

Preliminaries

```
> library(DAAG)
> library(rpart)
```

Exercise 1

Refer to the `head.injury` data frame.

- (a) Use the default setting in `rpart()` to obtain a tree-based model for predicting occurrence of clinically important brain injury, given the other variables.
- (b) How many splits gives the minimum cross-validation error? Prune the tree using the 1 standard error rule.

```
(a) > set.seed(29)
> injury.rpart <- rpart(clinically.important.brain.injury ~ .,
+   data = head.injury, method = "class", cp = 1e-04)
> plotcp(injury.rpart)
> printcp(injury.rpart)
```

Classification tree:

```
rpart(formula = clinically.important.brain.injury ~ ., data = head.injury,
      method = "class", cp = 1e-04)
```

Variables actually used in tree construction:

```
[1] GCS.13          GCS.15.2hours    age.65
[4] amnesia.before  basal.skull.fracture high.risk
[7] loss.of.consciousness vomiting
```

Root node error: 250/3121 = 0.080103

n= 3121

	CP	nsplit	rel error	xerror	xstd
1	0.0400	0	1.000	1.000	0.060660
2	0.0360	2	0.920	0.992	0.060438
3	0.0140	3	0.884	0.896	0.057678
4	0.0080	5	0.856	0.904	0.057915
5	0.0001	10	0.816	0.916	0.058268

The setting `cp=0.0001` was reached after some experimentation.

- (b) The minimum cross-validated relative error is for `nsplit=3`, i.e., for a tree size of 4.
- (c) The one-standard-error rule likewise chooses `nsplit=3`, with `cp=0.014`. Setting `cp=0.02`, i.e., larger than `cp` for the next smallest number of splits, will prune the tree back to this size. We have

```
> injury0.rpart <- prune(injury.rpart, cp = 0.02)
```

We plot the tree from (a) that shows the cross-validated relative error, and the tree obtained from (c).

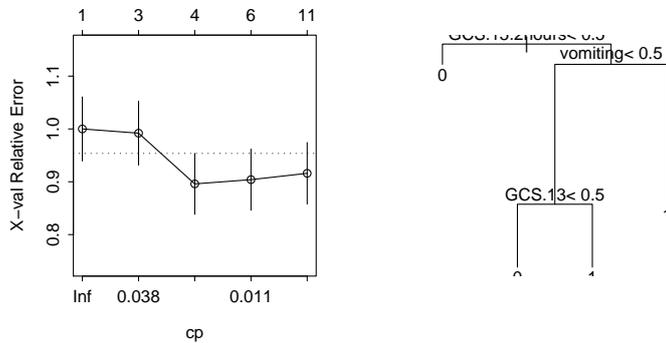


Figure 1: Plots from the `rpart` analysis of the head injury data: (i) cross-validated relative error versus `cp`; and (ii) the tree obtained in (c).

There can be substantial change from one run to the next.

Exercise 2

The data set `mifem` is part of the larger data set in the data frame `monica` that we have included in our `DAAG` package. Use tree-based regression to predict mortality in this larger data set. What is the most immediately striking feature of your analysis? Should this be a surprise?

```
> monica.rpart <- rpart(outcome ~ ., data = monica, method = "class")
```

```
> plot(monica.rpart)
> text(monica.rpart)
```



Figure 2: Classification tree for `monica` data.

Those who were not hospitalised were very likely to be dead! Check by examining the table:

```
> table(monica$hosp, monica$outcome)
```

```
   live dead
y 3522  920
n    3 1922
```

Exercise 3

Use tree-based regression to predict `re78` in the data frame `nsw74pred1` that is in our `DAAG` package. Compare the predictions with the multiple regression predictions in Chapter 6.

In order to reproduce the same results as given here, do:

```
> set.seed(21)
```

Code for the initial calculation is:

```
> nsw.rpart <- rpart(re78 ~ ., data = nsw74psid1, cp = 0.001)
> plotcp(nsw.rpart)
```

It is obvious that `cp=0.002` will be adequate. At this point, the following is a matter of convenience, to reduce the printed output:

```
> nsw.rpart <- prune(nsw.rpart, cp = 0.002)
> printcp(nsw.rpart)
```

Regression tree:

```
rpart(formula = re78 ~ ., data = nsw74psid1, cp = 0.001)
```

Variables actually used in tree construction:

```
[1] age educ re74 re75
```

Root node error: $6.5346e+11/2675 = 244284318$

n= 2675

	CP	nsplit	rel error	xerror	xstd
1	0.3446296	0	1.00000	1.00067	0.046287
2	0.1100855	1	0.65537	0.66461	0.038977
3	0.0409403	2	0.54528	0.55811	0.033004
4	0.0317768	3	0.50434	0.51821	0.035244
5	0.0158188	4	0.47257	0.50636	0.034622
6	0.0105727	5	0.45675	0.49139	0.034688
7	0.0105337	6	0.44618	0.48453	0.034527
8	0.0063341	7	0.43564	0.46901	0.032502
9	0.0056603	8	0.42931	0.46028	0.032969
10	0.0038839	9	0.42365	0.46133	0.033142
11	0.0035516	10	0.41976	0.46238	0.033095
12	0.0031768	11	0.41621	0.47329	0.033838
13	0.0028300	12	0.41304	0.47544	0.033675
14	0.0027221	13	0.41021	0.47495	0.033776
15	0.0023286	15	0.40476	0.47570	0.033783
16	0.0020199	16	0.40243	0.47642	0.033609
17	0.0020000	17	0.40041	0.47715	0.033851

The minimum cross-validated relative error is at `nsplit=12`. The one standard error limit is 0.498 ($=0.463+0.035$). The one standard error rule suggests taking `nsplit=5`.

If we go with the one standard error rule, we have a residual variance equal to $244284318 \times 0.49177 = 120131699$.

For the estimate of residual variance from the calculations of Section 6.x, we do the following.

```

> attach(nsw74psid1)
> here <- age <= 40 & re74 <= 5000 & re75 <= 5000 & re78 < 30000
> nsw74psidA <- nsw74psid1[here, ]
> detach(nsw74psid1)
> A1.lm <- lm(re78 ~ trt + (age + educ + re74 + re75) + (black +
+ hisp + marr + nodeg), data = nsw74psidA)
> summary(A1.lm)$sigma^2

```

```
[1] 40177577
```

The variance estimate is 40177577. This is about a third of the variance estimate that was obtained with tree-based regression.

Exercise 4

Copy down the email spam data set from the web site given in Section 10.2. Carry out a tree-based regression using all 57 available explanatory variables. Determine the change in the cross-validation estimate of predictive accuracy.

We set the random number seed to 21, to allow users to reproduce our results. In most other contexts, it will be best not to set a seed. The file `spam.shortnames` is available for copying from the web address <http://wwwmaths.anu.edu.au/~johnm/r-book/xtra-data>. The data frame `spam` is created thus:

```

> spam <- read.table("spambase.data", header = FALSE, sep = ",")
> nam <- scan("spam.shortnames", what = "")
> names(spam) <- nam

```

Now load `rpart` and proceed with the calculations.

```

> set.seed(21)
> spam.rpart <- rpart(yesno ~ ., data = spam, cp = 1e-04, method = "class")
> printcp(spam.rpart)

```

Classification tree:

```
rpart(formula = yesno ~ ., data = spam, method = "class", cp = 1e-04)
```

Variables actually used in tree construction:

```

[1] address    bang        crl.av      crl.long    crl.tot     data
[7] dollar     edu         email       font        free        george
[13] hp         internet   leftparen   money       n1999      n650
[19] our        over       re          remove     semicolon   technology
[25] will       you        your

```

Root node error: 1813/4601 = 0.39404

n= 4601

	CP	nsplit	rel error	xerror	xstd
1	0.47655819	0	1.00000	1.00000	0.0182819
2	0.14892443	1	0.52344	0.55819	0.0154972
3	0.04302261	2	0.37452	0.46001	0.0144131
4	0.03088803	4	0.28847	0.32212	0.0124547
5	0.01047987	5	0.25758	0.27910	0.0117052

6	0.00827358	6	0.24710	0.26586	0.0114576
7	0.00717044	7	0.23883	0.26089	0.0113626
8	0.00529509	8	0.23166	0.25317	0.0112121
9	0.00441258	14	0.19581	0.23552	0.0108559
10	0.00358522	15	0.19140	0.22835	0.0107060
11	0.00275786	19	0.17705	0.22559	0.0106475
12	0.00257400	22	0.16878	0.22228	0.0105767
13	0.00220629	25	0.16106	0.22228	0.0105767
14	0.00211436	27	0.15665	0.21897	0.0105052
15	0.00165472	33	0.14396	0.21180	0.0103477
16	0.00110314	36	0.13900	0.19857	0.0100476
17	0.00082736	43	0.13127	0.19581	0.0099834
18	0.00055157	47	0.12796	0.19581	0.0099834
19	0.00036771	53	0.12466	0.20077	0.0100985
20	0.00010000	62	0.12135	0.20077	0.0100985

Figure 3 shows the graph that is obtained by plotting this tree. For making a decision on the size of tree however, it is convenient to work from the information given by the function `printcp()`.

```
> plotcp(spam.rpart)
```

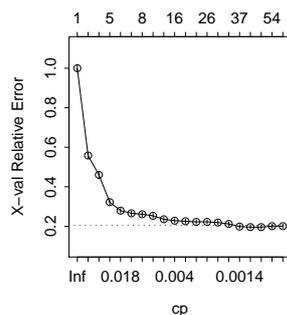


Figure 3: Plot of cross-validated relative error versus `cp`, for the full `spam` data set.

Setting `cp=0.0001` ensures, when the random number seed is set to 21, that the cross-validated relative error reaches a minimum, of 0.1958, at `nsplit=43`. Pruning to get the tree that is likely to have best predictive power can use `cp=0.001`. Adding the SE to the minimum cross-validated relative error gives 0.2. The smallest tree with an SE smaller than this is at `nsplit=36`; setting `cp=0.0012` will give this tree.

Here then are the two prunings:

```
> spam.rpart1 <- prune(spam.rpart, cp = 0.001)
> spam.rpart2 <- prune(spam.rpart, cp = 0.0012)
```

Additional Exercises A number of additional exercises are included in the laboratory exercises that are available from the web page <http://www.maths.anu.edu.au/~johnm/courses/dm>