# Multinomial Response - Categorical Data

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## Multicategory Responses

### **Multinomial Probability Distribution**

The categorical response variable Y now has multiple levels j = 1, ...J, where each category has  $\pi_j = P(Y = j)$  and  $N_j$  is the number of trails responding with category j. These may or may not be ordinal responses. The PMF is known as the "multinomial probability distribution". (This is for a *single sample* of n observations.)

$$P(N_1 = n_1, ...N_j = n_j) = \frac{n!}{\prod_{j=1}^J n_j!} \prod_{j=1}^J \pi_j^{n_j}$$

The MLE for each  $\pi_i$  is  $\hat{\pi}_i = n_i/n$  (ie. proportion in each category).

```
pi_j \leftarrow c(0.25, 0.35, 0.2, 0.1, 0.1) # assume these are the true probabilities of each class n_j \leftarrow rmultinom(n = 1, size = 1000, prob = pi_j) # just create 1 sample with 1000 trials (data.frame(n_j, pihat_j = n_j/1000, pi_j)) # shows n_j randomly generated from each class
```

#### Contingency Tables for mixture of two multinomial distributions

If you have *one* multinomial distribution that is characterized by levels of two random variables X, Y, the data can be represented in a contingency table with a row for each X and a column for each Y. The PMF for  $N_{11}, ... N_{IJ}$  is modified to:

$$P(N_{11} = n_{11}, ...N_{IJ} = n_{IJ}) = \frac{n!}{\prod_{i=1}^{I} \prod_{j=1}^{J} n_{ij}!} \prod_{i=1}^{I} \prod_{j=1}^{J} \pi_{ij}^{n_{ij}}$$

This isn't really a new equation, just new terminology to account for the multiple rows and columns in the contingency table.

```
##
              2
## X
          1
##
     1 0.2 0.2 0.1
     2 0.3 0.1 0.1
##
# simulate a sample with 1000 obsefvations
n_ij <- rmultinom(n = 1, size = 1000, prob = pi_ij)</pre>
(c_{table} \leftarrow array(data = n_{ij}, dim = c(2, 3), dimnames = list(X = 1:2,
    Y = 1:3)))
##
      Y
## X
              2
          1
                   3
##
     1 194 210
     2 288 100 111
(pi_hat_table <- c_table/sum(c_table))</pre>
##
## X
                         3
     1 0.194 0.21 0.097
##
     2 0.288 0.10 0.111
```

In some cases, instead of having one multinomial distribution, we may intentionally sample from I different groups (corresponding to the rows in the contingency table). This is called the "product multinomial model". In this case, we have to replace  $\pi_{ij}$  values with  $\pi_{j|i}$ , so that probabilities in each row add up to 1.

## Independence

If X does not have an effect on probabilities of outcomes of Y, they are *independent*. In this case, the probability of a joint outcome of X = i, Y = j is the product of the marginal probabilities P(X), P(Y). In our notation, that is  $\pi_{ij} = \pi_{i+} \cdot \pi_{+j}$ .

We often want to test for independence; ie. we want to know if  $\pi_{i|1} = ... = \pi_{i|I} = \pi_{+i}$  for all J.

Testing for independence:

 $H_0: \pi_{ij} = \pi_{i+} \cdot \pi_{+j}$  for all i and j.  $(H_a: \text{at least some aren't equal.})$ 

We use a Pearson  $\chi^2$  test. The test statistic is:

$$X^2 = \frac{(\text{observed count - estimated expected count})^2}{\text{estimated expected count}} = \sum_{i=1}^I \sum_{j=1}^J \frac{(n_{ij} - n\hat{\pi}_{i+}\hat{\pi}_{+j})^2}{n\hat{\pi}_{i+}\hat{\pi}_{+j}} = \sum_{i=1}^I \sum_{j=1}^J \frac{(n_{ij} - n_{i+}n_{+j}/n)^2}{n_{i+}n_{+j}/n}$$

Alternately, we can do a LRT with  $\Lambda = \frac{\text{Max likelihood under } H_0}{\text{Max likelihood under } H_0 \text{ or } H_a}$ . Our test statistic in this case is:

$$-2log(\Lambda) = 2\sum_{i=1}^{I} \sum_{j=1}^{J} n_{ij} log(\frac{n_{ij}}{n_{i+}n_{+j}/n})$$

With either test statistic, if  $H_0$  is true,  $X^2$  has an approximate  $\chi^2_{(I-1)(J-1)}$  distribution. Reject  $H_0$  if  $X^2$  or  $-2log(\Lambda)$  is great than  $\chi^2$ . NOTE: test results for  $X^2$  and  $-2log(\Lambda)$  may vary greatly for small sample sizes; generally we want an expected value in each cell to be > 1 or > 5.

```
# example of reading data in to a contingency table
diet <- read.csv("Fiber.csv")</pre>
head(diet)
##
     fiber bloat count
## 1 bran
             high
## 2
       gum
             high
                       5
## 3 both
             high
## 4 none
             high
                       0
## 5 bran medium
                       1
       gum medium
diet$fiber <- factor(diet$fiber, levels = c("none", "bran", "gum", "both"))</pre>
diet$bloat <- factor(diet$bloat, levels = c("none", "low", "medium", "high"))</pre>
(diet_table <- xtabs(count ~ fiber + bloat, data = diet))</pre>
##
         bloat
## fiber none low medium high
##
     none
                              0
##
     bran
             7
             2
                 2
                         3
##
     gum
                              5
##
     both
                 5
Now, here are 3 ways to test for independence:
(indep_test <- chisq.test(diet_table, correct = FALSE)) # don't let it apply extra correction
## Warning in chisq.test(diet_table, correct = FALSE): Chi-squared
## approximation may be incorrect
##
  Pearson's Chi-squared test
##
##
## data: diet_table
## X-squared = 16.943, df = 9, p-value = 0.04962
summary(diet_table)
## Call: xtabs(formula = count ~ fiber + bloat, data = diet)
## Number of cases in table: 48
## Number of factors: 2
## Test for independence of all factors:
## Chisq = 16.943, df = 9, p-value = 0.04962
## Chi-squared approximation may be incorrect
library(vcd) # calculates X^2 test only
## Loading required package: grid
```

### assocstats(diet\_table) # calculates both X2 and LRT tests ## $X^2 df P(> X^2)$ ## Likelihood Ratio 18.880 9 0.026230 ## Pearson 16.943 9 0.049621 ## ## Phi-Coefficient : NA ## Contingency Coeff.: 0.511 : 0.343 ## Cramer's V qchisq(p = 0.95, df = 9) # this is the critical value that the test stats are being compared ## [1] 16.91898 $(p_x2 \leftarrow 1 - pchisq(16.943, df = 9))$ # we can also validate the p-values reported above ## [1] 0.04961611 (p\_lrt <- 1 - pchisq(18.88, df = 9)) # we can also validate the p-values reported above ## [1] 0.02623278

Note that df = 9 = (I - 1)(J - 1) = (4 - 1)(4 - 1). All 3 methods give same result for  $X^2$  test. Small p-values mean we reject the null hypothesis of independence.

We can also get the expected counts to see whether they meet the > 1 or > 5 thresholds for  $\chi^2$  being a good approximation. The warning in some of the above methods is because some of these are < 5.

#### indep\_test\$expected

```
## bloat

## fiber none low medium high

## none 4.25 3.75 2.25 1.75

## bran 4.25 3.75 2.25 1.75

## gum 4.25 3.75 2.25 1.75

## both 4.25 3.75 2.25 1.75
```

## Nominal Response Regression Models

We can define odds as a comparison of any pair of response categories; a popular regression model for multinomial resonness is by forming the odds of the remaining J-1 categories against a base category. This multinomial regression model looks like:

$$log(\pi_j/\pi_1) = \beta_{j0} + \beta_{j1}x_1 + \dots + \beta_{jp}x_p \forall j = 2, \dots J$$

Note that there is a separate set of  $\beta$  parameters for each response category, so each response's log-odds can relate to the explanatory variables in a different way. It's easy to compare other categories:

$$log(\pi_2/\pi_3) = log(\pi_2/\pi_1) - log(\pi_3/\pi_1) = log(\pi_2) - log(\pi_3) = (\beta_{20} - \beta_{30}) + (\beta_{21} - \beta_{31})x_1 + \dots + (\beta_{2p} - \beta_{3p})x_p$$

### Calculating Probabilities

We can also calculate probabilities by maximum likelihood:

$$\pi_{i} = \pi_{1} exp(\beta_{i0} + \beta_{i1}x_{1} + ... + \beta_{ip}x_{p}) \forall j = 2, ...J$$

But we need to find  $\pi_1$  first. Since  $\pi_1 + \pi_2 + ... \pi_J = 1$ , we can get an expression for  $\pi_1$ :

$$\pi_1 = \frac{1}{1 + \sum_{j=2}^{J} exp(\beta_{j0} + \beta_{j1}x_1 + \dots + \beta_{jp}x_p)}$$

Combining, we get:

$$\pi_{j} = \frac{exp(\beta_{j0} + \beta_{j1}x_{1} + \dots + \beta_{jp}x_{p})}{1 + \sum_{j=2}^{J} exp(\beta_{j0} + \beta_{j1}x_{1} + \dots + \beta_{jp}x_{p})} \forall j = 2, \dots J$$

For a sample of size m observations, the likelihood function is the product of m multinomial distributions, with  $\pi_j$  as described in these equations. Iterative numerical procedures are used to find these MLEs using nnet::multinom(...):

```
wheat <- read.csv("wheat.csv")</pre>
head(wheat)
##
           density hardness
                                size weight moisture
                                                          type
       hrw 1.349253 60.32952 2.30274 24.6480 12.01538 Healthy
## 2
       hrw 1.287440 56.08972 2.72573 33.2985 12.17396 Healthy
## 3
      hrw 1.233985 43.98743 2.51246 31.7580 11.87949 Healthy
      hrw 1.336534 53.81704 2.27164 32.7060 12.11407 Healthy
## 4
      hrw 1.259040 44.39327 2.35478 26.0700 12.06487 Healthy
## 5
       hrw 1.300258 48.12066 2.49132 33.2985 12.18577 Healthy
levels(wheat$type) # observe that 'Healthy' is the base case
## [1] "Healthy" "Scab"
                           "Sprout"
library(nnet)
wheat_fit <- multinom(type ~ class + density + hardness + size + weight +
    moisture, data = wheat)
## # weights: 24 (14 variable)
## initial value 302.118379
## iter 10 value 234.991271
## iter 20 value 192.127549
## final value 192.112352
## converged
```

```
summary(wheat_fit)
## Call:
```

## class

0.964 2

```
## multinom(formula = type ~ class + density + hardness + size +
       weight + moisture, data = wheat)
##
## Coefficients:
##
          (Intercept)
                         classsrw
                                    density
                                               hardness
                                                              size
                                                                        weight
## Scab
             30.54650 -0.6481277 -21.59715 -0.01590741 1.0691139 -0.2896482
             19.16857 -0.2247384 -15.11667 -0.02102047 0.8756135 -0.0473169
## Sprout
##
             moisture
## Scab
           0.10956505
## Sprout -0.04299695
##
## Std. Errors:
##
          (Intercept)
                       classsrw density
                                             hardness
                                                            size
                                                                     weight
## Scab
             4.289865 0.6630948 3.116174 0.010274587 0.7722862 0.06170252
             3.767214 0.5009199 2.764306 0.008105748 0.5409317 0.03697493
## Sprout
##
           moisture
## Scab
          0.1548407
## Sprout 0.1127188
##
## Residual Deviance: 384.2247
## AIC: 412.2247
We interpret these results as:
```

```
log(\hat{\pi}_{scab}/\hat{\pi}_{healthy}) = 30.55 - 0.65I(class = SRW) - 21.60density - 0.016hardness + ...
```

$$log(\hat{\pi}_{sprout}/\hat{\pi}_{healthy}) = 19.17 - 0.22I(class = SRW) - 15.12 density - 0.021 hardness + \dots$$

NOTE: mcprofile(...) and confint(...) cannot be used for profile likelihood ratio intervals for multinomial models. Wald intervals are possible to calculate because multinom(...) gives you standard errors.

LRTs for significance of coefficients are straightforward using Anova(...). Typically we want to explore whether a particular explanatory variable as an effect on all response categories (not just one). That hypothesis test is  $H_0: \beta_{jr} = 0 \forall j = 2, ..., J$ .

```
library(car)
Anova(wheat_fit)
## Analysis of Deviance Table (Type II tests)
##
## Response: type
            LR Chisq Df Pr(>Chisq)
##
```

0.6175

```
90.555 2 < 2.2e-16 ***
## density
               7.074 2
## hardness
                            0.0291 *
               3.211 2
                            0.2008
## size
              28.230 2 7.411e-07 ***
## weight
## moisture
               1.193 2
                            0.5506
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
Here are two equivalent ways to see the estimated probabilities for each class for each observation:
head(wheat_fit$fitted.values)
##
       Healthy
                      Scab
                               Sprout
## 1 0.8552110 0.046396827 0.09839221
## 2 0.7492553 0.021572158 0.22917255
## 3 0.5172800 0.068979903 0.41374011
## 4 0.8982064 0.006740716 0.09505287
## 5 0.5103245 0.176260796 0.31341473
## 6 0.7924907 0.015304122 0.19220522
head(predict(wheat_fit, newdata = wheat, type = "probs"))
##
       Healthy
                      Scab
                               Sprout
## 1 0.8552110 0.046396827 0.09839221
## 2 0.7492553 0.021572158 0.22917255
## 3 0.5172800 0.068979903 0.41374011
## 4 0.8982064 0.006740716 0.09505287
## 5 0.5103245 0.176260796 0.31341473
## 6 0.7924907 0.015304122 0.19220522
head(predict(wheat_fit, newdata = wheat, type = "class"))
```

```
## [1] Healthy Healthy Healthy Healthy Healthy
## Levels: Healthy Scab Sprout
```

### Odds Ratios for Multinomial Models

Odds of a category j response vs. a category 1 response change by  $exp(c\beta_{jr})$  for every c-unit change in  $x_r$ , holding other variables constant.

Odds of a category j response vs. a category j' response change by  $exp[c(\beta_{jr} - \beta_{j'r})]$  for every c - unit change in  $x_r$ , holding other variables constant.

```
sd_wheat <- apply(wheat[, -c(1, 7)], MARGIN = 2, FUN = sd) # find st dev for continuous vars
(c_value <- c(1, sd_wheat))

## density hardness size weight moisture
## 1.0000000 0.1313021 27.3561563 0.4906125 7.9154398 2.0332132
beta_hat_scab <- coef(wheat_fit)[1, 2:7]
beta_hat_sprout <- coef(wheat_fit)[2, 2:7]</pre>
```

```
# We calculate the OR (according to a c = 1 * sd change) for all
# variables, but can only change one at a time
or_scab <- exp(c_value * beta_hat_scab)</pre>
round(or_scab, 2) # OR for a c-unit increase (scab vs healthy)
##
             density hardness
                                   size
                                           weight moisture
##
       0.52
                0.06
                                   1.69
                                             0.10
                          0.65
                                                      1.25
round(1/or_scab, 2) # OR for a c-unit descrease (scab vs healthy)
##
             density hardness
                                           weight moisture
                                   size
##
               17.04
                                   0.59
                                             9.90
       1.91
                          1.55
                                                      0.80
or_sprout <- exp(c_value * beta_hat_sprout)</pre>
round(or_sprout, 2) # OR for a c-unit increase (sprout vs healthy)
##
             density hardness
                                   size
                                           weight moisture
##
       0.80
                0.14
                          0.56
                                   1.54
                                             0.69
                                                      0.92
round(1/or sprout, 2) # OR for a c-unit descrease (sprout vs healthy)
##
             density hardness
                                           weight moisture
                                   size
##
       1.25
                7.28
                          1.78
                                   0.65
                                             1.45
                                                      1.09
```

#### Confidence Intervals for Parameters

The confint(...) function for multinomial regression does *not* use likelihood ratios; it uses Wald. We calculate CI for Odds Ratios similarly to binomial case, by finding the CI for the linear predictor first, and then exponentiating.

Note how confidence intervals are stored: [coefficients, lower:upper limits, class sequence from fit output].

```
## classsrw
                -1.94776958
                               0.651514098
## density
               -27.70474380 -15.489565975
## hardness
                -0.03604523
                               0.004230411
## size
                -0.44453927
                               2.582767006
## weight
                -0.41058295 -0.168713512
## moisture
                -0.19391723
                              0.413047326
##
## , , Sprout
##
##
                      2.5 %
                                   97.5 %
## (Intercept) 11.78496433 26.552173165
```

```
## classsrw
            -1.20652328 0.757046542
## density
             -20.53461137 -9.698731394
## hardness
              -0.03690744 -0.005133494
## size
               -0.18459306 1.935820104
## weight
               -0.11978643 0.025152642
## moisture
               -0.26392179 0.177927888
or_scab_ci <- exp(c_value * ci_betas[2:7, 1:2, 1]) # make sure to get the set for scab
round(or_scab_ci, 2) # OR CI for a c-unit increase
##
           2.5 % 97.5 %
## classsrw 0.14
                  1.92
## density
            0.03
                 0.13
## hardness 0.37
                 1.12
## size
            0.80
                  3.55
## weight
            0.04
                  0.26
## moisture 0.67
                  2.32
round(1/or_scab_ci, 2) # OR CI for a c-unit decrease
##
           2.5 % 97.5 %
## classsrw 7.01
                  0.52
## density 38.00
                  7.64
## hardness 2.68
                 0.89
## size
            1.24
                 0.28
## weight
           25.79
                   3.80
## moisture 1.48
                 0.43
or_sprout_ci <- exp(c_value * ci_betas[2:7, 1:2, 2]) # make sure to get the set for sprout
round(or_sprout_ci, 2) # OR CI for a c-unit increase
##
           2.5 % 97.5 %
## classsrw 0.30
                 2.13
                   0.28
## density
            0.07
## hardness 0.36
                 0.87
## size
            0.91
                   2.59
## weight
            0.39
                  1.22
## moisture 0.58
                  1.44
round(1/or_sprout_ci, 2) # OR CI for a c-unit decrease
##
           2.5 % 97.5 %
## classsrw 3.34 0.47
## density 14.82
                   3.57
## hardness 2.74
                  1.15
## size
            1.09
                 0.39
## weight
            2.58
                 0.82
## moisture 1.71 0.70
```

## Applying Multinomial Regression model to Contingency Table to Test Independence (Alternative Approach)

The multinomial regression model provides a convenient framework for performing a LRT to test for independence (alternative to method described earlier). We create an *indicator variable* for each level of X (ie.  $x_2 = I(x_{i=2})$ ). If we have *dependence*, the model looks like:

$$H_a: log(\pi_i/\pi_1) = \beta_{i0} + \beta_{i2}x_2 + ... + \beta_{ip}x_p \forall j = 2,...J$$

Note that for consistency, we start with j=2 and  $\beta_{j2}$  so that subscripts match class levels for non-base cases. When we have *independence*, the model becomes:

$$H_0: log(\pi_i/\pi_1) = \beta_{i0} \forall j = 2, ...J$$

This means that, while each category of Y can have a difference  $\pi_j$ , these values do *not* change as a function of X. These two models can be compared in a hypothesis test. This is the same as writing  $H_0: \beta_{j2} = ... = \beta_{jI} = 0 \forall j = 2, ... J$  and  $H_a:$  not all of these  $\beta$  parameters are 0, for some j.

In the following example, weights are the counts from each cell of the contingench table, even though we aren't using the crosstab-formatted matrix.

```
library(nnet)
head(diet)
##
     fiber
            bloat count
## 1
      bran
             high
                       5
## 2
             high
       gum
## 3 both
             high
                       2
## 4
                       0
     none
             high
## 5
      bran medium
                       1
## 6
       gum medium
diet_fit_nominal <- multinom(formula = bloat ~ fiber, weights = count,</pre>
    data = diet) # not in crosstab c-table format
## # weights:
               20 (12 variable)
## initial value 66.542129
## iter 10 value 54.519963
## iter
         20 value 54.197000
## final value 54.195737
## converged
summary(diet_fit_nominal)
## Call:
## multinom(formula = bloat ~ fiber, data = diet, weights = count)
##
## Coefficients:
                       fiberbran
          (Intercept)
                                    fibergum fiberboth
## low
           -0.4057626 -0.1538545
                                   0.4055575
                                              1.322135
## medium -1.0980713 -0.8481379
                                   1.5032639
                                              1.503764
```

```
-12.4401085 -4.1103893 13.3561038 12.440403
## high
##
## Std. Errors:
##
          (Intercept)
                          fiberbran
                                       fibergum
                                                 fiberboth
## low
            0.6455526
                          0.8997698
                                       1.190217
                                                  1.056797
## medium
            0.8163281
                          1.3451836
                                       1.224593
                                                  1.224649
## high
          205.2385583 1497.8087307 205.240263 205.240994
##
## Residual Deviance: 108.3915
## AIC: 132.3915
```

To perform LRT for independence, we simply use Anova(...):

```
Anova(diet_fit_nominal)
```

```
## # weights: 8 (3 variable)
## initial value 66.542129
## final value 63.635876
## converged

## Analysis of Deviance Table (Type II tests)
##
## Response: bloat
## LR Chisq Df Pr(>Chisq)
## fiber 18.88 9 0.02623 *
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
```

Note that the p-value (0.02623) is the same as that reported earlier in the  $-2log(\Lambda)$  test with vcd:assocstats(...).

## Proportional Odds Model for Ordinal Response data

Basically, when we have ordinal data, we just apply the logit transformation to the *cumulative* probability distribution (CDF). The *cumulative* probability for Y is  $P(Y|j) = \pi_1 + \pi_2 + ...\pi_j$ . Then we apply the logit transformation:

$$logit[P(Y \leq j)] = log\left[\frac{P(Y \leq j)}{1 - P(Y \leq j)}\right] = log\left[\frac{\pi_1 + ... \pi_j}{\pi_{j+1} + ... \pi_J}\right]$$

Then the "Proportional Odds Model" applies a big simplification by assuming the  $\beta$  parameters except for the intercept are the same for all explanatory variables (for all j = 1...J - 1). This means that effects of explanatory variables are the same, no matter which cumulative probabilities were used to form the log odds! If we were to plot these logit curves, they would have the identical shape, just be shifted on the x axis according to intercept terms.

$$logit[P(Y \le j)] = \beta_{j0} + \beta_1 x_1 + ... \beta_p x_p \forall J = 1, ... J - 1$$

Also, this can be rearranged to:

$$P(Y \le j) = \frac{exp(\beta_{j0} + \beta_1 x_1 + ... \beta_p x_p)}{1 + exp(\beta_{j0} + \beta_1 x_1 + ... \beta_p x_p)} \forall J = 1, ... J - 1$$

It's important to remember that because of the nature of cumulative probabilities,  $\beta_{J0} > ... > \beta_{10}$ .

If we wanted a model where the  $\beta$  parameters were allowed to vary for level of J, that would be the "cumulative probability model" (not covered in class). This model also gets more complicated when there are more than 2 categorical explanatory variables (because we need a m-dimensional contingency table).

We can also look at how we express  $\pi_j$  in terms of these cumulative probabilities:

$$\pi_j = P(Y=j) = P(Y \leq j) - P(Y \leq j - 1) = \frac{exp(\beta_{j0} + \beta_1x_1 + \ldots)}{1 + exp(\beta_{j0} + \beta_1x_1 + \ldots)} - \frac{exp(\beta_{j-1,0} + \beta_1x_1 + \ldots)}{1 + exp(\beta_{j-1,0} + \beta_1x_1 + \ldots)} \forall j = 2, \ldots J - 1$$

#### **Estimation and Inference**

Parameters are estimated using MLE, with the MASS::polr(...) function. If we have one explanatory variable, the hypotheses of interest are  $H_0: \beta_1 = 0$  and  $H_a: \beta_1 \neq 0$ .

If we reject  $H_0$ , then the ordering of log-odds comparing  $P(Y \leq j)$  and P(Y > j) holds. Log odds progressively grow larger (or smaller, depending on sign of  $\beta_1$ ).

If we fail to reject  $H_0$ , then the log-odds comparing  $P(Y \leq j)$  and P(Y > j) do not depend on this explanatory variable. For the case of two explanatory variables, this is equivalent to *independence*.

WARNING: levels of the Y variable must be properly ordered before using polr(...). Also, all parameters generated by polr(...) need to have signs reversed, except for  $\beta_{i0}$ .

```
levels(wheat$type)
## [1] "Healthy" "Scab"
                            "Sprout"
wheat$type <- factor(wheat$type, levels = c("Scab", "Sprout", "Healthy"))</pre>
levels(wheat$type)
## [1] "Scab"
                  "Sprout" "Healthy"
library(MASS)
wheat_fit_polr <- polr(formula = type ~ class + density + hardness + size +</pre>
    weight + moisture, data = wheat, method = "logistic")
summary(wheat_fit_polr)
##
## Re-fitting to get Hessian
## polr(formula = type ~ class + density + hardness + size + weight +
       moisture, data = wheat, method = "logistic")
##
##
```

```
## Coefficients:
##
                Value Std. Error t value
## classsrw 0.17370
                        0.391764
                                   0.4434
## density 13.50534
                         1.713009
                                   7.8840
## hardness 0.01039
                        0.005932 1.7522
             -0.29253
## size
                         0.413095 -0.7081
## weight
              0.12721
                        0.029996 4.2411
## moisture -0.03902
                         0.088396 -0.4414
## Intercepts:
##
                   Value
                            Std. Error t value
## Scab|Sprout
                   17.5724
                             2.2460
                                         7.8237
## Sprout | Healthy 20.0444 2.3395
                                         8.5677
##
## Residual Deviance: 422.4178
## AIC: 438.4178
We would interpret this as:
       logit[P(Y \leq j)] = \beta_{i0} - 0.17I(SRW) - 13.51 density - 0.01 hardness + 0.29 size...
with \beta_{10} = 17.57 and \beta_{20} = 20.04. The t value column is the Wald statistic (ie. compare to 1.96).
We can also do LRTs with Anova(...):
Anova(wheat_fit_polr)
## Analysis of Deviance Table (Type II tests)
##
## Response: type
##
            LR Chisq Df Pr(>Chisq)
## class
                0.197
                       1
                             0.65749
               98.437
                          < 2.2e-16 ***
## density
                       1
## hardness
                3.084
                       1
                             0.07908 .
## size
                0.499
                       1
                             0.47982
## weight
               18.965
                       1
                           1.332e-05 ***
## moisture
                0.195
                             0.65872
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
We can also predict probabilities and classes. ** However, predict(...) does not calculate the
standard errors that would be needed in order to calculate Wald confidence intervals for \pi_i^{**}.
wheat_classes_pi_hat <- predict(wheat_fit_polr, type = "probs")</pre>
wheat_classes <- predict(wheat_fit_polr, type = "class")</pre>
head(data.frame(wheat_classes_pi_hat, wheat_classes))
            Scab
                    Sprout
                              Healthy wheat_classes
## 1 0.03661601 0.2738502 0.6895338
                                             Healthy
## 2 0.03351672 0.2576769 0.7088064
                                             Healthy
## 3 0.08379891 0.4362428 0.4799583
```

Healthy

```
## 4 0.01694278 0.1526100 0.8304472 Healthy
## 5 0.11408176 0.4899557 0.3959626 Sprout
## 6 0.02874814 0.2308637 0.7403882 Healthy
```

#### Odds Ratio

Our odds ratio interpretation as  $exp(\beta_{j0} + \beta_1 x_1 + ... \beta_p x_p)$  is natural, except that we are using cumulative probabilities. Recall:

$$logit[P(Y \le j)] = log \left[ \frac{P(Y \le j)}{1 - P(Y \le j)} \right] = \beta_{j0} + \beta_1 x_1 + \dots + \beta_p x_p$$

As usual, for a c-unit increase in  $x_1$ :

class density hardness

or\_decrease <- 1/or\_increase

$$OR = \frac{Odds_{x_1+c}(Y \le j)}{Odds_{x_1}(Y \le j)} = exp(c\beta_1)$$

The interpretation is that the odds of  $Y \leq j$  vs. Y > j change by  $exp(c\beta_1)$  for every c-unit increase in  $x_1$ , holding all other variables constant. Due to the nature of the proportional odds model, this is the same result no matter which category is used for j.

size weight moisture

```
head(wheat)
```

##

```
hrw 1.349253 60.32952 2.30274 24.6480 12.01538 Healthy
      hrw 1.287440 56.08972 2.72573 33.2985 12.17396 Healthy
      hrw 1.233985 43.98743 2.51246 31.7580 11.87949 Healthy
      hrw 1.336534 53.81704 2.27164 32.7060 12.11407 Healthy
## 5
      hrw 1.259040 44.39327 2.35478 26.0700 12.06487 Healthy
      hrw 1.300258 48.12066 2.49132 33.2985 12.18577 Healthy
## 6
levels(wheat$type)
                 "Sprout" "Healthy"
## [1] "Scab"
c_value # we calculated the standard deviations for each parameter to use as our c-unit
##
                           hardness
                 density
                                                   weight
                                                            moisture
                                          size
   1.0000000
               0.1313021 27.3561563 0.4906125 7.9154398 2.0332132
```

or\_increase <- exp(c\_value \* (-coef(wheat\_fit\_polr))) # remember we have to take negative of

Interpretations: in all cases, "holding all other variables constant"

- the estimated odds of a scab vs. sprout or healthy are 0.84 times as large for soft rather than hard wheat.
- the estimated odds of a scab vs. sprout or healthy are 0.36 times as large for a c=1sd increase in weight.
- the estimated odds of a scab vs. sprout or healthy are 5.89 times as large for a c=1sd decrease in density.

• the estimated odds of a scab vs. sprout or healthy are 2.74 times as large for a c=1sd decrease in weight.

Also, due to definition of proportional odds model, each of the following statements could also apply to "... a scab or sprout vs. healthy" (ie. comparing different class threshholds).

It is possible do do Wald and LR Intervals for the Odds Ratio.

```
(ci_betas_polr <- confint(wheat_fit_polr, level = 0.95))</pre>
## Waiting for profiling to be done...
##
## Re-fitting to get Hessian
                   2.5 %
##
                             97.5 %
## classsrw -0.595305729
                         0.9435846
## density 10.315429541 17.0363926
## hardness -0.001207582 0.0221078
## size
            -1.103021561 0.5245184
## weight
             0.069318186
                         0.1872189
## moisture -0.213254701 0.1339876
(or_increase_ci <- exp(c_value * (-ci_betas_polr))) # note these are in reversed upper/lower
##
                2.5 %
                         97.5 %
## classsrw 1.8135853 0.3892301
## density 0.2580925 0.1067876
## hardness 1.0335865 0.5461922
## size
            1.7179919 0.7731090
## weight
            0.5777096 0.2272011
## moisture 1.5427897 0.7615303
(or_decrease_ci <- 1/exp(c_value * (-ci_betas_polr))) # note these are in reversed upper/lowe</pre>
                2.5 %
##
                        97.5 %
## classsrw 0.5513940 2.569174
## density 3.8745799 9.364385
## hardness 0.9675049 1.830858
## size
            0.5820749 1.293479
            1.7309734 4.401388
## weight
## moisture 0.6481765 1.313145
```

Interpretation: Holding all other variables constant, with 95% confidence, the odds of a scab instead of a sprout or healthy response increase by between 3.87 and 9.36 times for every c=1sd decrease in density.