

# MA678 Homework 4

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## 13.5 Interpreting logistic regression coefficients

Here is a fitted model from the Bangladesh analysis predicting whether a person with high-arsenic drinking water will switch wells, given the arsenic level in their existing well and the distance to the nearest safe well:

```
stan_glm(formula = switch ~ dist100 + arsenic, family=binomial(link="logit"), data=wells)
              Median MAD_SD
(Intercept)    0.00    0.08
dist100        -0.90    0.10
arsenic         0.46    0.04
```

Compare two people who live the same distance from the nearest well but whose arsenic levels differ, with one person having an arsenic level of 0.5 and the other person having a level of 1.0. You will estimate how much more likely this second person is to switch wells. Give an approximate estimate, standard error, 50% interval, and 95% interval, using two different methods:

(a)

Use the divide-by-4 rule, based on the information from this regression output.

```
# The approximate estimate is 0.46/4=11.5%
# The standard Error is 0.04, as shown
# The 50% interval is 0.115 +- 0.67*0.04 =[0.0882,0.142]
# The 95% interval is 0.115 +- 1.96*0.04 =[0.0366,0.193]
```

(b)

Use predictive simulation from the fitted model in R, under the assumption that these two people each live 50 meters from the nearest safe well.

```
library(rstanarm)

## Loading required package: Rcpp
## This is rstanarm version 2.21.3
## - See https://mc-stan.org/rstanarm/articles/priors for changes to default priors!
## - Default priors may change, so it's safest to specify priors, even if equivalent to the defaults.
## - For execution on a local, multicore CPU with excess RAM we recommend calling
##   options(mc.cores = parallel::detectCores())

wells <- read.csv("/Users/billg/Desktop/MA-678-Homework/MA678-HW4/wells.csv")
Reg13.5 <- stan_glm(formula = switch ~ dist100 + arsenic, family=binomial(link="logit"), data=wells, ref.
predic13.5 <- posterior_epred(Reg13.5,data=wells)
mean(predic13.5)
```

```
## [1] 0.5751749
```

```
sd(predic13.5)
```

```
## [1] 0.1203117
```

### 13.7 Graphing a fitted logistic regression

We downloaded data with weight (in pounds) and age (in years) from a random sample of American adults. We then defined a new variable:

```
heavy <- weight > 200
```

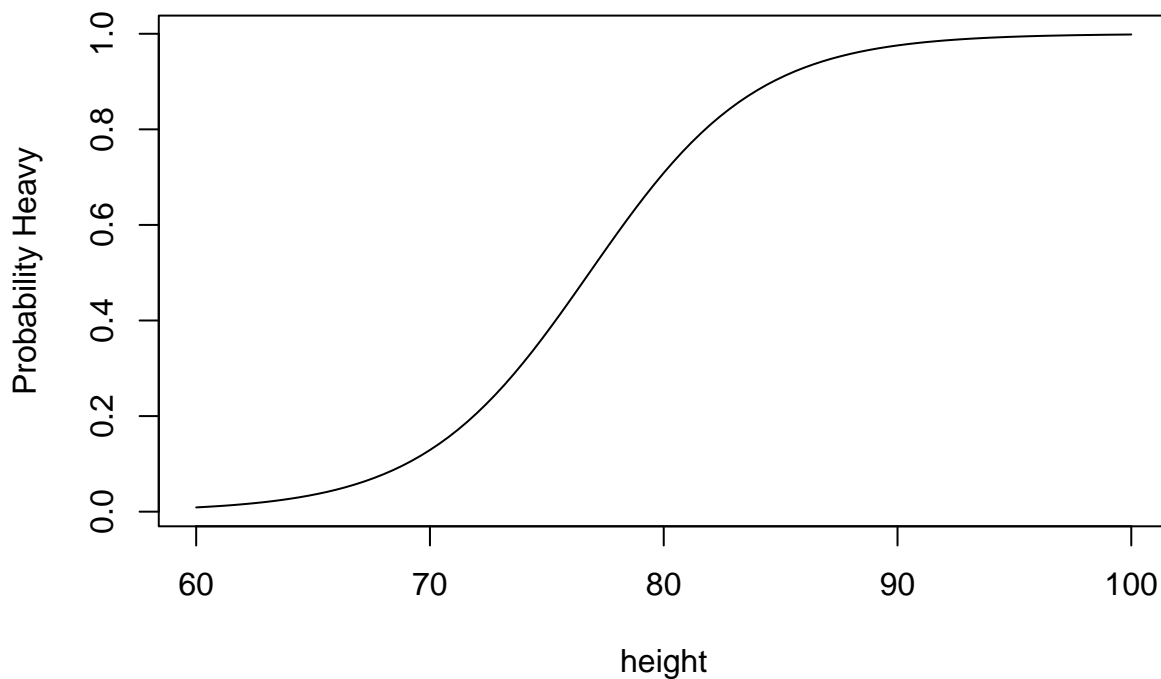
and fit a logistic regression, predicting heavy from height (in inches):

```
stan_glm(formula = heavy ~ height, family=binomial(link="logit"), data=health)
              Median MAD_SD
(Intercept)  -21.51    1.60
height         0.28    0.02
```

(a)

Graph the logistic regression curve (the probability that someone is heavy) over the approximate range of the data. Be clear where the line goes through the 50% probability point.

```
curve(invlogit(-21.51+0.28*x),xlab="height",ylab="Probability Heavy",xlim=c(60,100))
```



*# The line goes through the 50% point when  $-21.51+0.28*x=0$ . Thus  $x=79$*

(b)

Fill in the blank: near the 50% point, comparing two people who differ by one inch in height, you'll expect a difference of \_\_\_\_\_ in the probability of being heavy.

*# By using the divide by 4 rule, we will expect a difference of  $0.28/4 = 7\%$   
# in the probability of being heavy.*

## 13.8 Linear transformations

In the regression from the previous exercise, suppose you replaced height in inches by height in centimeters. What would then be the intercept and slope?

*# 1 inch is 2.54 centimeters. Therefore, after we replace it with height in centimeters,  
# the intercept is gonna be the same and the slope is gonna be 1/2.54 times the original slope*

## 13.10 Expressing a comparison of proportions as a logistic regression

A randomized experiment is performed within a survey, and 1000 people are contacted. Half the people contacted are promised a \$5 incentive to participate, and half are not promised an incentive. The result is a 50% response rate among the treated group and 40% response rate among the control group.

(a)

Set up these results as data in R. From these data, fit a logistic regression of response on the treatment indicator.

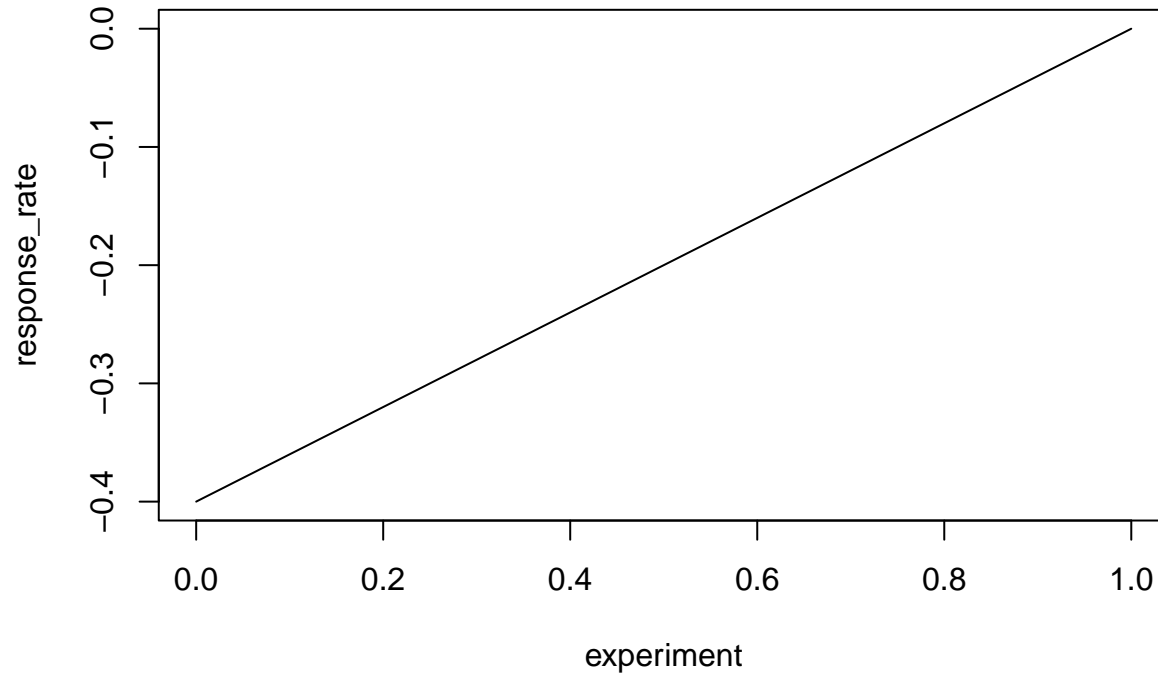
```
set.seed(114514)
experiment <- c(rep(1,500),rep(0,500))
response_rate <- c(rep(0,250),rep(1,250),rep(0,300),rep(1,200))
Reg13.10 <- stan_glm(response_rate~experiment,family =binomial(link=logit),refresh=0)
```

```
## Warning: Omitting the 'data' argument is not recommended and may not be allowed
## in future versions of rstanarm. Some post-estimation functions (in particular
## 'update', 'loo', 'kfold') are not guaranteed to work properly unless 'data' is
## specified as a data frame.
```

```
summary(Reg13.10)
```

```
##
## Model Info:
## function:      stan_glm
## family:        binomial [logit]
## formula:       response_rate ~ experiment
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  1000
## predictors:    2
##
## Estimates:
##           mean   sd  10%   50%   90%
## (Intercept) -0.4   0.1 -0.5  -0.4  -0.3
## experiment   0.4   0.1  0.2   0.4   0.6
##
## Fit Diagnostics:
##           mean   sd  10%   50%   90%
## mean_PPD 0.5    0.0  0.4   0.4   0.5
##
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
##
## MCMC diagnostics
##           mcse Rhat n_eff
## (Intercept) 0.0  1.0  2432
## experiment  0.0  1.0  2507
```

```
## mean_PPD      0.0  1.0  3194
## log-posterior 0.0  1.0  1674
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
curve(-0.4+0.4*x,xlab="experiment",ylab="response_rate")
```



(b)

Compare to the results from Exercise 4.1.

```
set.seed(114514)
experiment <- c(rep(1,500),rep(0,500))
response_rate <- c(rep(0,250),rep(1,250),rep(0,300),rep(1,200))
Reg13.10b <- lm(response_rate~experiment,family =binomial(link=logit),refresh=0)
```

```
## Warning: In lm.fit(x, y, offset = offset, singular.ok = singular.ok, ...) :
## extra arguments 'family', 'refresh' will be disregarded
```

```
summary(Reg13.10b)
```

```
##
## Call:
## lm(formula = response_rate ~ experiment, family = binomial(link = logit),
##     refresh = 0)
##
## Residuals:
```

	Min	1Q	Median	3Q	Max
	-0.500	-0.425	-0.400	0.500	0.600

```
##
## Coefficients:
```

	Estimate	Std. Error	t value	Pr(> t )
(Intercept)	0.40000	0.02216	18.052	< 2e-16 ***
experiment	0.10000	0.03134	3.191	0.00146 **

```
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 0.4955 on 998 degrees of freedom
## Multiple R-squared:  0.0101, Adjusted R-squared:  0.009109
## F-statistic: 10.18 on 1 and 998 DF,  p-value: 0.001461
```

*# The results are somewhat consistent*

## 13.11 Building a logistic regression model

The folder `Rodents` contains data on rodents in a sample of New York City apartments.

(a)

Build a logistic regression model to predict the presence of rodents (the variable `rodent2` in the dataset) given indicators for the ethnic groups (`race`). Combine categories as appropriate. Discuss the estimated coefficients in the model.

```
rodents <- read.table("/Users/billg/Desktop/MA-678-Homework/MA678-HW4/rodents.csv")
Reg13.11 <- stan_glm(rodent2~race,data=rodents,family =binomial(link =logit),refresh=0)
summary(Reg13.11)
```

```
##
## Model Info:
## function:      stan_glm
## family:        binomial [logit]
## formula:       rodent2 ~ race
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  1551
## predictors:    2
##
## Estimates:
##           mean   sd  10%   50%   90%
## (Intercept) -1.9   0.1 -2.1  -1.9  -1.8
## race         0.3   0.0  0.3   0.3   0.4
##
## Fit Diagnostics:
##           mean   sd  10%   50%   90%
## mean_PPD 0.2    0.0  0.2   0.2   0.3
##
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
##
## MCMC diagnostics
##           mcse Rhat n_eff
## (Intercept)  0.0  1.0  2419
## race         0.0  1.0  2642
## mean_PPD     0.0  1.0  2949
## log-posterior 0.0  1.0  1742
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
```

*# The estimate intercept coefficient is -1.9 meaning the average is -1.9 when race=0*

(b)

Add to your model some other potentially relevant predictors describing the apartment, building, and community district. Build your model using the general principles explained in Section 12.6. Discuss the coefficients for the ethnicity indicators in your model.

```
Reg13.11b <- stan_glm(rodent2~race+personrm+housewgt+sequenceno,data=rodents,family=binomial(link=logit))
summary(Reg13.11b)
```

```
##
## Model Info:
## function:      stan_glm
## family:        binomial [logit]
## formula:       rodent2 ~ race + personrm + housewgt + sequenceno
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  1551
## predictors:    5
##
## Estimates:
##              mean    sd   10%   50%   90%
## (Intercept) -1.3    0.3  -1.7  -1.3  -0.9
## race         0.2     0.0   0.2   0.2   0.3
## personrm     0.8     0.1   0.7   0.8   1.0
## housewgt     0.0     0.0   0.0   0.0   0.0
## sequenceno   0.0     0.0   0.0   0.0   0.0
##
## Fit Diagnostics:
##              mean    sd   10%   50%   90%
## mean_PPD 0.2     0.0   0.2   0.2   0.3
##
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
##
## MCMC diagnostics
##              mcse Rhat n_eff
## (Intercept)  0.0  1.0  4990
## race         0.0  1.0  4175
## personrm     0.0  1.0  4680
## housewgt     0.0  1.0  4902
## sequenceno   0.0  1.0  5528
## mean_PPD     0.0  1.0  4731
## log-posterior 0.0  1.0  1747
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
# The coefficients of the indicators are shown below.
```

### 14.3 Graphing logistic regressions

The well-switching data described in Section 13.7 are in the folder `Arsenic`.

(a)

Fit a logistic regression for the probability of switching using log (distance to nearest safe well) as a predictor.

```
wells <- read.csv("/Users/billg/Desktop/MA-678-Homework/MA678-HW4/wells.csv")
Reg14.3 <- stan_glm(switch~log(dist),data=wells,family =binomial(link=logit),refresh=0)
summary(Reg14.3)
```

```
##
## Model Info:
## function:      stan_glm
## family:        binomial [logit]
## formula:       switch ~ log(dist)
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  3020
## predictors:    2
##
## Estimates:
##              mean    sd   10%   50%   90%
## (Intercept)  1.0    0.2  0.8   1.0   1.2
## log(dist)   -0.2    0.0 -0.3  -0.2  -0.1
##
## Fit Diagnostics:
##              mean    sd   10%   50%   90%
## mean_PPD 0.6    0.0  0.6   0.6   0.6
##
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
##
## MCMC diagnostics
##              mcse Rhat n_eff
## (Intercept)  0.0  1.0  2974
## log(dist)    0.0  1.0  2959
## mean_PPD     0.0  1.0  3243
## log-posterior 0.0  1.0  1724
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
```

(b)

Make a graph similar to Figure 13.8b displaying  $\Pr(\text{switch})$  as a function of distance to nearest safe well, along with the data.

```
fit_1 <- stan_glm(switch ~ dist, family=binomial(link="logit"), data=wells)
```

```
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 1).
## Chain 1:
## Chain 1: Gradient evaluation took 6.1e-05 seconds
## Chain 1: 1000 transitions using 10 leapfrog steps per transition would take 0.61 seconds.
## Chain 1: Adjust your expectations accordingly!
## Chain 1:
## Chain 1:
## Chain 1: Iteration:    1 / 2000 [ 0%] (Warmup)
## Chain 1: Iteration:   200 / 2000 [ 10%] (Warmup)
## Chain 1: Iteration:   400 / 2000 [ 20%] (Warmup)
## Chain 1: Iteration:   600 / 2000 [ 30%] (Warmup)
## Chain 1: Iteration:   800 / 2000 [ 40%] (Warmup)
```

```

## Chain 1: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 1: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 1: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 1: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 1: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 1: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 1: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 1:
## Chain 1: Elapsed Time: 0.227938 seconds (Warm-up)
## Chain 1: 0.27525 seconds (Sampling)
## Chain 1: 0.503188 seconds (Total)
## Chain 1:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 2).
## Chain 2:
## Chain 2: Gradient evaluation took 4.9e-05 seconds
## Chain 2: 1000 transitions using 10 leapfrog steps per transition would take 0.49 seconds.
## Chain 2: Adjust your expectations accordingly!
## Chain 2:
## Chain 2:
## Chain 2: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 2: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 2: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 2: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 2: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 2: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 2: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 2: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 2: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 2: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 2: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 2: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 2:
## Chain 2: Elapsed Time: 0.249248 seconds (Warm-up)
## Chain 2: 0.232691 seconds (Sampling)
## Chain 2: 0.481939 seconds (Total)
## Chain 2:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 3).
## Chain 3:
## Chain 3: Gradient evaluation took 5.4e-05 seconds
## Chain 3: 1000 transitions using 10 leapfrog steps per transition would take 0.54 seconds.
## Chain 3: Adjust your expectations accordingly!
## Chain 3:
## Chain 3:
## Chain 3: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 3: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 3: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 3: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 3: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 3: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 3: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 3: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 3: Iteration: 1400 / 2000 [ 70%] (Sampling)

```

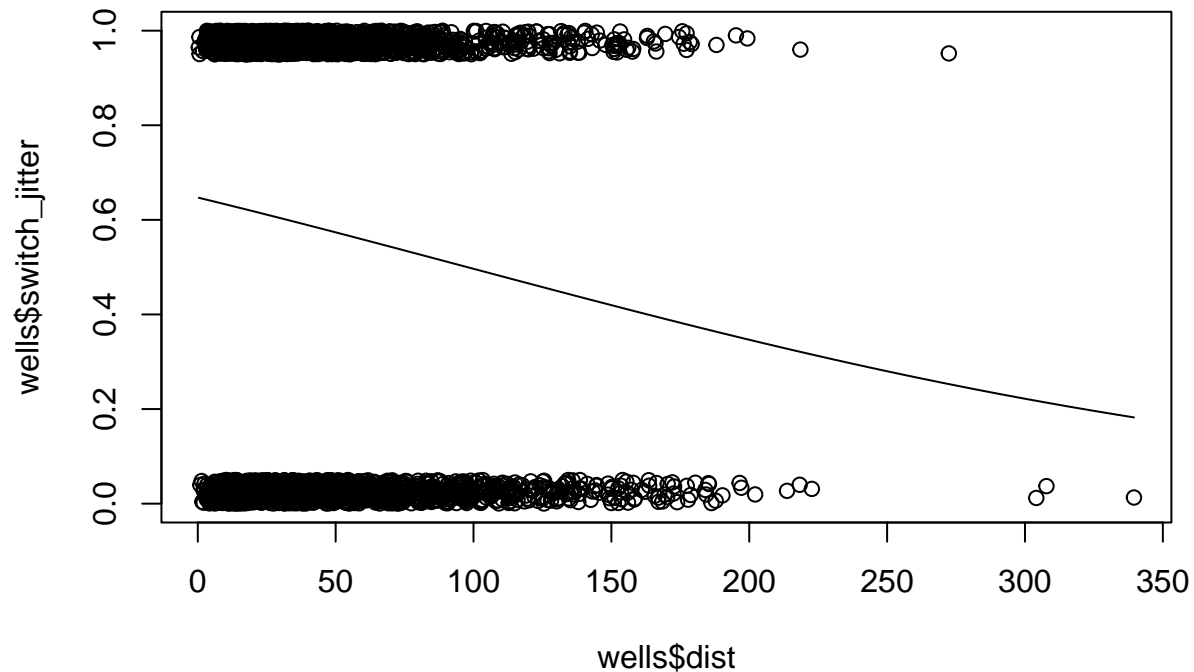


```

## Chain 3: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 3: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 3: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 3:
## Chain 3: Elapsed Time: 0.235996 seconds (Warm-up)
## Chain 3: 0.285556 seconds (Sampling)
## Chain 3: 0.521552 seconds (Total)
## Chain 3:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 4).
## Chain 4:
## Chain 4: Gradient evaluation took 5.5e-05 seconds
## Chain 4: 1000 transitions using 10 leapfrog steps per transition would take 0.55 seconds.
## Chain 4: Adjust your expectations accordingly!
## Chain 4:
## Chain 4:
## Chain 4: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 4: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 4: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 4: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 4: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 4: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 4: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 4: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 4: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 4: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 4: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 4: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 4:
## Chain 4: Elapsed Time: 0.252603 seconds (Warm-up)
## Chain 4: 0.275057 seconds (Sampling)
## Chain 4: 0.52766 seconds (Total)
## Chain 4:

jitter_binary <- function(a, jitt=0.05){
  ifelse(a==0, runif(length(a), 0, jitt), runif(length(a), 1 - jitt, 1))
}
wells$switch_jitter <- jitter_binary(wells$switch)
plot(wells$dist, wells$switch_jitter)
curve(invlogit(coef(fit_1)[1] + coef(fit_1)[2]*x), add=TRUE)

```

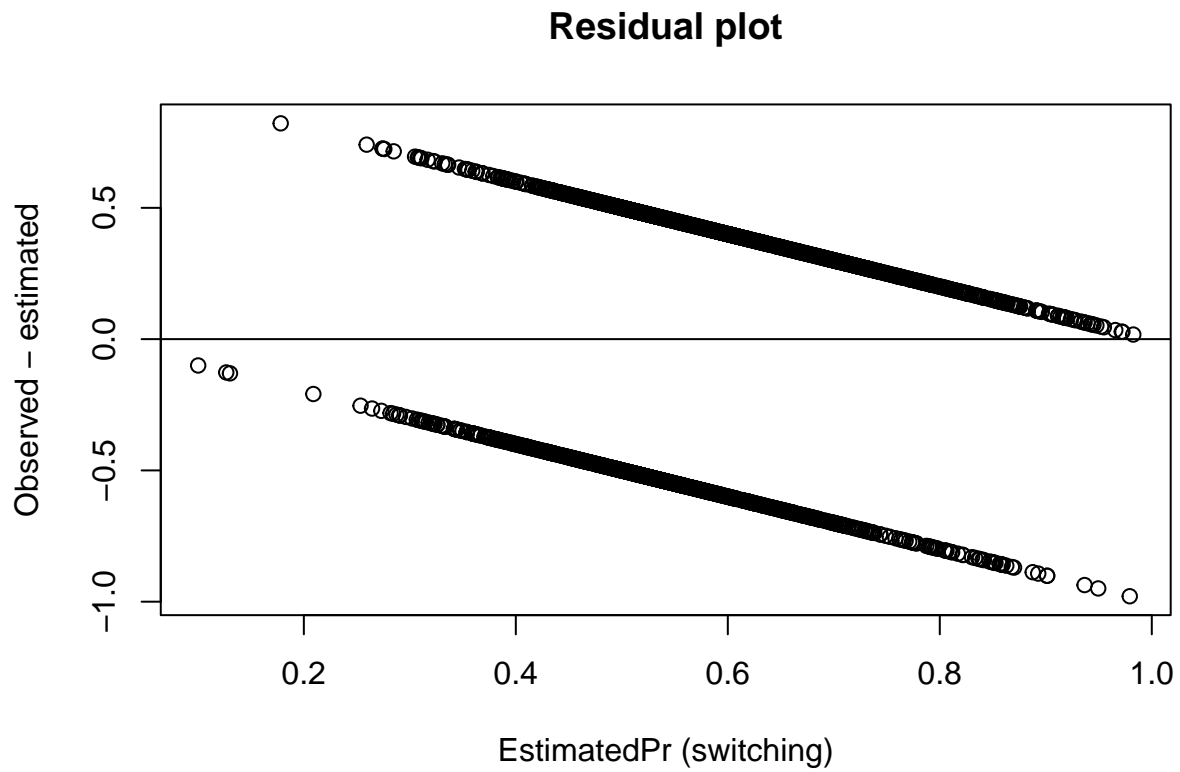


(c)

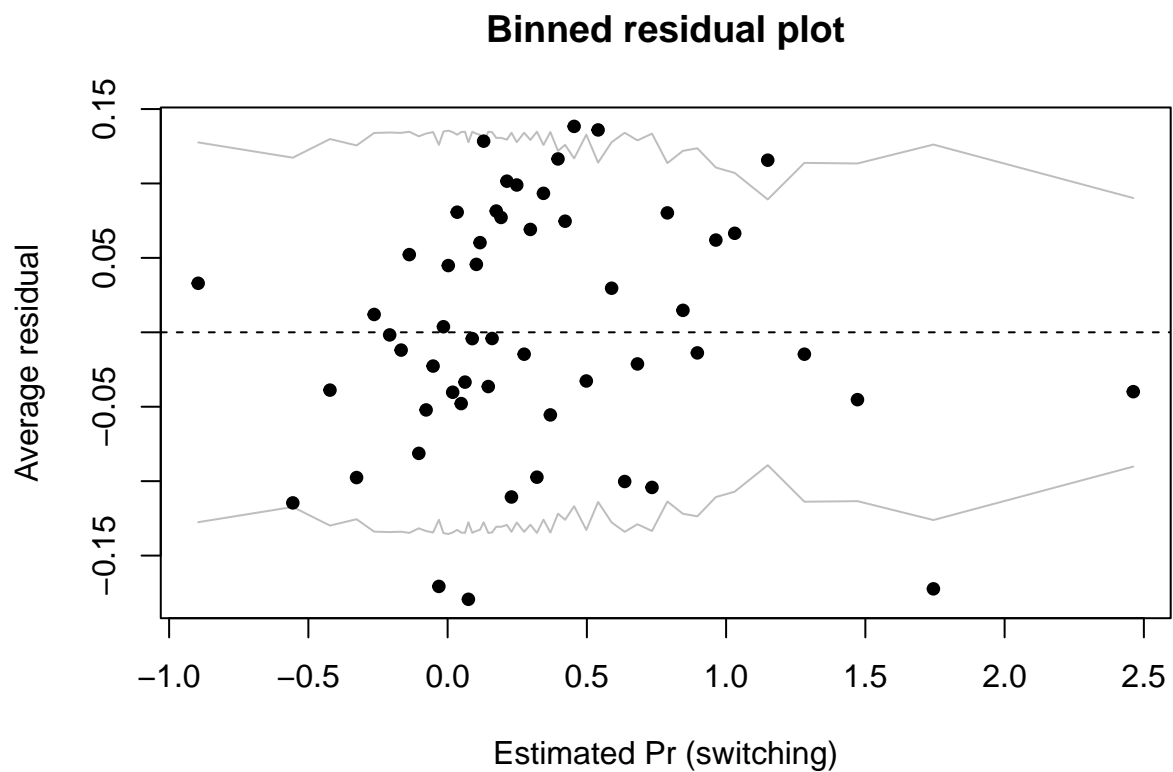
Make a residual plot and binned residual plot as in Figure 14.8.

```
library(arm)

## Loading required package: MASS
## Loading required package: Matrix
## Loading required package: lme4
##
## arm (Version 1.13-1, built: 2022-8-25)
## Working directory is /Users/billg/Desktop/MA-678-Homework/MA678-HW4
##
## Attaching package: 'arm'
## The following objects are masked from 'package:rstanarm':
##
##   invlogit, logit
fit_4 <- stan_glm(switch ~ dist100 + arsenic + dist100:arsenic,
  family=binomial(link="logit"), data=wells,refresh=0)
res <- resid(fit_4)
plot(fitted(fit_4), res,xlab="EstimatedPr (switching)",ylab="Observed - estimated",main="Residual plot",
abline(0,0)
```



```
x <- predict(fit_4)
binnedplot(x,res,xlab="Estimated Pr (switching)")
```



(d)

Compute the error rate of the fitted model and compare to the error rate of the null model.

```
fitted <- fitted(fit_4)
error <- mean(abs(wells$switch-mean(fitted)))
round(error,3)
```

```
## [1] 0.489
```

```
error_null <- mean(abs(wells$switch-fitted))
round(error_null,3)
```

```
## [1] 0.459
```

(e)

Create indicator variables corresponding to `dist < 100`; `dist` between 100 and 200; and `dist > 200`. Fit a logistic regression for `Pr(switch)` using these indicators. With this new model, repeat the computations and graphs for part (a) of this exercise.

```
variable <- NULL
variable[wells$dist<100] <- 0
variable[wells$dist>100&wells$dist<200] <- 1
variable[wells$dist>200] <- 2
Reg14.3e <- stan_glm(wells$switch~variable,family =binomial(link =logit),refresh=0)
```

```
## Warning: Omitting the 'data' argument is not recommended and may not be allowed
## in future versions of rstanarm. Some post-estimation functions (in particular
## 'update', 'loo', 'kfold') are not guaranteed to work properly unless 'data' is
## specified as a data frame.
```

```
summary(Reg14.3e)
```

```
##
```

```
## Model Info:
```

```
## function:      stan_glm
## family:        binomial [logit]
## formula:       wells$switch ~ variable
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  3020
## predictors:    2
```

```
##
```

```
## Estimates:
```

```
##           mean    sd  10%   50%   90%
## (Intercept)  0.4    0.0  0.3   0.4   0.4
## variable    -0.7    0.1 -0.8  -0.7  -0.5
```

```
##
```

```
## Fit Diagnostics:
```

```
##           mean    sd  10%   50%   90%
## mean_PPD  0.6    0.0  0.6   0.6   0.6
```

```
##
```

```
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
```

```
##
```

```
## MCMC diagnostics
```

```
##           mcse Rhat n_eff
```

```
## (Intercept)    0.0  1.0  3094
## variable      0.0  1.0  2927
## mean_PPD      0.0  1.0  3258
## log-posterior 0.0  1.0  1707
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
```

## 14.7 Model building and comparison

Continue with the well-switching data described in the previous exercise.

(a)

Fit a logistic regression for the probability of switching using, as predictors, distance, log(arsenic), and their interaction. Interpret the estimated coefficients and their standard errors.

```
wells <- read.csv("wells.csv")
Reg14.7 <- stan_glm(switch~dist+log(arsenic)+dist*log(arsenic),data=wells, family=binomial(link="logit").

##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 1).
## Chain 1:
## Chain 1: Gradient evaluation took 6.3e-05 seconds
## Chain 1: 1000 transitions using 10 leapfrog steps per transition would take 0.63 seconds.
## Chain 1: Adjust your expectations accordingly!
## Chain 1:
## Chain 1:
## Chain 1: Iteration:    1 / 2000 [  0%] (Warmup)
## Chain 1: Iteration:   200 / 2000 [ 10%] (Warmup)
## Chain 1: Iteration:   400 / 2000 [ 20%] (Warmup)
## Chain 1: Iteration:   600 / 2000 [ 30%] (Warmup)
## Chain 1: Iteration:   800 / 2000 [ 40%] (Warmup)
## Chain 1: Iteration:  1000 / 2000 [ 50%] (Warmup)
## Chain 1: Iteration:  1001 / 2000 [ 50%] (Sampling)
## Chain 1: Iteration:  1200 / 2000 [ 60%] (Sampling)
## Chain 1: Iteration:  1400 / 2000 [ 70%] (Sampling)
## Chain 1: Iteration:  1600 / 2000 [ 80%] (Sampling)
## Chain 1: Iteration:  1800 / 2000 [ 90%] (Sampling)
## Chain 1: Iteration:  2000 / 2000 [100%] (Sampling)
## Chain 1:
## Chain 1: Elapsed Time: 0.612742 seconds (Warm-up)
## Chain 1:                   0.706211 seconds (Sampling)
## Chain 1:                   1.31895 seconds (Total)
## Chain 1:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 2).
## Chain 2:
## Chain 2: Gradient evaluation took 5.9e-05 seconds
## Chain 2: 1000 transitions using 10 leapfrog steps per transition would take 0.59 seconds.
## Chain 2: Adjust your expectations accordingly!
## Chain 2:
## Chain 2:
## Chain 2: Iteration:    1 / 2000 [  0%] (Warmup)
## Chain 2: Iteration:   200 / 2000 [ 10%] (Warmup)
## Chain 2: Iteration:   400 / 2000 [ 20%] (Warmup)
```

```

## Chain 2: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 2: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 2: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 2: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 2: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 2: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 2: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 2: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 2: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 2:
## Chain 2: Elapsed Time: 0.669141 seconds (Warm-up)
## Chain 2: 0.653382 seconds (Sampling)
## Chain 2: 1.32252 seconds (Total)
## Chain 2:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 3).
## Chain 3:
## Chain 3: Gradient evaluation took 6.5e-05 seconds
## Chain 3: 1000 transitions using 10 leapfrog steps per transition would take 0.65 seconds.
## Chain 3: Adjust your expectations accordingly!
## Chain 3:
## Chain 3:
## Chain 3: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 3: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 3: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 3: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 3: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 3: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 3: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 3: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 3: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 3: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 3: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 3: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 3:
## Chain 3: Elapsed Time: 0.639812 seconds (Warm-up)
## Chain 3: 0.667639 seconds (Sampling)
## Chain 3: 1.30745 seconds (Total)
## Chain 3:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 4).
## Chain 4:
## Chain 4: Gradient evaluation took 7.3e-05 seconds
## Chain 4: 1000 transitions using 10 leapfrog steps per transition would take 0.73 seconds.
## Chain 4: Adjust your expectations accordingly!
## Chain 4:
## Chain 4:
## Chain 4: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 4: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 4: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 4: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 4: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 4: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 4: Iteration: 1001 / 2000 [ 50%] (Sampling)

```

```
## Chain 4: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 4: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 4: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 4: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 4: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 4:
## Chain 4: Elapsed Time: 0.649931 seconds (Warm-up)
## Chain 4: 0.661741 seconds (Sampling)
## Chain 4: 1.31167 seconds (Total)
## Chain 4:
```

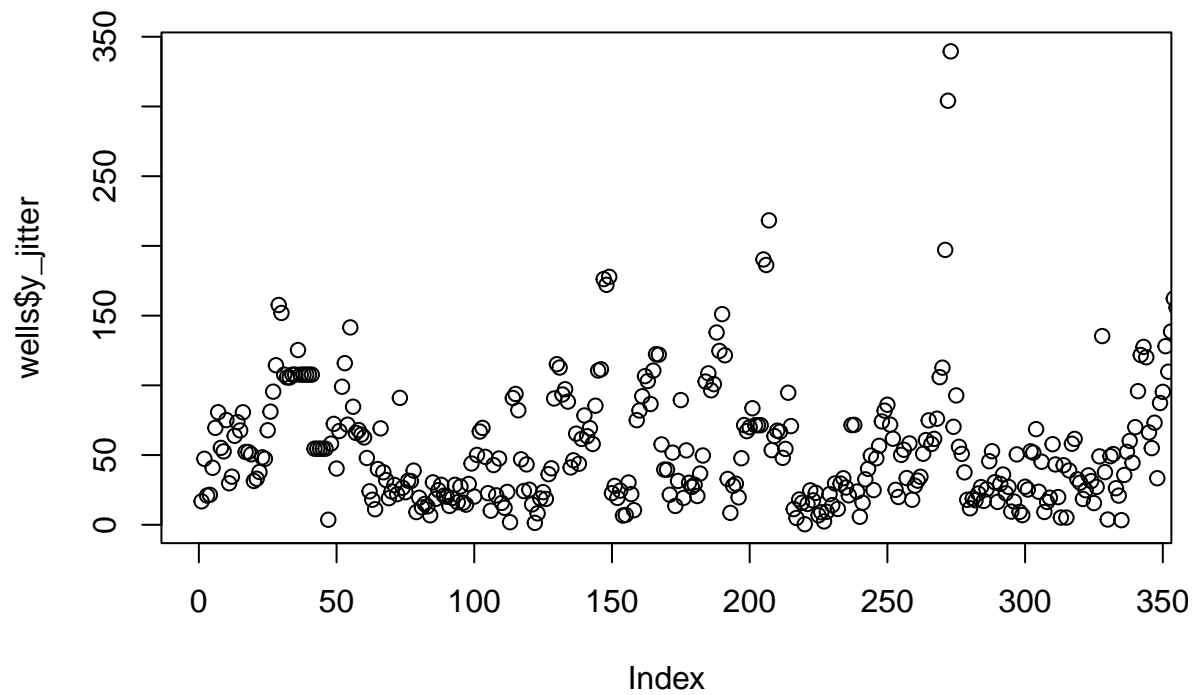
```
summary(Reg14.7)
```

```
##
## Model Info:
## function:      stan_glm
## family:        binomial [logit]
## formula:       switch ~ dist + log(arsenic) + dist * log(arsenic)
## algorithm:     sampling
## sample:        4000 (posterior sample size)
## priors:        see help('prior_summary')
## observations:  3020
## predictors:    4
##
## Estimates:
##              mean    sd  10%  50%  90%
## (Intercept)    0.5    0.1  0.4   0.5   0.6
## dist           0.0    0.0  0.0   0.0   0.0
## log(arsenic)    1.0    0.1  0.8   1.0   1.1
## dist:log(arsenic) 0.0    0.0  0.0   0.0   0.0
##
## Fit Diagnostics:
##              mean    sd  10%  50%  90%
## mean_PPD 0.6    0.0  0.6   0.6   0.6
##
## The mean_ppd is the sample average posterior predictive distribution of the outcome variable (for de
##
## MCMC diagnostics
##              mcse Rhat n_eff
## (Intercept)    0.0  1.0  2344
## dist           0.0  1.0  1965
## log(arsenic)    0.0  1.0  1606
## dist:log(arsenic) 0.0  1.0  1396
## mean_PPD        0.0  1.0  3047
## log-posterior    0.0  1.0  1242
##
## For each parameter, mcse is Monte Carlo standard error, n_eff is a crude measure of effective sample
```

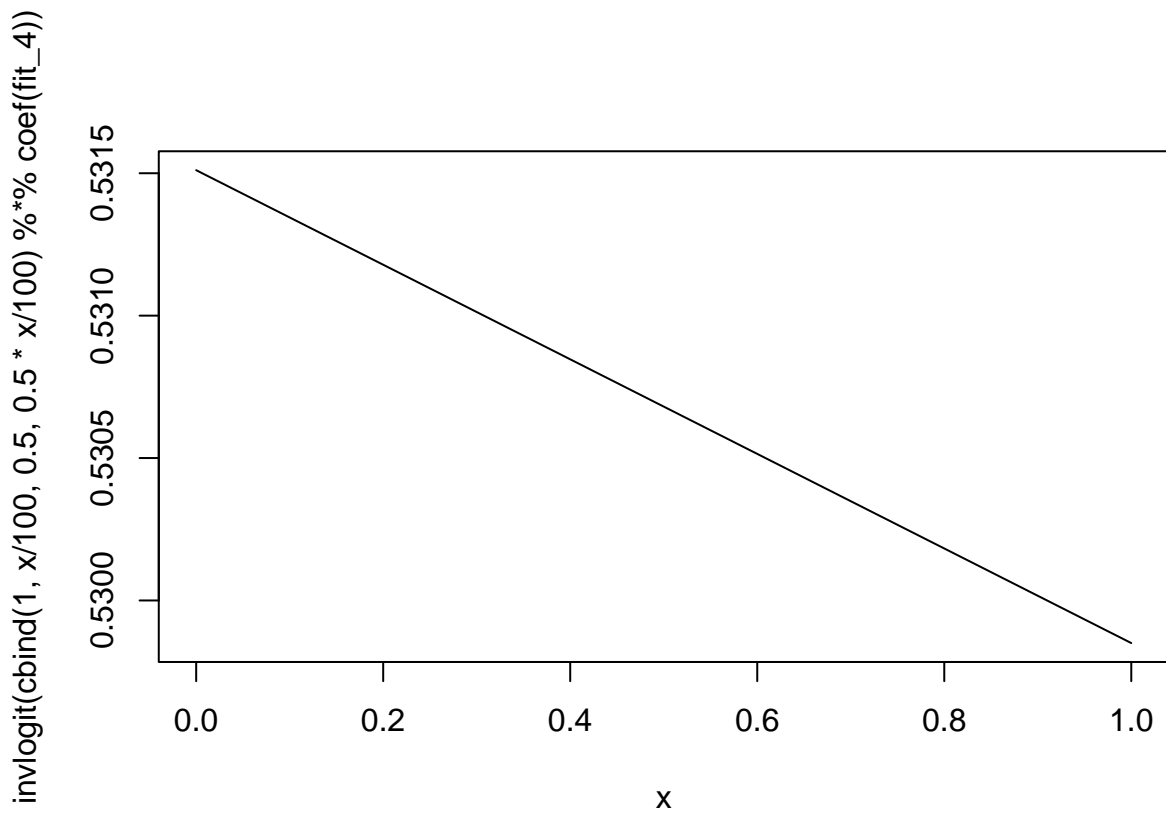
(b)

Make graphs as in Figure 14.3 to show the relation between probability of switching, distance, and arsenic level.

```
plot(wells$dist, wells$y_jitter, xlim=c(0,max(wells$dist)))
```

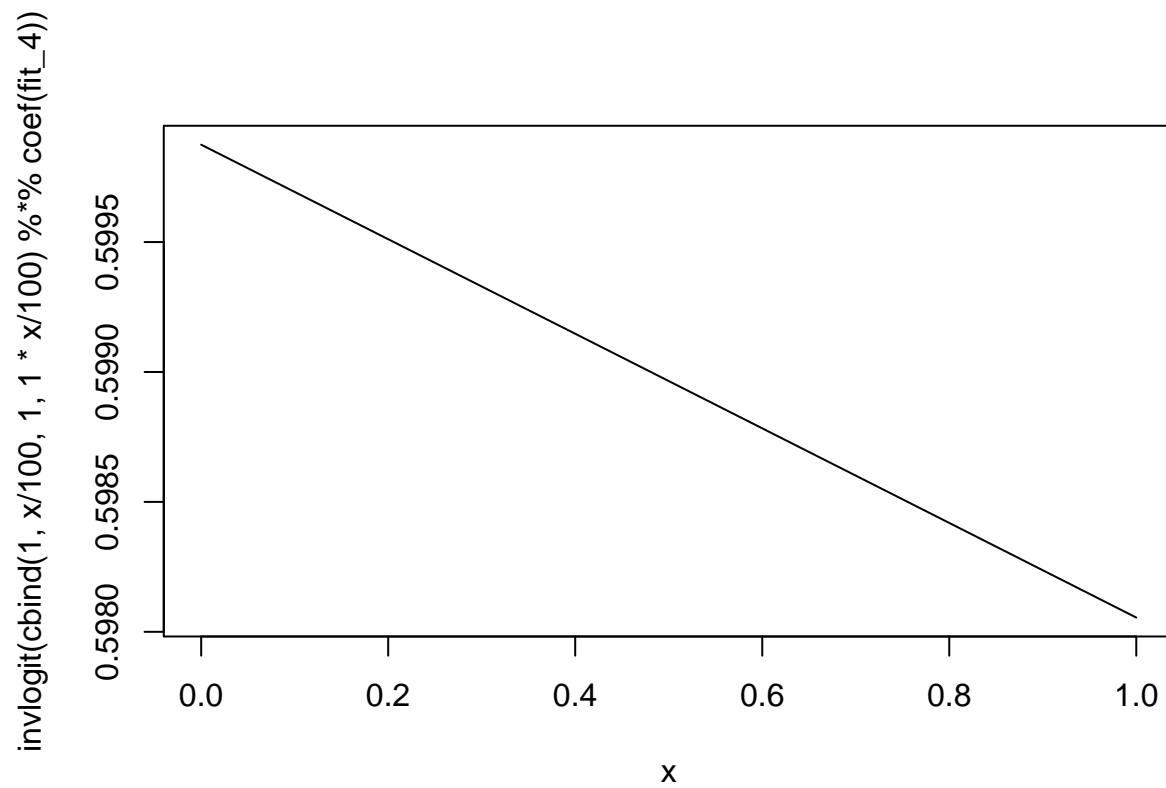


```
curve(invlogit(cbind(1, x/100, 0.5, 0.5*x/100) %*% coef(fit_4)))
```

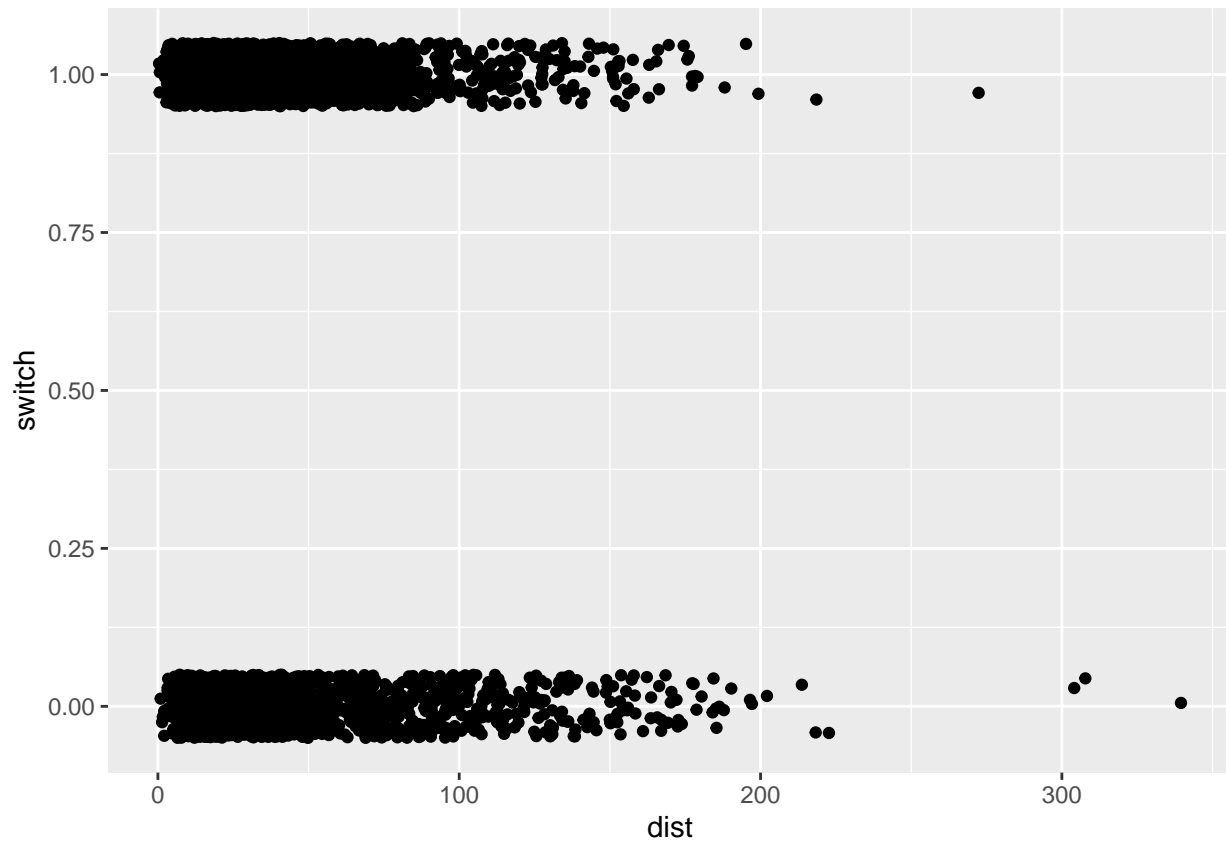


```
curve(invlogit(cbind(1, x/100, 1.0, 1.0*x/100) %*% coef(fit_4)))
```





```
library(ggplot2)
ggplot(wells,aes(dist,switch))+
  geom_jitter(position=position_jitter(height=0.05))
```



(c)

Following the procedure described in Section 14.4, compute the average predictive differences corresponding to:

- i. A comparison of `dist = 0` to `dist = 100`, with `arsenic` held constant.
- ii. A comparison of `dist = 100` to `dist = 200`, with `arsenic` held constant.
- iii. A comparison of `arsenic = 0.5` to `arsenic = 1.0`, with `dist` held constant.
- iv. A comparison of `arsenic = 1.0` to `arsenic = 2.0`, with `dist` held constant.

Discuss these results.

```
Reg14.7 <- stan_glm(switch~dist+log(arsenic)+dist*log(arsenic),data=wells, family=binomial(link="logit"))
```

```
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 1).
## Chain 1:
## Chain 1: Gradient evaluation took 7.2e-05 seconds
## Chain 1: 1000 transitions using 10 leapfrog steps per transition would take 0.72 seconds.
## Chain 1: Adjust your expectations accordingly!
## Chain 1:
## Chain 1:
## Chain 1: Iteration:    1 / 2000 [ 0%] (Warmup)
## Chain 1: Iteration:  200 / 2000 [10%] (Warmup)
## Chain 1: Iteration:  400 / 2000 [20%] (Warmup)
## Chain 1: Iteration:  600 / 2000 [30%] (Warmup)
```

```

## Chain 1: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 1: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 1: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 1: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 1: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 1: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 1: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 1: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 1:
## Chain 1: Elapsed Time: 0.653142 seconds (Warm-up)
## Chain 1: 0.695213 seconds (Sampling)
## Chain 1: 1.34835 seconds (Total)
## Chain 1:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 2).
## Chain 2:
## Chain 2: Gradient evaluation took 6.4e-05 seconds
## Chain 2: 1000 transitions using 10 leapfrog steps per transition would take 0.64 seconds.
## Chain 2: Adjust your expectations accordingly!
## Chain 2:
## Chain 2:
## Chain 2: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 2: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 2: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 2: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 2: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 2: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 2: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 2: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 2: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 2: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 2: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 2: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 2:
## Chain 2: Elapsed Time: 0.661922 seconds (Warm-up)
## Chain 2: 0.699523 seconds (Sampling)
## Chain 2: 1.36145 seconds (Total)
## Chain 2:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 3).
## Chain 3:
## Chain 3: Gradient evaluation took 6e-05 seconds
## Chain 3: 1000 transitions using 10 leapfrog steps per transition would take 0.6 seconds.
## Chain 3: Adjust your expectations accordingly!
## Chain 3:
## Chain 3:
## Chain 3: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 3: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 3: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 3: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 3: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 3: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 3: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 3: Iteration: 1200 / 2000 [ 60%] (Sampling)

```

```

## Chain 3: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 3: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 3: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 3: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 3:
## Chain 3: Elapsed Time: 0.612145 seconds (Warm-up)
## Chain 3: 0.65821 seconds (Sampling)
## Chain 3: 1.27035 seconds (Total)
## Chain 3:
##
## SAMPLING FOR MODEL 'bernoulli' NOW (CHAIN 4).
## Chain 4:
## Chain 4: Gradient evaluation took 6.1e-05 seconds
## Chain 4: 1000 transitions using 10 leapfrog steps per transition would take 0.61 seconds.
## Chain 4: Adjust your expectations accordingly!
## Chain 4:
## Chain 4:
## Chain 4: Iteration: 1 / 2000 [ 0%] (Warmup)
## Chain 4: Iteration: 200 / 2000 [ 10%] (Warmup)
## Chain 4: Iteration: 400 / 2000 [ 20%] (Warmup)
## Chain 4: Iteration: 600 / 2000 [ 30%] (Warmup)
## Chain 4: Iteration: 800 / 2000 [ 40%] (Warmup)
## Chain 4: Iteration: 1000 / 2000 [ 50%] (Warmup)
## Chain 4: Iteration: 1001 / 2000 [ 50%] (Sampling)
## Chain 4: Iteration: 1200 / 2000 [ 60%] (Sampling)
## Chain 4: Iteration: 1400 / 2000 [ 70%] (Sampling)
## Chain 4: Iteration: 1600 / 2000 [ 80%] (Sampling)
## Chain 4: Iteration: 1800 / 2000 [ 90%] (Sampling)
## Chain 4: Iteration: 2000 / 2000 [100%] (Sampling)
## Chain 4:
## Chain 4: Elapsed Time: 0.660538 seconds (Warm-up)
## Chain 4: 0.654715 seconds (Sampling)
## Chain 4: 1.31525 seconds (Total)
## Chain 4:

b <- coef(Reg14.7)
#i
hi <- 100
lo <- 0
delta <- invlogit(b[1] + b[2]*hi + b[3]*log(wells$arsenic) + b[4]*log(wells$arsenic)*hi) -
  invlogit(b[1] + b[2]*lo + b[3]*log(wells$arsenic) + b[4]*log(wells$arsenic)*lo)
round(mean(delta), 2)

## [1] -0.21

#i
hi <- 200
lo <- 100
delta2 <- invlogit(b[1] + b[2]*hi + b[3]*log(wells$arsenic) + b[4]*log(wells$arsenic)*hi) -
  invlogit(b[1] + b[2]*lo + b[3]*log(wells$arsenic) + b[4]*log(wells$arsenic)*lo)
round(mean(delta2), 2)

## [1] -0.21

#i i i
hi <- 1.0

```

```

lo <- 0.5
delta3 <- invlogit(b[1] + b[2]*wells$dist + b[3]*hi + b[4]*log(wells$dist)*hi) -
  invlogit(b[1] + b[2]*wells$dist + b[3]*lo + b[4]*log(wells$dist)*lo)
round(mean(delta3), 2)

```

```
## [1] 0.1
```

```
#iv
```

```

hi <- 2.0
lo <- 1.0
delta4 <- invlogit(b[1] + b[2]*wells$dist + b[3]*hi + b[4]*log(wells$dist)*hi) -
  invlogit(b[1] + b[2]*wells$dist + b[3]*lo + b[4]*log(wells$dist)*lo)
round(mean(delta4), 2)

```

```
## [1] 0.14
```