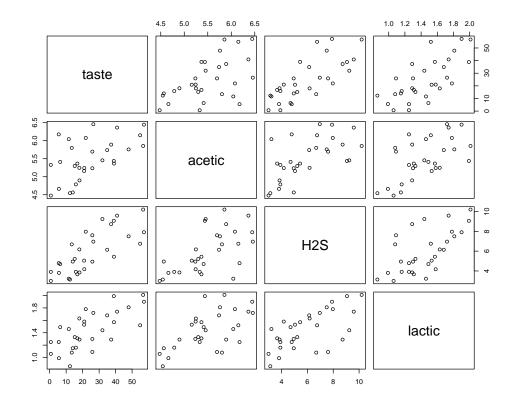
Solutions for 7.5 Exercises

1. Cheese

(a) > cheese <- read.csv("cheese.csv")
> plot(cheese)



(b) > taste.lm1 <- lm(taste ~ acetic, data = cheese)
> summary(taste.lm1)

Call:

lm(formula = taste ~ acetic, data = cheese)

Residuals:

Min 1Q Median 3Q Max -29.642 -7.443 2.082 6.597 26.581

Coefficients:

Estimate Std. Error t value Pr(>|t|)
(Intercept) -61.499 24.846 -2.475 0.01964 *
acetic 15.648 4.496 3.481 0.00166 **

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Residual standard error: 13.82 on 28 degrees of freedom Multiple R-squared: 0.302, Adjusted R-squared: 0.2771 F-statistic: 12.11 on 1 and 28 DF, p-value: 0.001658

```
> taste.lm2 <- lm(taste ~ H2S, data = cheese)</pre>
   > summary(taste.lm2)
   Call:
   lm(formula = taste ~ H2S, data = cheese)
   Residuals:
       Min
                1Q Median
                                 3Q
                                        Max
   -15.426 -7.611 -3.491
                              6.420
                                     25.687
   Coefficients:
               Estimate Std. Error t value Pr(>|t|)
   (Intercept) -9.7868
                             5.9579 -1.643
                                               0.112
                                      6.107 1.37e-06 ***
   H2S
                 5.7761
                             0.9458
   Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1
   Residual standard error: 10.83 on 28 degrees of freedom
   Multiple R-squared: 0.5712, Adjusted R-squared: 0.5558
   F-statistic: 37.29 on 1 and 28 DF, p-value: 1.374e-06
   > taste.lm3 <- lm(taste ~ lactic, data = cheese)</pre>
   > summary(taste.lm3)
   Call:
   lm(formula = taste ~ lactic, data = cheese)
   Residuals:
                       Median
        Min
                   1Q
                                     3Q
                                             Max
   -19.9439 -8.6839 -0.1095
                                 8.9998 27.4245
   Coefficients:
               Estimate Std. Error t value Pr(>|t|)
                             10.582 -2.822 0.00869 **
   (Intercept)
                -29.859
   lactic
                 37.720
                              7.186
                                      5.249 1.41e-05 ***
   Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1
   Residual standard error: 11.75 on 28 degrees of freedom
   Multiple R-squared: 0.4959, Adjusted R-squared: 0.4779
   F-statistic: 27.55 on 1 and 28 DF, p-value: 1.405e-05
   From these results, we conclude that there is a statistically significant association (at the
   0.05 significance level) between taste and each of the three chemical levels.
(c) > taste.lm4 <- lm(taste ~ acetic + H2S + lactic, data = cheese)
   > summary(taste.lm4)
   lm(formula = taste ~ acetic + H2S + lactic, data = cheese)
```

Residuals:

```
Min 1Q Median 3Q Max -17.391 -6.612 -1.009 4.908 25.449
```

Coefficients:

```
Estimate Std. Error t value Pr(>|t|)
(Intercept) -28.8768
                        19.7354
                                 -1.463
                                         0.15540
acetic
              0.3277
                         4.4598
                                  0.073 0.94198
H2S
              3.9118
                         1.2484
                                  3.133 0.00425 **
lactic
             19.6705
                         8.6291
                                  2.280 0.03108 *
```

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Residual standard error: 10.13 on 26 degrees of freedom Multiple R-squared: 0.6518, Adjusted R-squared: 0.6116 F-statistic: 16.22 on 3 and 26 DF, p-value: 3.81e-06

The multiple regression is highly significant and accounts for 65% of the variability in taste. However, not all of the terms are needed. In particular, the P-value for acetic acid is 0.94 which implies that it should be removed from the model.

```
> taste.lm5 <- lm(taste ~ H2S + lactic, data = cheese)
> summary(taste.lm5)
```

Call:

lm(formula = taste ~ H2S + lactic, data = cheese)

Residuals:

```
Min 1Q Median 3Q Max -17.343 -6.530 -1.164 4.844 25.618
```

Coefficients:

```
Estimate Std. Error t value Pr(>|t|)
(Intercept) -27.592 8.982 -3.072 0.00481 **
H2S 3.946 1.136 3.475 0.00174 **
lactic 19.887 7.959 2.499 0.01885 *
```

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Residual standard error: 9.942 on 27 degrees of freedom Multiple R-squared: 0.6517, Adjusted R-squared: 0.6259 F-statistic: 25.26 on 2 and 27 DF, p-value: 6.551e-07

Removing acetic acid from the model results in a lower residual standard error and a higher adjusted R^2 . Both lactic acid and H_2S are still significant in this new model. It is the best model.

(Using the step() function in R should give the same result — check this.)

(d) > cor(cheese)

```
taste acetic H2S lactic taste 1.0000000 0.5495393 0.7557523 0.7042362
```

```
acetic 0.5495393 1.0000000 0.6179559 0.6037826

H2S 0.7557523 0.6179559 1.0000000 0.6448123

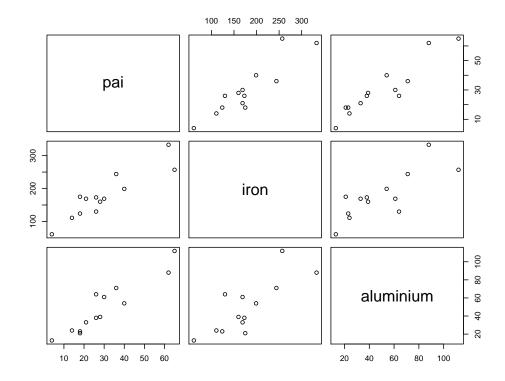
lactic 0.7042362 0.6037826 0.6448123 1.0000000
```

The relatively high correlations between taste and each of the explanatory variables are consistent with the findings in part (b). There is also high correlation amongst the explanatory variables, which explains why not all three were required in the model in part (c). Given the values of H_2S and lactic acid, we can obtain quite a reasonable estimate of the value of acetic acid The concentration of acetic acid has little additional information about taste, beyond that of H_2S and lactic acid.

(e) Taste is predicted to increase by 3.95 units for each additional unit of H₂S concentration, provided lactic acid concentration remains constant.

2. Soil Data

```
(a) > soil < -data.frame(pai = c(4,18,14,18,26,26,21,30,28,36,65,62,40), + iron = c(61,175,111,124,130,173,169,169,160,244,257,333,199), + aluminium = c(13,21,24,23,64,38,33,61,39,71,112,88,54)) <math>> plot(soil)
```



The relationships between the response pai and both variables iron and aluminium appears quite linear, as does the relationship between iron and aluminium.

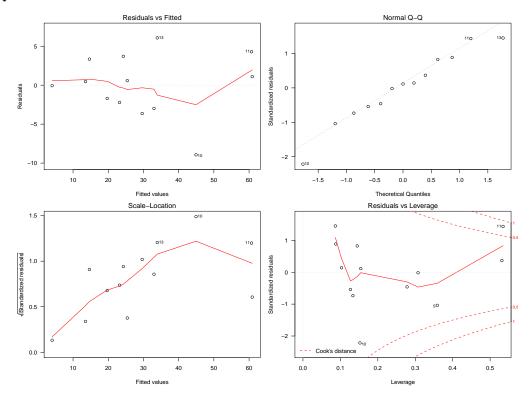
```
(b) > soil.lm.iron <- lm(pai ~ iron, data = soil)
> summary(soil.lm.iron)
```

```
Call:
   lm(formula = pai ~ iron, data = soil)
   Residuals:
        Min
                  1Q
                      Median
                                   3Q
                                           Max
                     0.3174
   -11.3196 -3.3714
                              2.1030 16.9700
   Coefficients:
                Estimate Std. Error t value Pr(>|t|)
   (Intercept) -10.61117
                           6.00940 -1.766
                           0.03168 7.202 1.75e-05 ***
   iron
                 0.22818
   Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1
   Residual standard error: 7.694 on 11 degrees of freedom
   Multiple R-squared: 0.825,
                                 Adjusted R-squared: 0.8091
   F-statistic: 51.86 on 1 and 11 DF, p-value: 1.749e-05
   > soil.lm.1 <- lm(pai ~ iron + aluminium, data = soil)</pre>
   > summary(soil.lm.1)
   Call:
   lm(formula = pai ~ iron + aluminium, data = soil)
   Residuals:
       Min
                1Q Median
                                3Q
   -8.9352 -2.2182 0.4613 3.3448 6.0708
   Coefficients:
               Estimate Std. Error t value Pr(>|t|)
   (Intercept) -7.35066
                          3.48467 -2.109 0.061101 .
                0.11273
                          0.02969
                                   3.797 0.003504 **
   iron
   aluminium
                0.34900
                          0.07131
                                   4.894 0.000628 ***
   Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1
   Residual standard error: 4.379 on 10 degrees of freedom
  Multiple R-squared: 0.9485,
                                     Adjusted R-squared: 0.9382
   F-statistic: 92.03 on 2 and 10 DF, p-value: 3.634e-07
(c) > predict(soil.lm.iron, newdata = data.frame(iron = 150), interval = "prediction")
         fit
                  lwr
                           upr
   1 23.6152 5.938094 41.29230
   > predict(soil.lm.1, newdata = data.frame(iron = 150,
            aluminium=mean(soil$aluminium)), interval = "prediction")
          fit
                   lwr
   1 26.76768 16.48159 37.05377
```

The predicted value is similar, but the prediction interval is wider, which is not surprising for a model which doesn't fit as well.

(d) >
$$par(mfrow = c(2, 2), las = 1, mar = c(4, 4, 2, 1))$$

> $plot(soil.lm.1)$



There is some evidence that the variance increases with the fitted values, but with only 13 points it is difficult to be sure. The Cook's distance plot suggests that point 11 has a substantial impact on the model, but it looks OK on the other plots.

```
(e) > step(soil.lm.1, ~.)
Start: AIC=40.99
```

pai ~ iron + aluminium

Df Sum of Sq RSS AIC
<none> 191.79 40.989
- iron 1 276.49 468.28 50.594
- aluminium 1 459.43 651.22 54.881

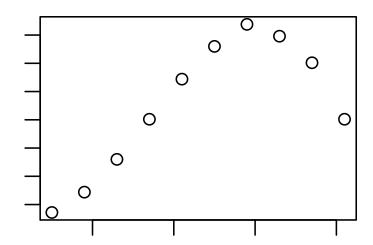
Call:

lm(formula = pai ~ iron + aluminium, data = soil)

Coefficients:

(Intercept) iron aluminium -7.3507 0.1127 0.3490 Omission of either explanatory variable substantially increases AIC, so both need to remain in the model.

3. Girths of horses



> cor(girth)

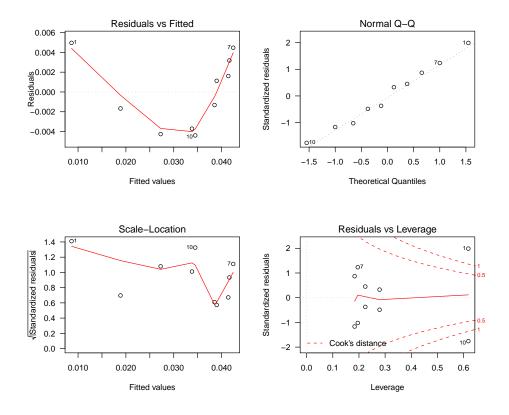
weight stretch weight 1.0000000 0.7404735 stretch 0.7404735 1.0000000

r=0.740. This is of limited usefulness here, because it measures the strength of the linear association, and the relationship is not linear.

```
weight 6.725e-03 1.035e-03 6.496 0.000336 ***
I(weight^2) -2.299e-04 4.396e-05 -5.230 0.001212 **
---
Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Residual standard error: 0.00404 on 7 degrees of freedom
Multiple R-squared: 0.908, Adjusted R-squared: 0.8817
F-statistic: 34.53 on 2 and 7 DF, p-value: 0.0002365

> par(mfrow = c(2, 2), las = 1)
> plot(girth.lm.2)
```

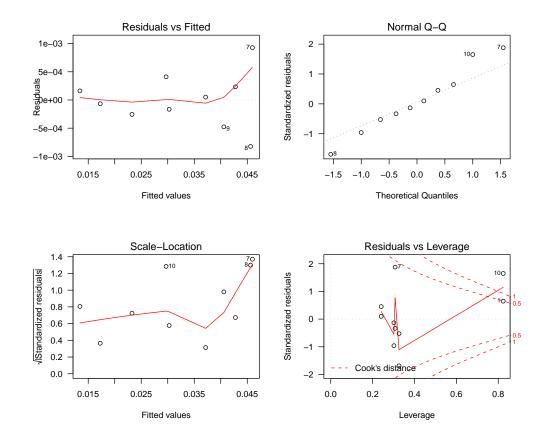


The quadratic term is highly significant, but the residuals suggest a cubic term is needed.

Coefficients:

```
Estimate Std. Error t value Pr(>|t|)
(Intercept)
            1.344e-02
                        1.366e-03
                                    9.833 6.37e-05 ***
                        4.765e-04
weight
            -1.330e-03
                                   -2.791
                                            0.0315 *
                        4.653e-05
                                   12.721 1.45e-05 ***
I(weight^2) 5.919e-04
                        1.336e-06 -17.836 2.00e-06 ***
I(weight^3) -2.382e-05
Signif. codes:
                0 *** 0.001 ** 0.01 * 0.05 . 0.1 1
Residual standard error: 0.0005938 on 6 degrees of freedom
Multiple R-squared: 0.9983, Adjusted R-squared: 0.9974
F-statistic: 1172 on 3 and 6 DF, p-value: 1.081e-08
```

> par(mfrow = c(2, 2), las = 1) > plot(girth.lm.3)

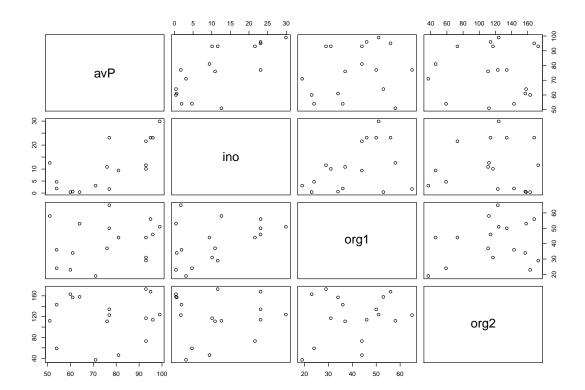


The cubic term makes a further significant improvement, and the residuals don't suggest a 4th power. The fit is amazingly good (too good?), judging from the \mathbb{R}^2 and the residual standard deviation.

(c) > predict(girth.lm.3, newdata = data.frame(weight = 10.5), interval = "confidence") fit lwr upr 1 0.03714918 0.03643587 0.0378625 Residual = 0.0372 - 0.03715 = 0.00005.

4. Phosphorus content in corn.

(a) > Pcorn <- read.csv("Pcorn.csv")
> plot(Pcorn)



> cor(Pcorn)

```
    avP
    ino
    org1
    org2

    avP
    1.00000000
    0.720086602
    0.2118376
    0.029934165

    ino
    0.72008660
    1.000000000
    0.3989231
    -0.006425396

    org1
    0.21183758
    0.398923129
    1.0000000
    0.222479243

    org2
    0.02993417
    -0.006425396
    0.2224792
    1.000000000
```

Available P seems to have a reasonably strong relationship with inorganic P, but little relationship with the two organic P measurements.

```
(b) > Pcorn.lm <- lm(avP ~ ino + org1 + org2, data = Pcorn)
   > step(Pcorn.lm, ~.)
   Start: AIC=89.78
   avP ~ ino + org1 + org2
          Df Sum of Sq
                          RSS
                                  AIC
   - org2 1
                 14.12 2101.3 87.891
                 38.78 2125.9 88.089
   - org1
          1
                        2087.2 89.776
   <none>
   - ino
               2139.28 4226.5 99.770
           1
```

```
Step: AIC=87.89
avP ~ ino + org1
      Df Sum of Sq RSS
         29.95 2131.2 86.131
- org1 1
<none>
                  2101.3 87.891
+ org2 1
           14.12 2087.2 89.776
- ino 1 2126.54 4227.8 97.776
Step: AIC=86.13
avP ~ ino
      Df Sum of Sq RSS
<none>
                  2131.2 86.131
            29.95 2101.3 87.891
+ org1 1
+ org2 1
            5.29 2125.9 88.089
- ino 1
           2295.23 4426.5 96.556
Call:
lm(formula = avP ~ ino, data = Pcorn)
Coefficients:
(Intercept)
                   ino
    62.569
                 1.229
> Pcorn.lm1 <- lm(avP ~ 1, data = Pcorn)</pre>
> step(Pcorn.lm1, ~. + ino + org1 + org2)
Start: AIC=96.56
avP ~ 1
      Df Sum of Sq RSS
                            AIC
+ ino 1 2295.23 2131.2 86.131
<none>
                  4426.5 96.556
          198.64 4227.8 97.776
+ org1 1
+ org2 1 3.97 4422.5 98.541
Step: AIC=86.13
avP ~ ino
      Df Sum of Sq
                   RSS
                            AIC
<none>
                  2131.2 86.131
+ org1 1
             29.95 2101.3 87.891
+ org2 1
            5.29 2125.9 88.089
- ino 1 2295.23 4426.5 96.556
```

Call:

11

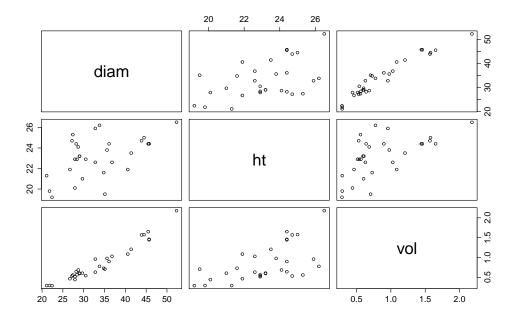


Figure 1: > plot(timber)

```
lm(formula = avP ~ ino, data = Pcorn)
Coefficients:
(Intercept) ino
62.569 1.229
```

For either backward elimination or forward selection, the smallest AIC is for the model with only inorganic P, confirming the impressions from the graphs.

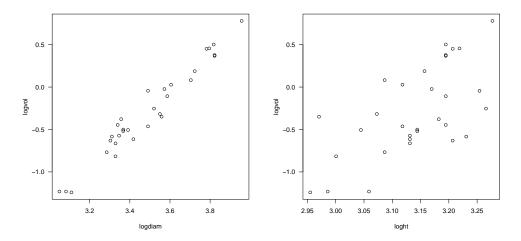
5. Estimating timber volume

```
> timber <- read.csv("timber.csv")
> plot(timber)
```

From looking at the graphs (and based on the geometry of the problem), taking logs of each variable might be a good place to start.

```
> timber$logvol = log(timber$vol)
> timber$logdiam = log(timber$diam)
> timber$loght = log(timber$ht)

> par(mfrow = c(1, 2), las = 1, mar = c(4, 6, 2, 1))
> plot(logvol ~ logdiam, data = timber)
> plot(logvol ~ loght, data = timber)
```



log(volume) vs log(diameter) looks very promising.

log(volume) vs log(height) has more noise, but also looks linear.

Call:

lm(formula = logvol ~ logdiam + loght, data = timber)

Residuals:

Min 1Q Median 3Q Max -0.171144 -0.048326 0.006864 0.061369 0.128931

Coefficients:

Signif. codes: 0 *** 0.001 ** 0.01 * 0.05 . 0.1 1

Residual standard error: 0.08148 on 28 degrees of freedom Multiple R-squared: 0.9776, Adjusted R-squared: 0.976 F-statistic: 611.3 on 2 and 28 DF, p-value: < 2.2e-16

 $R^2 = 0.978$, which is larger than the R^2 for models fitted without transforming.