# Ch8-baggboost-lab

# Lab: Decision Trees

# Fitting Classification Trees

The tree library is used to construct classification and regression trees.

```
library(tree)
```

We first use classification trees to analyze the Carseats data set. In these data, Sales is a continuous variable, and so we begin by recoding it as a binary variable. We use the ifelse() function to create a variable, called High, which takes on a value of Yes if the Sales variable exceeds 8, and takes on a value of No otherwise.

```
library(ISLR2)
attach(Carseats)
High <- factor(ifelse(Sales <= 8, "No", "Yes"))</pre>
```

Finally, we use the data.frame() function to merge High with the rest of the Carseats data.

```
Carseats <- data.frame(Carseats, High)</pre>
```

We now use the tree() function to fit a classification tree in order to predict High using all variables but Sales. The syntax of the tree() function is quite similar to that of the lm() function.

```
tree.carseats <- tree(High ~ . - Sales, Carseats)
```

The summary() function lists the variables that are used as internal nodes in the tree, the number of terminal nodes, and the (training) error rate.

```
summary(tree.carseats)
```

```
##
## Classification tree:
## tree(formula = High ~ . - Sales, data = Carseats)
## Variables actually used in tree construction:
## [1] "ShelveLoc" "Price" "Income" "CompPrice" "Population"
## [6] "Advertising" "Age" "US"
## Number of terminal nodes: 27
## Residual mean deviance: 0.4575 = 170.7 / 373
## Misclassification error rate: 0.09 = 36 / 400
```

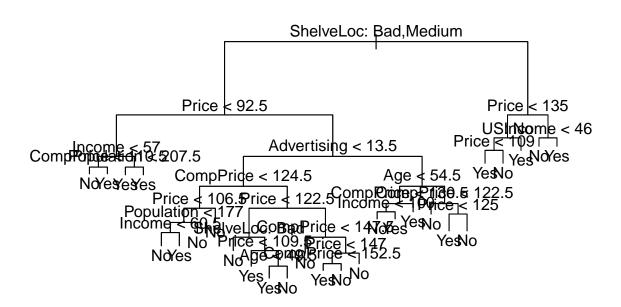
We see that the training error rate is 9%. For classification trees, the deviance reported in the output of summary() is given by

$$-2\sum_{m}\sum_{k}n_{mk}\log\hat{p}_{mk},$$

where  $n_{mk}$  is the number of observations in the *m*th terminal node that belong to the *k*th class. This is closely related to the entropy, defined in (8.7). A small deviance indicates a tree that provides a good fit to the (training) data. The *residual mean deviance* reported is simply the deviance divided by  $n - |T_0|$ , which in this case is 400 - 27 = 373.

One of the most attractive properties of trees is that they can be graphically displayed. We use the plot() function to display the tree structure, and the text() function to display the node labels. The argument pretty = 0 instructs R to include the category names for any qualitative predictors, rather than simply displaying a letter for each category.

```
plot(tree.carseats)
text(tree.carseats, pretty = 0)
```



The most important indicator of Sales appears to be shelving location, since the first branch differentiates Good locations from Bad and Medium locations.

If we just type the name of the tree object, R prints output corresponding to each branch of the tree. R displays the split criterion (e.g. Price < 92.5), the number of observations in that branch, the deviance, the overall prediction for the branch (Yes or No), and the fraction of observations in that branch that take on values of Yes and No. Branches that lead to terminal nodes are indicated using asterisks.

```
## node), split, n, deviance, yval, (yprob)
         * denotes terminal node
##
     1) root 400 541.500 No ( 0.59000 0.41000 )
##
       2) ShelveLoc: Bad, Medium 315 390.600 No (0.68889 0.31111)
##
         4) Price < 92.5 46 56.530 Yes ( 0.30435 0.69565 )
##
##
          8) Income < 57 10 12.220 No ( 0.70000 0.30000 )
##
            16) CompPrice < 110.5 5
                                    0.000 No ( 1.00000 0.00000 ) *
                                     6.730 Yes ( 0.40000 0.60000 ) *
##
            17) CompPrice > 110.5 5
##
           9) Income > 57 36 35.470 Yes (0.19444 0.80556)
##
            18) Population < 207.5 16 21.170 Yes (0.37500 0.62500) *
##
            19) Population > 207.5 20
                                       7.941 Yes ( 0.05000 0.95000 ) *
##
         5) Price > 92.5 269 299.800 No ( 0.75465 0.24535 )
##
          10) Advertising < 13.5 224 213.200 No ( 0.81696 0.18304 )
##
            20) CompPrice < 124.5 96 44.890 No ( 0.93750 0.06250 )
##
              40) Price < 106.5 38 33.150 No ( 0.84211 0.15789 )
##
                80) Population < 177 12 16.300 No ( 0.58333 0.41667 )
##
                 160) Income < 60.5 6 0.000 No (1.00000 0.00000) *
                 161) Income > 60.5 6 5.407 Yes ( 0.16667 0.83333 ) *
##
##
                81) Population > 177 26 8.477 No ( 0.96154 0.03846 ) *
##
              41) Price > 106.5 58
                                   0.000 No ( 1.00000 0.00000 ) *
            21) CompPrice > 124.5 128 150.200 No ( 0.72656 0.27344 )
##
##
              42) Price < 122.5 51 70.680 Yes ( 0.49020 0.50980 )
##
                                       6.702 No ( 0.90909 0.09091 ) *
                84) ShelveLoc: Bad 11
##
                85) ShelveLoc: Medium 40 52.930 Yes (0.37500 0.62500)
                                       7.481 Yes ( 0.06250 0.93750 ) *
##
                 170) Price < 109.5 16
##
                 171) Price > 109.5 24 32.600 No ( 0.58333 0.41667 )
##
                   342) Age < 49.5 13 16.050 Yes ( 0.30769 0.69231 ) *
##
                   343) Age > 49.5 11
                                       6.702 No ( 0.90909 0.09091 ) *
##
              43) Price > 122.5 77 55.540 No ( 0.88312 0.11688 )
##
                86) CompPrice < 147.5 58 17.400 No ( 0.96552 0.03448 ) *
##
                87) CompPrice > 147.5 19 25.010 No ( 0.63158 0.36842 )
##
                 174) Price < 147 12 16.300 Yes ( 0.41667 0.58333 )
##
                   348) CompPrice < 152.5 7
                                              5.742 Yes ( 0.14286 0.85714 ) *
##
                                              5.004 No ( 0.80000 0.20000 ) *
                   349) CompPrice > 152.5 5
##
                 175) Price > 147 7
                                    0.000 No (1.00000 0.00000) *
##
          11) Advertising > 13.5 45 61.830 Yes ( 0.44444 0.55556 )
##
            22) Age < 54.5 25 25.020 Yes ( 0.20000 0.80000 )
##
              44) CompPrice < 130.5 14 18.250 Yes ( 0.35714 0.64286 )
##
                88) Income < 100 9 12.370 No ( 0.55556 0.44444 ) *
##
                                   0.000 Yes ( 0.00000 1.00000 ) *
                89) Income > 100 5
##
              45) CompPrice > 130.5 11
                                       0.000 Yes ( 0.00000 1.00000 ) *
##
            23) Age > 54.5 20 22.490 No ( 0.75000 0.25000 )
##
                                        0.000 No ( 1.00000 0.00000 ) *
              46) CompPrice < 122.5 10
##
              47) CompPrice > 122.5 10 13.860 No ( 0.50000 0.50000 )
##
                                    0.000 Yes ( 0.00000 1.00000 ) *
                94) Price < 125 5
##
                95) Price > 125 5
                                    0.000 No ( 1.00000 0.00000 ) *
##
       3) ShelveLoc: Good 85 90.330 Yes ( 0.22353 0.77647 )
##
         6) Price < 135 68 49.260 Yes (0.11765 0.88235)
##
          12) US: No 17 22.070 Yes ( 0.35294 0.64706 )
##
            24) Price < 109 8  0.000 Yes ( 0.00000 1.00000 ) *
            25) Price > 109 9 11.460 No ( 0.66667 0.33333 ) *
##
```

```
## 13) US: Yes 51 16.880 Yes ( 0.03922 0.96078 ) *
## 7) Price > 135 17 22.070 No ( 0.64706 0.35294 )
## 14) Income < 46 6 0.000 No ( 1.00000 0.00000 ) *
## 15) Income > 46 11 15.160 Yes ( 0.45455 0.54545 ) *
```

In order to properly evaluate the performance of a classification tree on these data, we must estimate the test error rather than simply computing the training error. We split the observations into a training set and a test set, build the tree using the training set, and evaluate its performance on the test data. The predict() function can be used for this purpose. In the case of a classification tree, the argument type = "class" instructs R to return the actual class prediction. This approach leads to correct predictions for around 77% of the locations in the test data set.

```
set.seed(2)
train <- sample(1:nrow(Carseats), 200)</pre>
Carseats.test <- Carseats[-train, ]</pre>
High.test <- High[-train]</pre>
tree.carseats <- tree(High ~ . - Sales, Carseats,</pre>
    subset = train)
tree.pred <- predict(tree.carseats, Carseats.test,</pre>
    type = "class")
table(tree.pred, High.test)
##
             High.test
## tree.pred No Yes
##
          No
             104
                    33
##
          Yes
              13
                    50
```

```
## [1] 0.77
```

(104 + 50) / 200

(If you re-run the predict() function then you might get slightly different results, due to "ties": for instance, this can happen when the training observations corresponding to a terminal node are evenly split between Yes and No response values.)

Next, we consider whether pruning the tree might lead to improved results. The function cv.tree() performs cross-validation in order to determine the optimal level of tree complexity; cost complexity pruning is used in order to select a sequence of trees for consideration. We use the argument FUN = prune.misclass in order to indicate that we want the classification error rate to guide the cross-validation and pruning process, rather than the default for the cv.tree() function, which is deviance. The cv.tree() function reports the number of terminal nodes of each tree considered (size) as well as the corresponding error rate and the value of the cost-complexity parameter used (k, which corresponds to  $\alpha$  in (8.4)).

```
set.seed(7)
cv.carseats <- cv.tree(tree.carseats, FUN = prune.misclass)
names(cv.carseats)

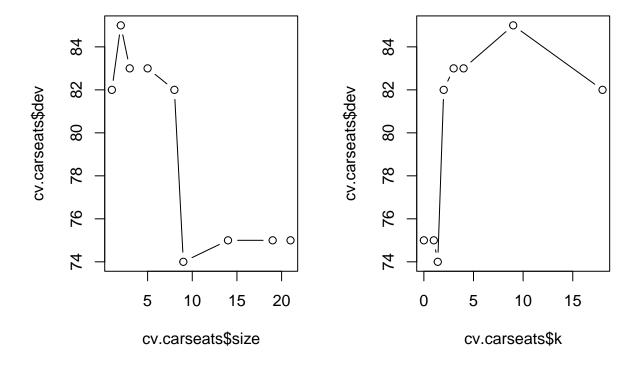
## [1] "size" "dev" "k" "method"

cv.carseats</pre>
```

```
## $size
  [1] 21 19 14 9
                  8 5
                         3
                            2 1
##
## $dev
  [1] 75 75 75 74 82 83 83 85 82
##
##
## $k
            0.0 1.0 1.4 2.0 3.0 4.0 9.0 18.0
## [1] -Inf
##
## $method
  [1] "misclass"
##
## attr(,"class")
## [1] "prune"
                       "tree.sequence"
```

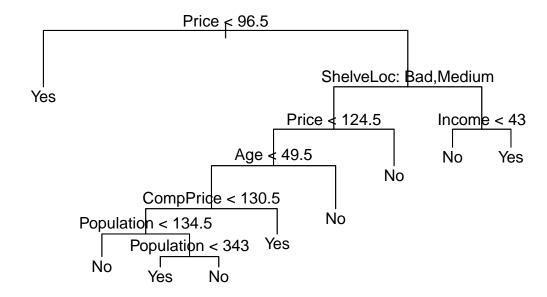
Despite its name, dev corresponds to the number of cross-validation errors. The tree with 9 terminal nodes results in only 74 cross-validation errors. We plot the error rate as a function of both size and k.

```
par(mfrow = c(1, 2))
plot(cv.carseats$size, cv.carseats$dev, type = "b")
plot(cv.carseats$k, cv.carseats$dev, type = "b")
```



We now apply the prune.misclass() function in order to prune the tree to obtain the nine-node tree.

```
prune.carseats <- prune.misclass(tree.carseats, best = 9)
plot(prune.carseats)
text(prune.carseats, pretty = 0)</pre>
```



How well does this pruned tree perform on the test data set? Once again, we apply the predict() function.

```
tree.pred <- predict(prune.carseats, Carseats.test,
    type = "class")
table(tree.pred, High.test)

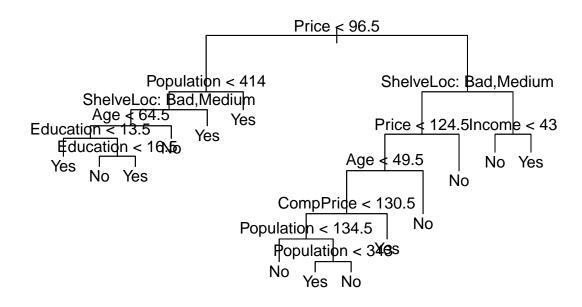
## High.test
## tree.pred No Yes
## No 97 25
## Yes 20 58</pre>
(97 + 58) / 200
```

## [1] 0.775

Now 77.5% of the test observations are correctly classified, so not only has the pruning process produced a more interpretable tree, but it has also slightly improved the classification accuracy.

If we increase the value of best, we obtain a larger pruned tree with lower classification accuracy:

```
prune.carseats <- prune.misclass(tree.carseats, best = 14)
plot(prune.carseats)
text(prune.carseats, pretty = 0)</pre>
```



## [1] 0.77

# Fitting Regression Trees

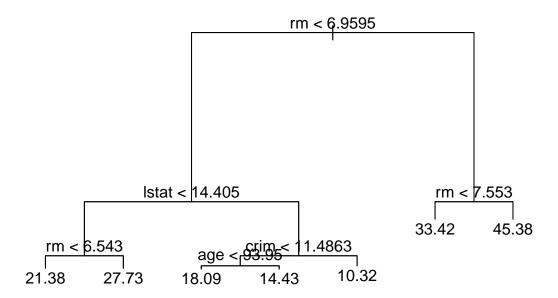
Here we fit a regression tree to the Boston data set. First, we create a training set, and fit the tree to the training data.

```
set.seed(1)
train <- sample(1:nrow(Boston), nrow(Boston) / 2)
tree.boston <- tree(medv ~ ., Boston, subset = train)
summary(tree.boston)</pre>
```

```
##
## Regression tree:
## tree(formula = medv ~ ., data = Boston, subset = train)
## Variables actually used in tree construction:
## [1] "rm"
               "lstat" "crim"
## Number of terminal nodes: 7
## Residual mean deviance: 10.38 = 2555 / 246
## Distribution of residuals:
##
       Min.
             1st Qu.
                       Median
                                        3rd Qu.
                                                     Max.
                                  Mean
## -10.1800 -1.7770
                     -0.1775
                                0.0000
                                          1.9230
                                                  16.5800
```

Notice that the output of summary() indicates that only four of the variables have been used in constructing the tree. In the context of a regression tree, the deviance is simply the sum of squared errors for the tree. We now plot the tree.

```
plot(tree.boston)
text(tree.boston, pretty = 0)
```



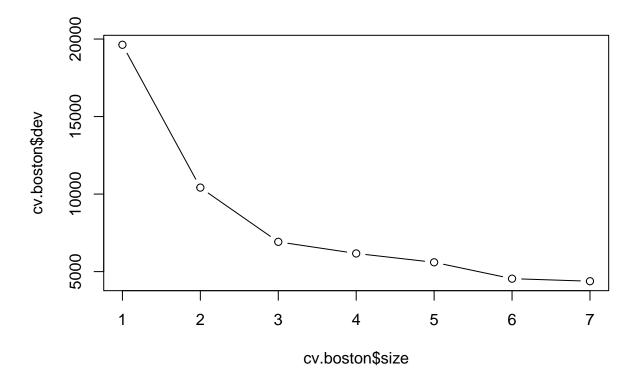
The variable lstat measures the percentage of individuals with {lower socioeconomic status}, while the variable rm corresponds to the average number of rooms. The tree indicates that larger values of rm, or lower

values of lstat, correspond to more expensive houses. For example, the tree predicts a median house price of \$45,400 for homes in census tracts in which rm >= 7.553.

It is worth noting that we could have fit a much bigger tree, by passing control = tree.control(nobs = length(train), mindev = 0) into the tree() function.

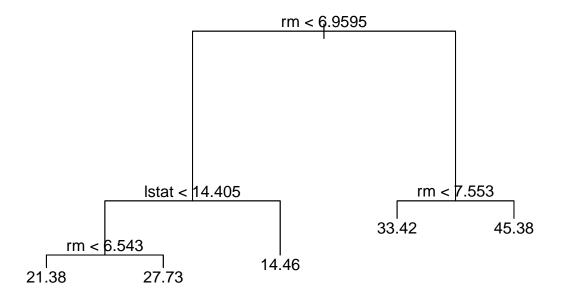
Now we use the cv.tree() function to see whether pruning the tree will improve performance.

```
cv.boston <- cv.tree(tree.boston)
plot(cv.boston$size, cv.boston$dev, type = "b")</pre>
```



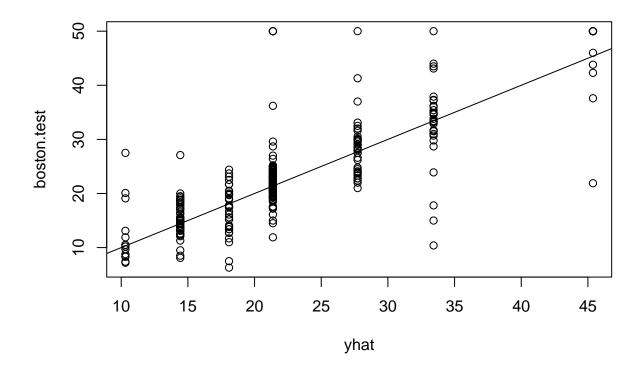
In this case, the most complex tree under consideration is selected by cross-validation. However, if we wish to prune the tree, we could do so as follows, using the prune.tree() function:

```
prune.boston <- prune.tree(tree.boston, best = 5)
plot(prune.boston)
text(prune.boston, pretty = 0)</pre>
```



In keeping with the cross-validation results, we use the unpruned tree to make predictions on the test set.

```
yhat <- predict(tree.boston, newdata = Boston[-train, ])
boston.test <- Boston[-train, "medv"]
plot(yhat, boston.test)
abline(0, 1)</pre>
```



mean((yhat - boston.test)^2)

## [1] 35.28688

In other words, the test set MSE associated with the regression tree is 35.29. The square root of the MSE is therefore around 5.941, indicating that this model leads to test predictions that are (on average) within approximately \$5.941 of the true median home value for the census tract.

# **Bagging and Random Forests**

Here we apply bagging and random forests to the Boston data, using the randomForest package in R. The exact results obtained in this section may depend on the version of R and the version of the randomForest package installed on your computer. Recall that bagging is simply a special case of a random forest with m=p. Therefore, the randomForest() function can be used to perform both random forests and bagging. We perform bagging as follows:

#### library(randomForest)

## randomForest 4.7-1.1

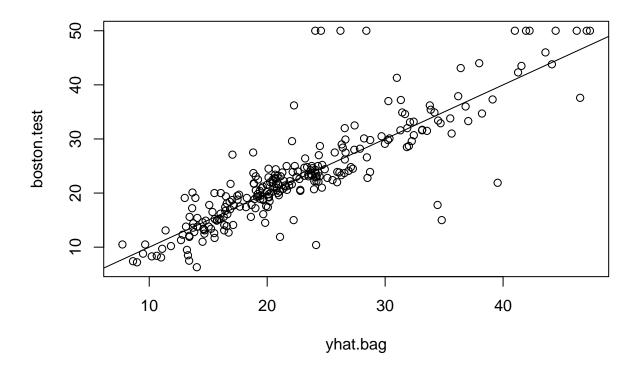
## Type rfNews() to see new features/changes/bug fixes.

```
set.seed(1)
bag.boston <- randomForest(medv ~ ., data = Boston,
    subset = train, mtry = 12, importance = TRUE)
bag.boston</pre>
```

```
##
##
  Call:
##
    randomForest(formula = medv ~ ., data = Boston, mtry = 12, importance = TRUE,
                                                                                          subset = train)
##
                  Type of random forest: regression
                        Number of trees: 500
##
## No. of variables tried at each split: 12
##
##
             Mean of squared residuals: 11.40162
##
                       \% Var explained: 85.17
```

The argument mtry = 12 indicates that all 12 predictors should be considered for each split of the tree—in other words, that bagging should be done. How well does this bagged model perform on the test set?

```
yhat.bag <- predict(bag.boston, newdata = Boston[-train, ])
plot(yhat.bag, boston.test)
abline(0, 1)</pre>
```



```
mean((yhat.bag - boston.test)^2)
```

#### ## [1] 23.41916

The test set MSE associated with the bagged regression tree is 23.42, about two-thirds of that obtained using an optimally-pruned single tree. We could change the number of trees grown by randomForest() using the ntree argument:

```
bag.boston <- randomForest(medv ~ ., data = Boston,
    subset = train, mtry = 12, ntree = 25)
yhat.bag <- predict(bag.boston, newdata = Boston[-train, ])
mean((yhat.bag - boston.test)^2)</pre>
```

#### ## [1] 25.75055

Growing a random forest proceeds in exactly the same way, except that we use a smaller value of the mtry argument. By default, randomForest() uses p/3 variables when building a random forest of regression trees, and  $\sqrt{p}$  variables when building a random forest of classification trees. Here we use mtry = 6.

```
set.seed(1)
rf.boston <- randomForest(medv ~ ., data = Boston,
    subset = train, mtry = 6, importance = TRUE)
yhat.rf <- predict(rf.boston, newdata = Boston[-train, ])
mean((yhat.rf - boston.test)^2)</pre>
```

#### ## [1] 20.06644

The test set MSE is 20.07; this indicates that random forests yielded an improvement over bagging in this case.

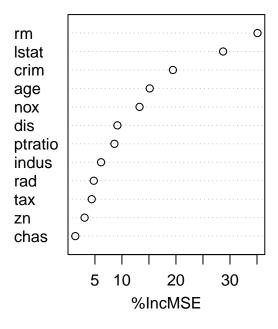
Using the importance() function, we can view the importance of each variable.

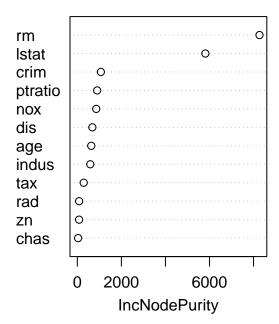
```
importance(rf.boston)
```

```
##
             %IncMSE IncNodePurity
## crim
           19.435587
                         1070.42307
                           82.19257
## zn
            3.091630
## indus
            6.140529
                          590.09536
## chas
            1.370310
                           36.70356
## nox
           13.263466
                          859.97091
                         8270.33906
           35.094741
## rm
           15.144821
                          634.31220
## age
            9.163776
                          684.87953
## dis
            4.793720
## rad
                           83.18719
## tax
            4.410714
                          292.20949
            8.612780
## ptratio
                          902.20190
## lstat
           28.725343
                         5813.04833
```

Two measures of variable importance are reported. The first is based upon the mean decrease of accuracy in predictions on the out of bag samples when a given variable is permuted. The second is a measure of the total decrease in node impurity that results from splits over that variable, averaged over all trees (this was plotted in Figure 8.9). In the case of regression trees, the node impurity is measured by the training RSS, and for classification trees by the deviance. Plots of these importance measures can be produced using the varImpPlot() function.

# rf.boston





The results indicate that across all of the trees considered in the random forest, the wealth of the community (lstat) and the house size (rm) are by far the two most important variables.

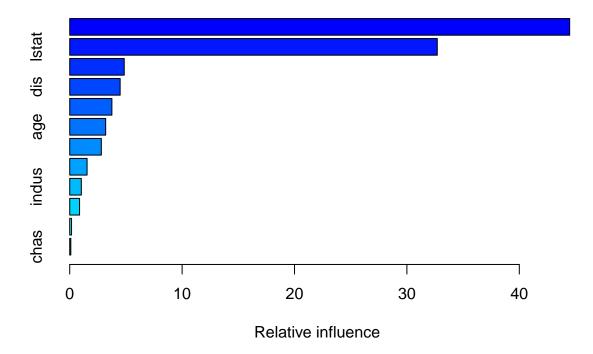
# **Boosting**

Here we use the gbm package, and within it the gbm() function, to fit boosted regression trees to the Boston data set. We run gbm() with the option distribution = "gaussian" since this is a regression problem; if it were a binary classification problem, we would use distribution = "bernoulli". The argument n.trees = 5000 indicates that we want 5000 trees, and the option interaction.depth = 4 limits the depth of each tree.

```
library(gbm)
```

```
## Loaded gbm 2.1.8.1
```

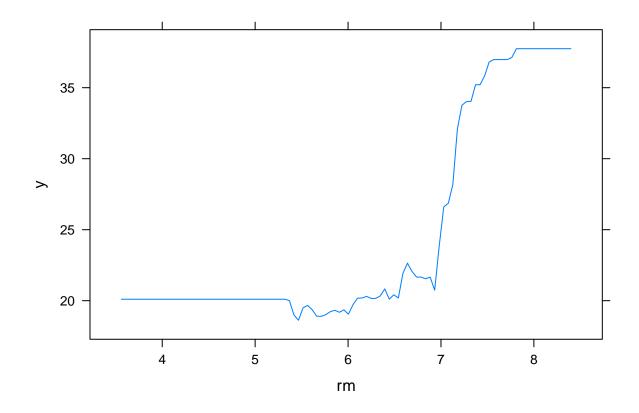
The summary() function produces a relative influence plot and also outputs the relative influence statistics.



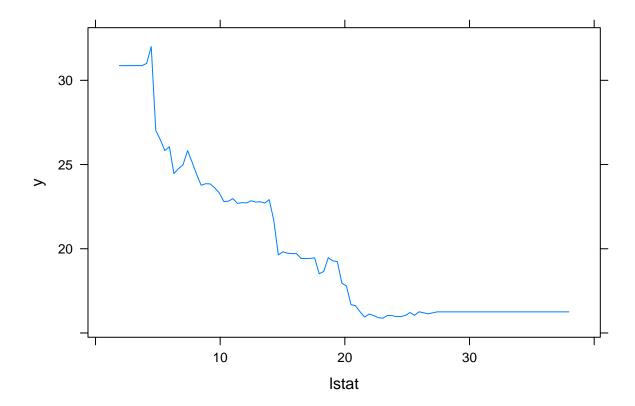
```
##
                var
                        rel.inf
                 rm 44.48249588
## rm
             1stat 32.70281223
## 1stat
## crim
               crim
                     4.85109954
## dis
                dis
                     4.48693083
## nox
                     3.75222394
                nox
## age
                     3.19769210
                age
## ptratio ptratio
                     2.81354826
## tax
                     1.54417603
                tax
## indus
              indus
                     1.03384666
                     0.87625748
## rad
                rad
## zn
                     0.16220479
                 zn
## chas
                     0.09671228
               chas
```

We see that lstat and rm are by far the most important variables. We can also produce partial dependence plots for these two variables. These plots illustrate the marginal effect of the selected variables on the response after integrating out the other variables. In this case, as we might expect, median house prices are increasing with rm and decreasing with lstat.

```
plot(boost.boston, i = "rm")
```



plot(boost.boston, i = "lstat")



We now use the boosted model to predict med on the test set:

```
yhat.boost <- predict(boost.boston,
    newdata = Boston[-train, ], n.trees = 5000)
mean((yhat.boost - boston.test)^2)</pre>
```

### ## [1] 18.39057

The test MSE obtained is 18.39: this is superior to the test MSE of random forests and bagging. If we want to, we can perform boosting with a different value of the shrinkage parameter  $\lambda$  in (8.10). The default value is 0.001, but this is easily modified. Here we take  $\lambda = 0.2$ .

### ## [1] 16.54778

In this case, using  $\lambda = 0.2$  leads to a lower test MSE than  $\lambda = 0.001$ .

# Bayesian Additive Regression Trees

## done 1000 (out of 1100)

In this section we use the BART package, and within it the gbart() function, to fit a Bayesian additive regression tree model to the Boston housing data set. The gbart() function is designed for quantitative outcome variables. For binary outcomes, lbart() and pbart() are available.

To run the gbart() function, we must first create matrices of predictors for the training and test data. We run BART with default settings.

```
library(BART)
## Loading required package: nlme
## Loading required package: nnet
## Loading required package: survival
x <- Boston[, 1:12]
y <- Boston[, "medv"]
xtrain <- x[train, ]</pre>
ytrain <- y[train]</pre>
xtest <- x[-train, ]</pre>
ytest <- y[-train]</pre>
set.seed(1)
bartfit <- gbart(xtrain, ytrain, x.test = xtest)</pre>
## *****Calling gbart: type=1
## ****Data:
## data:n,p,np: 253, 12, 253
## y1,yn: 0.213439, -5.486561
## x1,x[n*p]: 0.109590, 20.080000
## xp1,xp[np*p]: 0.027310, 7.880000
## *****Number of Trees: 200
## *****Number of Cut Points: 100 ... 100
## ****burn,nd,thin: 100,1000,1
## ****Prior:beta,alpha,tau,nu,lambda,offset: 2,0.95,0.795495,3,3.71636,21.7866
## ****sigma: 4.367914
## ****w (weights): 1.000000 ... 1.000000
## *****Dirichlet:sparse,theta,omega,a,b,rho,augment: 0,0,1,0.5,1,12,0
## ****printevery: 100
##
## MCMC
## done 0 (out of 1100)
## done 100 (out of 1100)
## done 200 (out of 1100)
## done 300 (out of 1100)
## done 400 (out of 1100)
## done 500 (out of 1100)
## done 600 (out of 1100)
## done 700 (out of 1100)
## done 800 (out of 1100)
## done 900 (out of 1100)
```

```
## time: 2s
```

## trcnt,tecnt: 1000,1000

Next we compute the test error.

```
yhat.bart <- bartfit$yhat.test.mean
mean((ytest - yhat.bart)^2)</pre>
```

```
## [1] 15.91912
```

On this data set, the test error of BART is lower than the test error of random forests and boosting. Now we can check how many times each variable appeared in the collection of trees.

```
ord <- order(bartfit$varcount.mean, decreasing = T)
bartfit$varcount.mean[ord]</pre>
```

```
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          21.653 21.638 20.725 20.021 19.615 19.283 19.278 19.073 15.576
##
   22.973
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             crim
##
  13.800 11.607
```