

Midterm Comments

Statistics 135

Autumn 2005



Midterm Comments

1. Puffin nesting frequency

$$\text{nesting}_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \beta_3 \text{angle}_i + \beta_4 \text{distance}_i + \epsilon_i$$

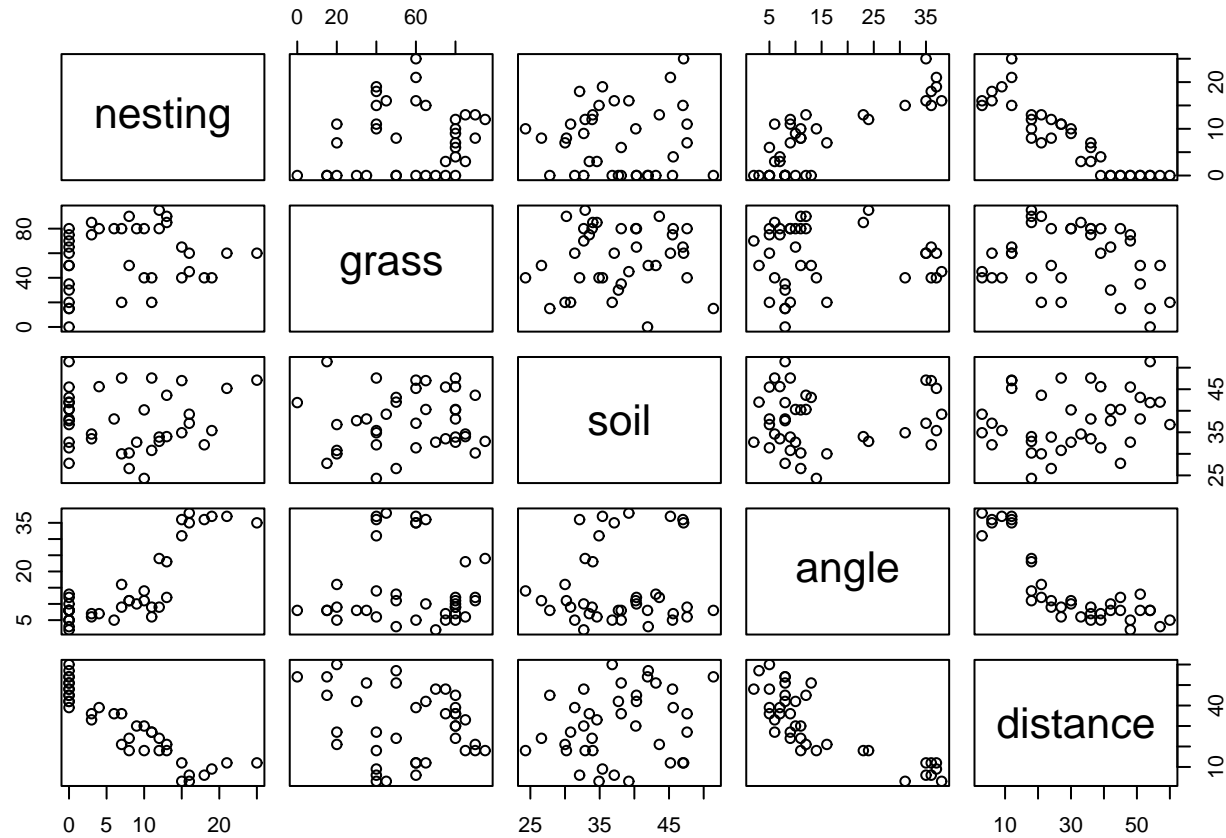
- (a) (5 points) The data from this study is available on the datasets page in the file `puffin.txt`. Read in the data, calculate the standard summary statistics (mean, standard deviation, and 5 figure summary) for each of the variables and create a scatter plot matrix of the data. Does it appear that any of the potential predictor variables are associated with nesting frequency?

```

> summary(puffin)
      nesting      grass      soil      angle
Min.   : 0.000   Min.   : 0.00   Min.   :24.30   Min.   : 2.00
1st Qu.: 0.000   1st Qu.:40.00   1st Qu.:32.75   1st Qu.: 7.25
Median : 7.500   Median :60.00   Median :37.40   Median :10.00
Mean   : 7.684   Mean   :56.45   Mean   :37.72   Mean   :15.00
3rd Qu.:12.750   3rd Qu.:80.00   3rd Qu.:42.83   3rd Qu.:21.25
Max.   :25.000   Max.   :95.00   Max.   :51.40   Max.   :38.00

      distance
Min.   : 3.00
1st Qu.:18.00
Median :30.00
Mean   :30.39
3rd Qu.:44.25
Max.   :60.00
> sapply(puffin, function(x) sqrt(var(x)))
      nesting      grass      soil      angle      distance
7.192734 25.306443  6.653978 11.691993 16.581558
> sqrt(diag(var(puffin)))
      nesting      grass      soil      angle      distance
7.192734 25.306443  6.653978 11.691993 16.581558

```



```
plot(puffin) # splom(~puffin) is fine as well
```

It appears that angle and distance are associated with nesting and possibly soil as well. Note that there are correlations between some of the predictors, particularly between angle and distance.

While not asked for, the correlation matrix for all the variables is

```
> cor(puffin)
      nesting      grass      soil      angle      distance
nesting 1.00000000 0.15848477 0.02167879 0.83558202 -0.9078590
grass   0.15848477 1.00000000 0.06935884 -0.01735530 -0.2052506
soil    0.02167879 0.06935884 1.00000000 0.06579734 0.2116592
angle   0.83558202 -0.01735530 0.06579734 1.00000000 -0.8146941
distance -0.90785903 -0.20525060 0.21165916 -0.81469413 1.0000000
```

- (b) (5 points) Run the linear regression model for the above model and give the standard summaries (parameter estimates and standard errors, ANOVA table, etc). What evidence is there that some of the variables are useful in describing nesting frequency?

```
> puffin.lm <- lm(nesting ~ grass + soil + angle + distance, data=puffin)
```

```

> summary(puffin.lm)
Call:
lm(formula = nesting ~ grass + soil + angle + distance, data = puffin)

Residuals:
    Min       1Q   Median       3Q      Max
-4.0166 -2.1088  0.2293  1.2505  6.9881

Coefficients:
            Estimate Std. Error t value Pr(>|t|)
(Intercept) 10.117840   3.185028   3.177  0.00323 **
grass       -0.007408   0.019459  -0.381  0.70586
soil        0.209211   0.077238   2.709  0.01062 *
angle       0.082389   0.077796   1.059  0.29727
distance   -0.366571   0.057473  -6.378 3.18e-07 ***
---
Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1

Residual standard error: 2.647 on 33 degrees of freedom
Multiple R-Squared: 0.8792,    Adjusted R-squared: 0.8645
F-statistic: 60.03 on 4 and 33 DF,  p-value: 1.113e-14

```

```
> anova(puffin.lm)
```

```
Analysis of Variance Table
```

```
Response: nesting
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
grass	1	48.08	48.08	6.8599	0.01321	*
soil	1	0.22	0.22	0.0313	0.86057	
angle	1	1349.50	1349.50	192.5410	2.506e-15	***
distance	1	285.12	285.12	40.6802	3.184e-07	***
Residuals	33	231.29	7.01			

```
---
```

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

The clearest evidence is given by the F test from the summary output which examines the hypotheses

$$H_0 : \beta_1 = \beta_2 = \beta_3 = \beta_4 = 0$$

$$H_A : \text{at least one } \beta_i \neq 0, i = 1, \dots, 4$$

F-statistic: 60.03 on 4 and 33 DF, p-value: 1.113e-14

which is highly significant. At least one of the predictors appears to be useful in predicting nesting frequency. To determine the most likely ones are distance and soil based on the t tests from the summary output. After these two variables have been accounted for, the other two variables don't add much.

The F tests from the ANOVA table are not particularly useful in this case as they do not give tests that we are interested in this case.

Aside: Common Tests in Regression

For a regression model of the form

$$y_i = \beta_0 + \beta_1 x_{1i} + \beta_2 x_{2i} + \dots + \beta_p x_{pi} + \epsilon_i$$

there are three common testing situations about the predictors

1. Testing all β s simultaneously:

$$H_0 : \beta_1 = \beta_2 = \dots = \beta_p = 0$$

$$H_A : \text{at least one } \beta_i \neq 0, i = 1, \dots, p$$

This is equivalent to comparing the models

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 x_{1i} + \beta_2 x_{2i} + \dots + \beta_p x_{pi} + \epsilon_i$$

2. Testing one β :

$$H_0 : \beta_j = 0$$

$$H_A : \beta_j \neq 0$$

This is equivalent to comparing the models (when $j = 1$)

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \beta_2 x_{2i} + \dots + \beta_p x_{pi} + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 x_{1i} + \beta_2 x_{2i} + \dots + \beta_p x_{pi} + \epsilon_i$$

3. Testing a subset of the β s e.g.

$$H_0 : \beta_1 = \beta_2 = 0$$

$$H_A : \beta_1 \neq 0 \text{ or } \beta_2 \neq 0$$

The alternative hypothesis corresponds to the three situations

1. $\beta_1 = 0, \beta_2 \neq 0, \beta_3, \dots, \beta_p$ arbitrary
2. $\beta_1 \neq 0, \beta_2 = 0, \beta_3, \dots, \beta_p$ arbitrary
3. $\beta_1 \neq 0, \beta_2 \neq 0, \beta_3, \dots, \beta_p$ arbitrary

This is equivalent to comparing the models

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \beta_3 x_{1i} + \dots + \beta_p x_{pi} + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 x_{1i} + \beta_2 x_{2i} + \dots + \beta_p x_{pi} + \epsilon_i$$

These situations can all be examined within the same framework. Let $SSE(m)$ be the error sum of squares for model m ,

$$SSE(m) = \sum_{i=1}^n (y_i - \hat{y}_i(m))^2$$

where $\hat{y}_i(m)$ is the fitted value for observation i under model m . Now let $df(m)$ be the error degrees of freedom for model m ,

$$df(m) = n - k(m)$$

where $k(m)$ is the number of predictors in model m .

The usual test statistic in all three cases is

$$F = \frac{(SSE(R) - SSE(F)) / (df(R) - df(F))}{SSE(F) / df(F)}$$

which is compared to an F distribution with $df(R) - df(F)$ and $df(F)$ degrees of freedom. The numerator degrees of freedom $df(R) - df(F)$ is the number of parameters given in H_0 .

This is exactly what the **S** `anova` function does when given two `lm` objects.

It is possible to show that in the case of testing a single β , the t given in the summary output is equivalent to the F test as

$$t^2 = F$$

and if $t \sim t_{df}$, then $F = t^2 \sim F_{1,df}$.

The output given by `anova(puffin.lm)` earlier doesn't test what we really want to do

```
> anova(puffin.lm)
```

```
Analysis of Variance Table
```

```
Response: nesting
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
grass	1	48.08	48.08	6.8599	0.01321	*
soil	1	0.22	0.22	0.0313	0.86057	
angle	1	1349.50	1349.50	192.5410	2.506e-15	***
distance	1	285.12	285.12	40.6802	3.184e-07	***
Residuals	33	231.29	7.01			

What **S** does in this case examines the following sequence of models

1. grass

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
grass	1	48.08	48.08	6.8599	0.01321 *

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 \text{grass}_i + \epsilon_i$$

2. soil

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
soil	1	0.22	0.22	0.0313	0.86057

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \beta_1 \text{grass}_i + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \epsilon_i$$

3. angle

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
angle	1	1349.50	1349.50	192.5410	2.506e-15 ***

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \beta_3 \text{angle}_i + \epsilon_i$$

4. distance

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
distance	1	285.12	285.12	40.6802	3.184e-07 ***

$$\text{Reduced Model}(H_0) : y_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \beta_3 \text{angle}_i + \epsilon_i$$

$$\text{Full Model}(H_A) : y_i = \beta_0 + \beta_1 \text{grass}_i + \beta_2 \text{soil}_i + \beta_3 \text{angle}_i + \beta_4 \text{distance}_i + \epsilon_i$$

So the order that the variables are entered into the model affects the results given by anova **unless** $\text{Var}(X)$ is a diagonal matrix, i.e.

$$\text{Corr}(X_i, X_j) = 0 \quad \text{for all pairs } i \text{ \& } j$$

A partial justification of this can be seen by looking at the following regressions

```
> puff.lm <- lm(nesting ~ distance, data=puffin)
> angle.lm <- lm(angle ~ distance, data=puffin)

> nestres <- resid(puff.lm)
> angleres <- resid(angle.lm)

> puffang.lm <- lm(nestres ~ angleres)
```



```
> summary(puffang.lm)
```

Coefficients:

	Estimate	Std. Error	t value	Pr(> t)
(Intercept)	-2.045e-17	4.557e-01	-4.49e-17	1.0000
angleres	1.755e-01	6.811e-02	2.577	0.0142 *

```
> puffboth.lm <- lm(nesting ~ distance + angle, data=puffin)
```

```
> summary(puffboth.lm)
```

Coefficients:

	Estimate	Std. Error	t value	Pr(> t)
(Intercept)	13.95582	2.44522	5.707	1.88e-06 ***
distance	-0.29297	0.04871	-6.015	7.39e-07 ***
angle	0.17554	0.06908	2.541	0.0156 *

```
> puffa.lm <- lm(nesting ~ angle, data=puffin)
> summary(puffa.lm)
```

Coefficients:

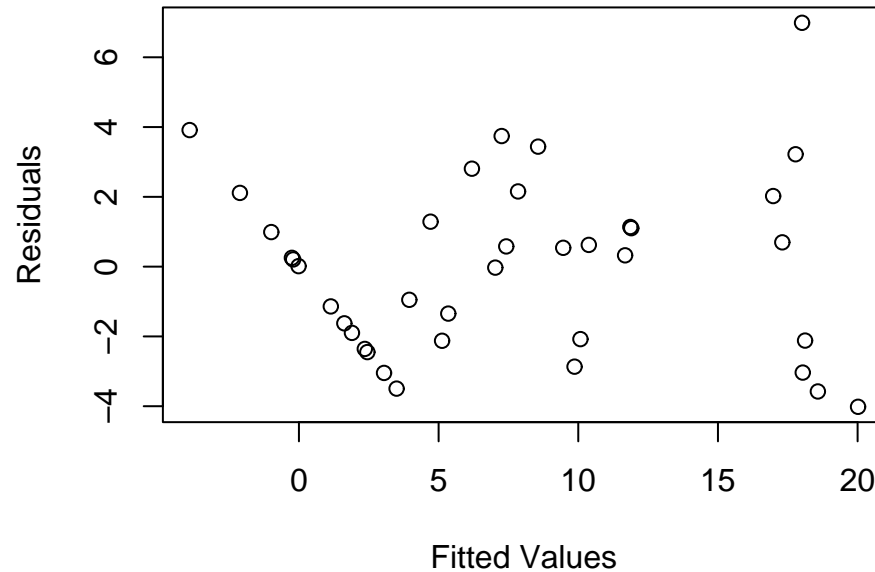
	Estimate	Std. Error	t value	Pr(> t)
(Intercept)	-0.02635	1.06591	-0.025	0.98
angle	0.51404	0.05633	9.126	6.75e-11 ***

So in multiple regression, the parameter estimate you get for a single predictor is the same as you get by regressing the residuals from a model with other predictors with the residuals you get from trying to predict the predictor of interest with the other predictors. (Sorry, I know this looks ugly).

So the bottom line is that what you get out of a multiple regression model for a variable, is what you get if you enter it last in a sequence of single variable regression models.

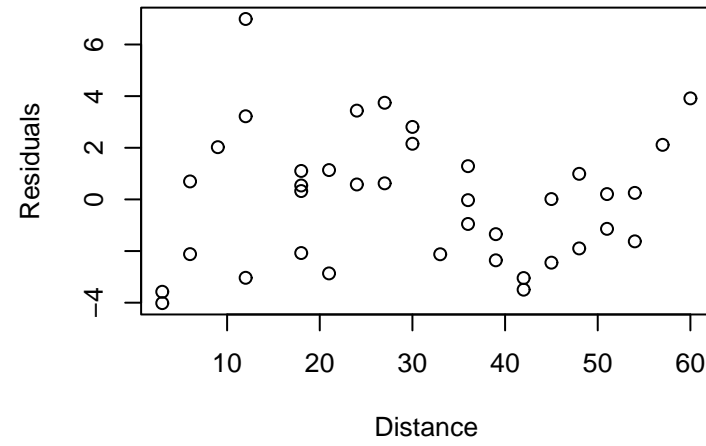
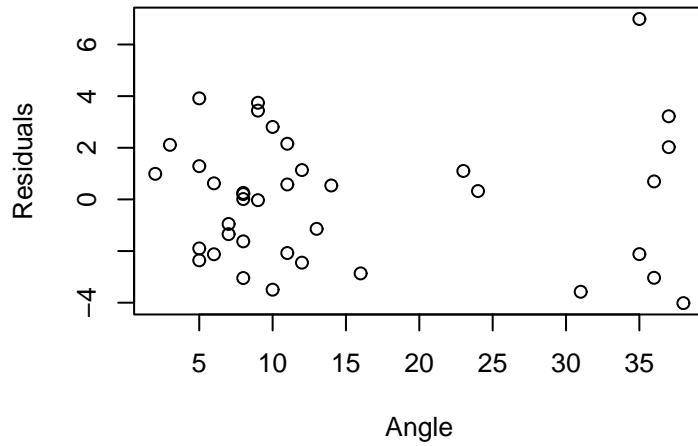
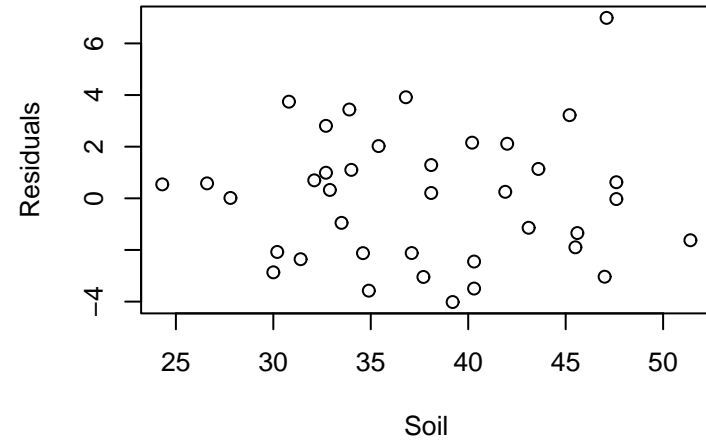
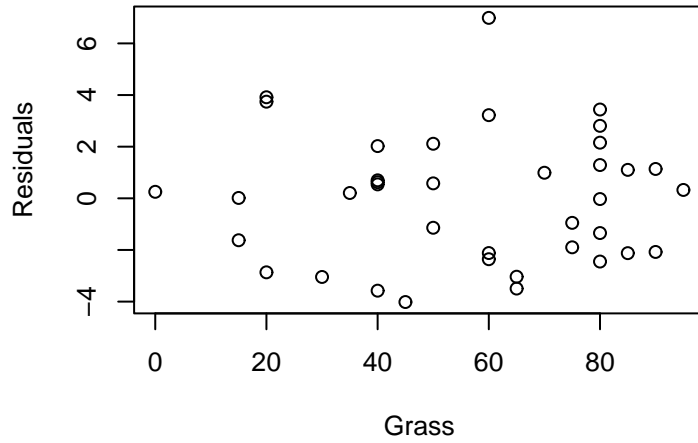
End Aside

- (c) (5 points) Create the residual plot of the residuals against the fitted values. In addition, in a single figure, plot the residuals against each of the predictor variables. Do any of these figures suggest a problem with the regression model?



```
plot(fitted(puffin.lm), resid(puffin.lm), xlab="Fitted Values",  
     ylab="Residuals")
```

Residuals vs Predictors



```
attach(puffin)
postscript("../Assignments/puffinres.eps", horiz=F, width=8, height=6.5)
par(mfrow=c(2,2), oma=c(0,0,4,0))
plot(grass, resid(puffin.lm), xlab="Grass", ylab="Residuals")
plot(soil, resid(puffin.lm), xlab="Soil", ylab="Residuals")
plot(angle, resid(puffin.lm), xlab="Angle", ylab="Residuals")
plot(distance, resid(puffin.lm), xlab="Distance", ylab="Residuals")
mtext(side=3, line=0, cex=1.5, outer=T, "Residuals vs Predictors")
```

The plots of the residuals versus the predictors don't show major problems. One outlier is suggested, though with 40 observations, not much to worry about, except for the fact that it is one the edge of the predictor space (high angle, low distance, high soil).

The plot of residuals versus the fits shows one problem, the decreasing line on the left of the plot. While it looks like curvature, that really isn't an accurate description of the problem. These are actually observations having the same value for nesting, in this case 0. It stands out in this case since it occurs with the minimum possible value of nesting.

(d) (5 points) Suppose that a new site was found where `grass = 95`, `soil = 25`, `angle = 5`, and `distance = 60`. Predict the number of nests for this site based on the original 38 sites. Any comments about this prediction?

```
> newdata <- data.frame(grass=95, soil=25,  
                        angle=5, distance=60)  
  
> predict(puffin.lm,newdata)  
[1] -6.938013
```

Since the `nesting` can't be negative, this suggests a problem with the model. In fact we probably should have fit a Poisson regression model as the model fit allows for negative nesting values and the `nesting` values are counts.

- (e) (5 points) Calculate the F test for comparing the model with all four predictors in the model with the model having only soil and distance in the model. What does this F test imply about the predictors nesting?

This test examines the hypotheses

$$H_0 : \beta_1 = \beta_3 = 0$$

$$H_A : \beta_1 \neq 0 \text{ or } \beta_3 \neq 0$$

```
> puffinred.lm <- lm(nesting ~ soil + distance, data=puffin)
```

```
> anova(puffinred.lm, puffin.lm)
```

```
Model 1: nesting ~ soil + distance
```

```
Model 2: nesting ~ grass + soil + angle + distance
```

	Res.Df	RSS	Df	Sum of Sq	F	Pr(>F)
1	35	244.869				
2	33	231.293	2	13.576	0.9685	0.3902

In this case the p -value is large, suggesting that grass and angle do not add anything to the predictions after soil and distance have been accounted for.

(f) (5 points) Now fit the model

$$\text{nesting}_i = \beta_0 + \beta_1 \text{soil}_i + \beta_2 \text{distance}_i + \beta_3 \text{soil}_i \times \text{distance}_i + \epsilon_i$$

Is there any evidence of an interaction between distance and soil on the nesting frequency?

```
> puffinint.lm <- lm(nesting ~ soil * distance, data=puffin)
```

```
> summary(puffinint.lm)
```

Coefficients:

	Estimate	Std. Error	t value	Pr(> t)	
(Intercept)	3.742393	5.679834	0.659	0.51440	
soil	0.435530	0.148838	2.926	0.00608	**
distance	-0.157837	0.178889	-0.882	0.38380	
soil:distance	-0.006575	0.004535	-1.450	0.15625	

```
> anova(puffinint.lm)
Analysis of Variance Table
```

```
Response: nesting
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
soil	1	0.90	0.90	0.1326	0.7180	
distance	1	1668.44	1668.44	245.9860	<2e-16	***
soil:distance	1	14.26	14.26	2.1022	0.1563	
Residuals	34	230.61	6.78			

The test on the interaction (either t or F is fine here) clearly isn't significant

(g) (5 points) Now fit the model

$$\text{nesting}_i = \beta_0 + \beta_1 \text{soil}_i + \beta_2 \text{distance}_i + \beta_3 \text{soil}_i^2 + \beta_4 \text{distance}_i^2 + \epsilon_i$$

Is there any evidence of a nonlinearity in the relationship of distance or soil on the nesting frequency?

```
> puffinquad.lm <- lm(nesting ~ soil + distance + I(soil^2)
+ I(distance^2), data=puffin)
```

```
> summary(puffinquad.lm)
```

Coefficients:

	Estimate	Std. Error	t value	Pr(> t)	
(Intercept)	16.220574	14.035610	1.156	0.256114	
soil	0.024878	0.732079	0.034	0.973095	
distance	-0.495445	0.113372	-4.370	0.000116	***
I(soil^2)	0.002714	0.009538	0.285	0.777772	
I(distance^2)	0.001348	0.001826	0.738	0.465510	

```

> anova(puffinquad.lm)
Response: nesting
      Df Sum Sq Mean Sq F value Pr(>F)
soil   1   0.90   0.90   0.1233 0.7277
distance 1 1668.44 1668.44 228.7155 <2e-16 ***
I(soil^2) 1   0.16   0.16   0.0223 0.8822
I(distance^2) 1   3.98   3.98   0.5452 0.4655
Residuals 33 240.73   7.29

```

Since both t tests on the quadratic terms are clearly not significant, the linearity of the relationship seems ok. However a better test of this (which agrees with the conclusion) is

```

> anova(puffinred.lm, puffinquad.lm)
Model 1: nesting ~ soil + distance Model 2: nesting ~ soil +
distance + I(soil^2) + I(distance^2)
  Res.Df  RSS Df Sum of Sq    F Pr(>F)
1     35 244.87
2     33 240.73  2     4.14 0.2837 0.7548

```

2. (30 points) The Ryan-Joiner test considers the following hypotheses

$H_0 : \{x_1, x_2, \dots, x_n\}$ is a random sample from a normal population

$H_A : \{x_1, x_2, \dots, x_n\}$ is not a random sample from a normal population

The test statistic is r , the correlation coefficient of the coordinate pairs,

$$\left(\Phi^{-1} \left(\frac{i - 0.375}{n + 0.25} \right), x_{(i)} \right), \quad i = 1, \dots, n$$

of a normal Q-Q plot, where $\Phi^{-1}(\cdot)$ denotes the inverse CDF of a standard normal RV and $x_{(1)} \leq x_{(2)} \leq \dots \leq x_{(n)}$ are the ordered data values. Under H_0 , r should be close to 1.

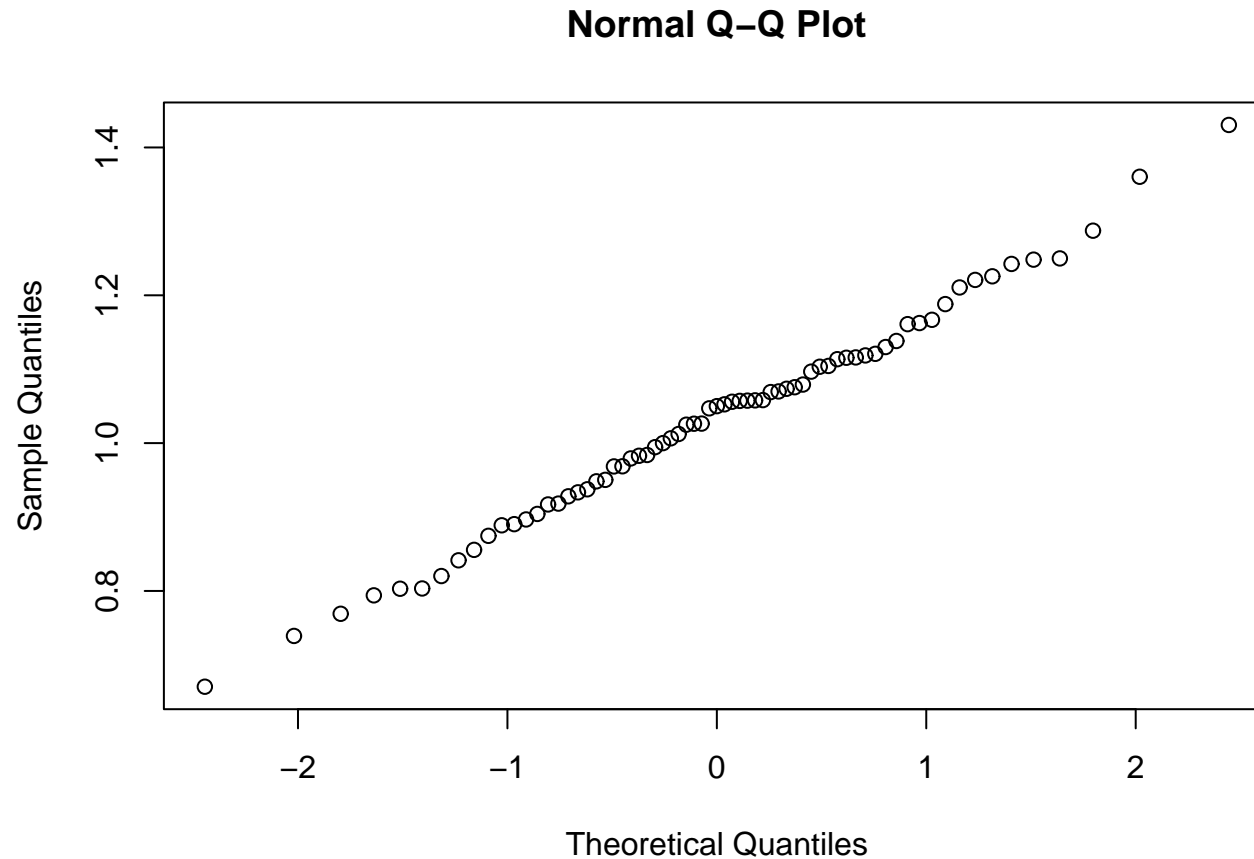
- (a) (5 points) For the dataset `douglasfir.txt`, available on the datasets page for the course, calculate in **R**, the value of the statistic r described above.

```
rjstat <- function(x) {  
  n <- length(x)  
  cor(qnorm(((1:n)-0.375)/(n+0.25)), sort(x))  
}  
>  
> rjstat(douglasfir)  
[1] 0.996571
```

- (b) (5 points) Create the Q-Q plot (aka Normal Scores plot) for this dataset.

The plot can be created by

```
qqnorm(douglasfir)
```



- (c) (5 points) The function `qqline` will add a straight line to Q-Q plots created by either `qqnorm` or `qqplot`. This line is a description of the main trend in the plot. By examining the code for `qqline`, what line is added to the `qqnorm` plot (i.e. when `datax = FALSE`).

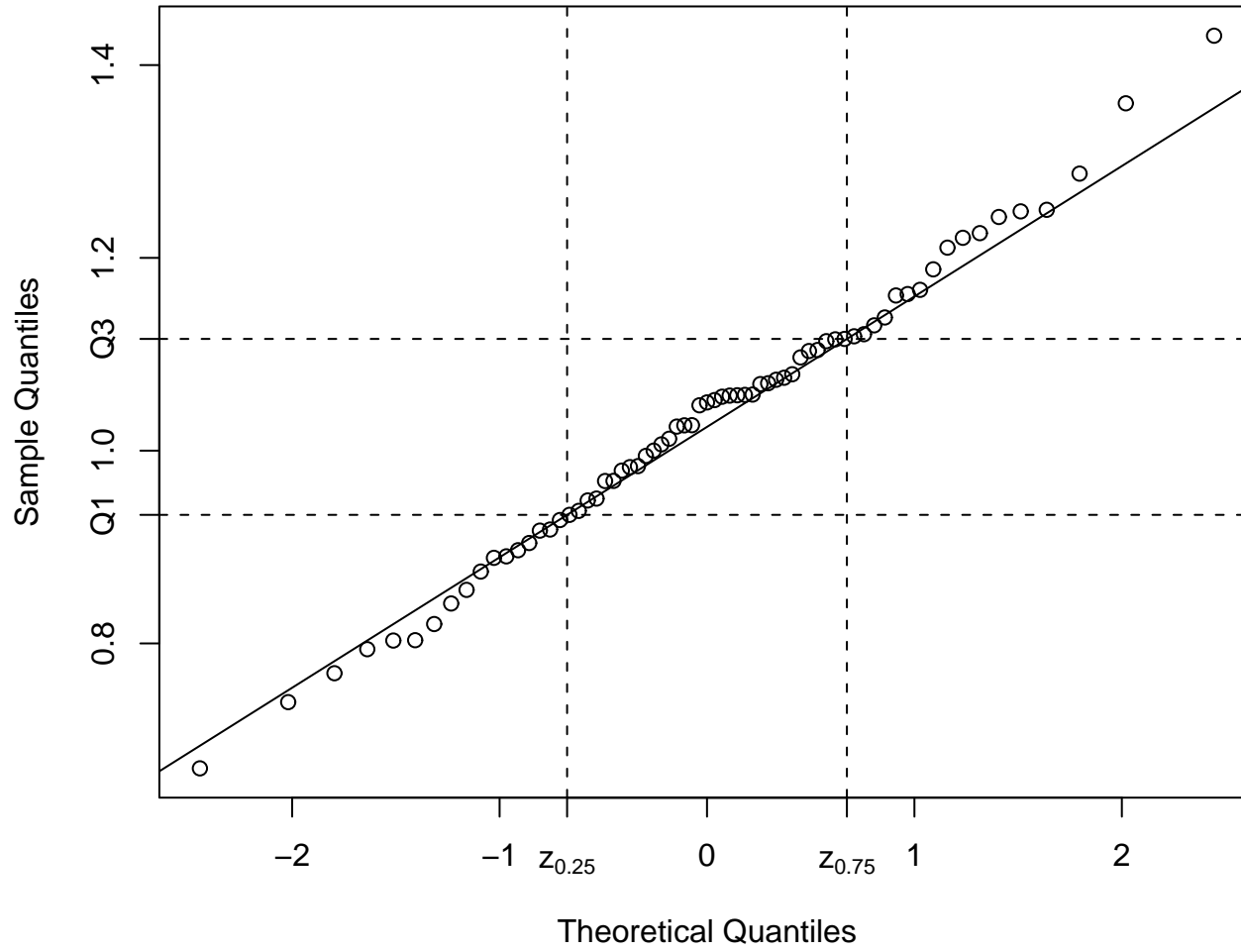
```

> qqline
function (y, datax = FALSE, ...) {
  y <- quantile(y[!is.na(y)], c(0.25, 0.75))
  x <- qnorm(c(0.25, 0.75))
  if (datax) {
    slope <- diff(x)/diff(y)
    int <- x[1] - slope * y[1]
  }
  else {
    slope <- diff(y)/diff(x)
    int <- y[1] - slope * x[1]
  }
  abline(int, slope, ...)
}

```

Let $Q1$ and $Q3$ be the sample quartiles of the dataset and $z_{0.25}$ and $z_{0.75}$ be the population quartiles of the $N(0, 1)$ distribution. Then `qqline` draws the line through the points $(z_{0.25}, Q1)$ and $(z_{0.75}, Q3)$. The slope of this line is an estimate of the standard deviation, though different than the sample standard deviation s .

Normal Q-Q Plot



(d) (10 points) Get p -value of test statistic by simulation

```
rjtest <- function(x, niter = 1000) {  
  n <- length(x)  
  r <- rjstat(douglasfir)  
  h0mat <- matrix(rnorm(n*niter), ncol=n)  
  rjh0 <- apply(h0mat, 1, rjstat)  
  pval <- mean(rjh0 <= r)  
  list(pval=pval, r=r, niter=niter)  
}
```

```
> rjtest(douglasfir)  
$pval [1] 0.934  
$r [1] 0.996571  
$niter [1] 1000
```

```
> rjtest(douglasfir,10000)  
$pval [1] 0.9514  
$r [1] 0.996571  
$niter [1] 10000
```

(e) (5 points) Based on the above analysis, is there any evidence to suggest that the Douglas fir data is not normally distributed?

No. The Q-Q plot is very linear and the p -value from the Ryan-Joiner test is large.

One comment about calculating the p -value. One way of thinking of a p -value is

$$p\text{-value} = P[\text{a test stat as or more extreme than observed} \mid H_0]$$

In this case, only small values of r should be considered extreme. r s close to 1 are highly consistent with H_0 , so a “one-sided” p -value should be calculated here, which is why

$$\hat{p} = \frac{1}{m} \sum_{j=1}^m I(r_j \leq r)$$

was used.