756286
gre 0.001679963 0.005552748
```

Plotting the Fitted Model

Step 1: Create a data frame of the predictors in the model.

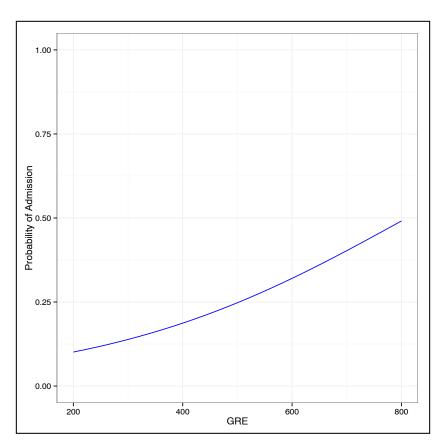
```
> plotdata <- data.frame(
    gre = seq(from = 200, to = 800, by = 20)
)</pre>
```

Step 2: Use the predict() function to produce a fitted value for each row in the newly created data frame.



Step 3: Plot the fitted values versus the *X*s

```
> ggplot(data = plotdata, aes(x = gre, y = fitted)) +
     geom_line(color = "blue") +
     xlab("GRE") +
     ylab("Probability of Admission") +
     ylim(c(0, 1)) +
     theme_bw()
```



Shortcut when there is only one predictor

```
> ggplot(data = grad, aes(x = gre, y = admit)) +
    geom_smooth(method = "glm", family = "binomial) +
    xlab("GRE") +
    ylab("Probability of Admission") +
    ylim(c(0, 1)) +
    theme_bw()
```

Adding a Confidence Envelope to the Model

Use the se.fit = TRUE argument in the predict() function

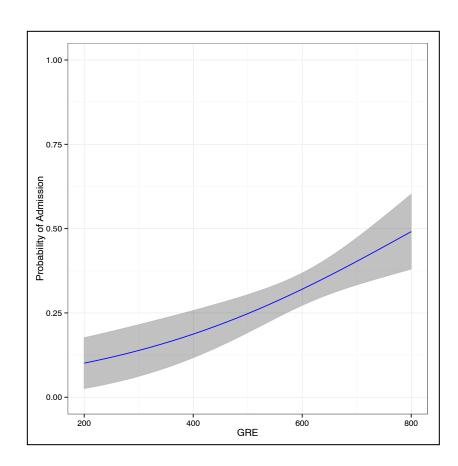
```
> fitted <- predict(glm.a, newdata = plotdata,
    type = "response", se.fit = TRUE)</pre>
```

Create a data frame that includes the predictor, the fitted values and the SEs

```
> plotdata2 <- data.frame(
    gre = plotdata$gre,
    fitted = fitted$fit,
    se = fitted$se.fit
)</pre>
```

Compute the lower and upper bounds for the confidence envelope

```
> plotdata2$lowerLimit <- plotdata2$fitted - 2 * plotdata2$se
> plotdata2$upperLimit <- plotdata2$fitted + 2 * plotdata2$se</pre>
```



Use geom_ribbon() to draw the confidence envelope

```
> ggplot(data = plotdata2, aes(x = gre, y = fitted)) +
    geom_ribbon(aes(ymin = lowerLimit, ymax = upperLimit),
        color = "grey", alpha = 0.3) +
    geom_line(color = "blue") +
    xlab("GRE") +
    ylab("Probability of Admission") +
    ylim(c(0, 1)) +
    theme_bw()
```