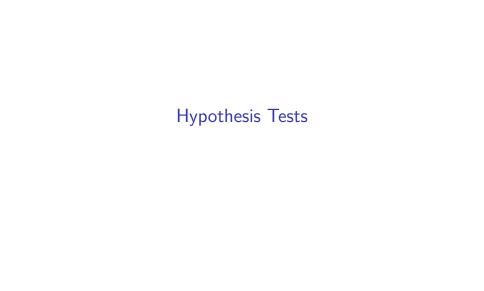
STAT 210 Applied Statistics and Data Analysis Experimental Design II

Joaquin Ortega



In an experiment involving k treatment levels, a hypothesis of interest is whether the treatments are different in terms of their effect on the response variable. Thus, the null hypothesis would be

$$H_0: \tau_1 = \tau_2 = \cdots = \tau_k$$
 vs. $H_1:$ at least two of the τ_i differ.

Even though it looks like the null hypothesis involves nonestimable parameters, it can easily be recasted in terms of estimable contrasts:

$$H_0: \tau_2 - \tau_1 = 0$$
, and $\tau_3 - \tau_1 = 0$ and ... and $\tau_k - \tau_1 = 0$.

There are other ways of rewriting H_0 in terms of estimable contrasts, but they will always depend on k-1 distinct contrasts, which is the number of degrees of freedom for treatment.

As we saw in the previous lecture, the basic idea of the analysis of variance is that the error sum of squares measures how well the model fits the data.

One way of testing whether the treatments have an effect is to compare the sums of squares for the complete model with that obtained with a model that assumes that the null hypothesis is true.

This last model is known as the reduced model and is

$$y_{ij} = \mu + \epsilon_{ij}$$

with the same hypothesis as before for the noise.

The least squares estimator for μ in this model is the overall mean:

$$\hat{\mu} = \bar{y}_{\bullet \bullet}$$

and the error sum of squares for the reduced model is

$$SST = \sum_{i} \sum_{j} (y_{ij} - \bar{y}_{\bullet \bullet})^2 = \sum_{i} \sum_{j} (y_{ij}^2 - 2y_{ij}\bar{y}_{\bullet \bullet} + \bar{y}_{\bullet \bullet}^2)$$
$$= \sum_{i} \sum_{j} y_{ij}^2 - n\bar{y}_{\bullet \bullet}^2. \tag{1}$$

If the null hypothesis is false and treatment effects differ, *SSE* should be smaller than *SST*.

The test is based on the difference SST - SSE, relative to the size of the SSE, i.e. (SST - SSE)/SSE and we reject H_0 if this quantity is large.

We used before the notation SSA for SST - SSE and called it the treatment sum of squares. Using (??) and (1), it is given by

$$SSA = \sum_{i} \sum_{j} y_{ij}^2 - n\bar{y}_{\bullet\bullet}^2 - \left(\sum_{i} \sum_{j} y_{ij}^2 - \sum_{i} r_i \bar{y}_{i\bullet}^2\right) = \sum_{i} r_i \bar{y}_{i\bullet}^2 - n\bar{y}_{\bullet\bullet}^2.$$

Equivalently,

$$SSA = \sum_{i} r_i (\bar{y}_{i\bullet} - \bar{y}_{\bullet\bullet})^2. \tag{2}$$

As we mentioned before, this sum represents the difference between treatment means and corresponds to the comparison between groups.

It can be shown that

- ► SSE/σ^2 has a χ^2_{n-k} distribution
- ► SSA/σ^2 has a χ^2_{k-1} distribution under H_0 ,
- these variables are independent.

In consequence, under H_0 the quotient

$$\frac{SSA/\sigma^2(k-1)}{SSE/\sigma^2(n-k)} \sim F_{k-1,n-k} \tag{3}$$

and we can use this relation to test H_0 .

Recall that we defined MSE = SSE/(n-k) and define MSA = SSA/(k-1), then (3) becomes

$$\frac{MSA}{MSE} \sim F_{k-1,n-k}$$

and if msE and msA represent the observed values of these variables, the decision rule for testing H_0 at level of significance α is

reject
$$H_0$$
 if $\frac{msA}{msE} > F_{1-\alpha,k-1,n-k}$ (4)

As we have seen, the values for the sums of squares, degrees of freedom, mean squares and F test are usually written in an Analysis of Variance table, such as Table 3 below.

Table 3: Anova table for the one-way analysis of variance	Table 3:	Anova	table	for the	one-way	analys	is of	variance
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Source	SS	d.f.	MS	F _{obs}	Critical <i>F</i>				
Treatment	SSA	k-1	$MSA = \frac{SSA}{k-1}$	$F = \frac{MSA}{MSE}$	qf(1- α , k-1, n-k)				
Error	SSE	n-k	$MSE = \frac{SSE}{n-k}$						
Total	SST	n-1							
Computational Formulae									
$SSA = \sum_{i} r_{i} \bar{y}_{i \bullet}^{2} - n \bar{y}_{\bullet \bullet}^{2}$				$SSE = \sum_{i} \sum_{j} y_{ij}^{2} - \sum_{i} r_{i} \bar{y}_{i\bullet}^{2}$					
$SST = \sum_{i} \sum_{j} y_{ij}^2 - n \bar{y}_{ullet}^2$									

Loading required package: e1071

library(PASWR)

Signif. codes:

An Anova table can be obtained in R using the aov function. The inputs for this function are the same as those for the 1m function we used previously, but the summary of an object created with the aov function is the Anova table. For the tire tread example the code is

'***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' '

The results for treatment are in the line labeled tire while the results corresponding to the errors are in the line labeled Residuals.

Results for *total* can be obtained adding up the corresponding terms in the table.

The F value is the ratio MSA/MSE and the last column labeled Pr(>F) is the probability of exceeding the calculated F-value when the null hypothesis is true, i.e. it is the p-value for the F test.

In this example we would conclude that there are significant differences among the average braking distances for different treads at the 0.05 level and also at the 0.01 level.

Residuals 20 7098.8 354.94

anova (mod0)

The same table is produced using the function anova() on the model we obtained with the lm function (mod0).

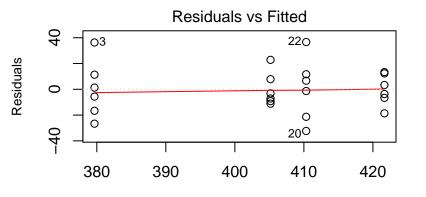
```
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.3
```

To see the complete list of quantities calculated by the aov function type names(mod1):

```
##
    [1] "coefficients"
                        "residuals"
                                        "effects"
                        "fitted.values" "assign"
##
   [4] "rank"
                                        "contrasts"
   [7] "qr"
                        "df.residual"
##
   [10] "xlevels"
                        "call"
                                        "terms"
## [13] "model"
                 "Sum Sq" "Mean Sq" "F value" "Pr(>F)"
## [1] "Df"
```

Look at diagnostic plots to check the assumptions of the model.

```
plot(mod1, which=1, cex.lab=0.8, cex.sub=0.8)
```



Fitted values aov(StopDist ~ tire)

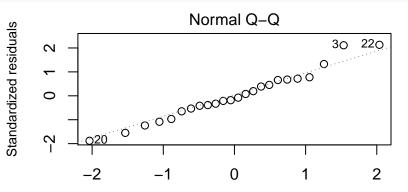
Since we have six replications for each treatment level and there are only four x-values, the points appear vertically aligned at these values.

We look in this graphs for constant variance. We see that values in some cases appear to be more spread than in others, and this may be a sign of non-constant variance.

However, we only have a few points and this is difficult to determine.

We look for departures from normality in the standardized residuals. Considering the fit we observe, the normality assumption seems justified.

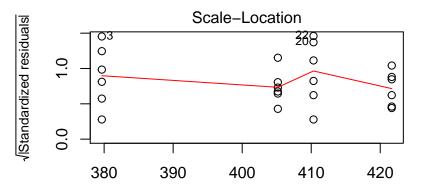
plot(mod1, which=2, cex.lab=0.8, cex.sub=0.8)



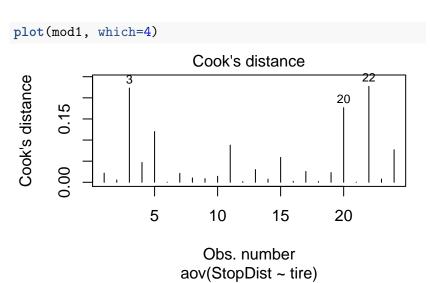
Theoretical Quantiles aov(StopDist ~ tire)

The third graph plots the square root of standardized residuals versus fitted values, and again we look for changes in the variance. As in the first graph, we see that some points seem to have more spread than others, perhaps pointing to heterocedasticity. However, there are very few points to be certain.

plot(mod1, which=3, cex.lab=0.8, cex.sub=0.8)



Cittod voluce



Another plot that is frequently useful is the plot of residuals versus the order in which the experiments have been performed.

This may help detecting if there is tendency in the results that should be taken into account.

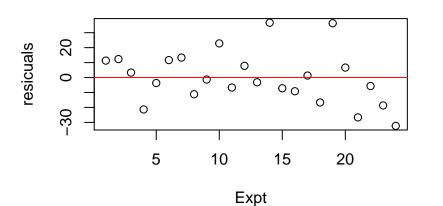
In this case that information is not available in the data set.

To illustrate the procedure, we will do the graph using a random ordering of the experiments.

To get the residuals for the model use the command residuals (mode1).

```
plot(residuals(mod1) ~ sample(1:24,24), xlab='Expt',
        ylab='resicuals', main='Residuals vs. Exp. Unit',
        font.main=1)
abline(h=0, col='red')
```

Residuals vs. Exp. Unit



Effect Sizes

replic.

To see the effect sizes in tabular form use model.tables

```
model.tables(mod1, se=T)
## Tables of effects
##
## tire
## tire
##
            В
## -24.542 0.958 17.458 6.125
##
## Standard errors of effects
##
           tire
         7.691
##
```

Effect Sizes

Specifying means you get

```
model.tables(mod1, 'means', se=T)
## Tables of means
## Grand mean
##
## 404.2083
##
## tire
## tire
##
      A B C
## 379.7 405.2 421.7 410.3
##
## Standard errors for differences of means
##
           tire
##
          10.88
## replic.
```

If the result of the F test is to reject the null hypothesis of no treatment effects one is naturally interested in determining where the difference lies. For this, it becomes necessary to compare the individual groups.

This can be (partly) done looking at the regression coefficients using summary and lm.

```
summary(mod0)
##
## Call:
## lm(formula = StopDist ~ tire, data = Tire)
##
## Residuals:
      Min 1Q Median 3Q
                                  Max
##
## -32.333 -9.667 -2.250 11.417 36.667
##
## Coefficients:
             Estimate Std. Error t value Pr(>|t|)
##
## (Intercept) 379.667 7.691 49.363 < 2e-16 ***
## tireB 25.500 10.877 2.344 0.029497 *
## tireC 42.000 10.877 3.861 0.000973 ***
         30.667 10.877 2.819 0.010594 *
## tireD
## ---
## Signif. codes:
## 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 18.84 on 20 degrees of freedom
## Multiple R-squared: 0.4442, Adjusted R-squared: 0.3608
## F-statistic: 5.328 on 3 and 20 DF, p-value: 0.007316
```

Recall that default coding for treatment in R means that the first value (Intercept) corresponds to the average value for the first treatment level $\hat{\mu} + \hat{\tau}_1$ while tireB = $\hat{\tau}_2 - \hat{\tau}_1$, tireC = $\hat{\tau}_3 - \hat{\tau}_1$, and tireD = $\hat{\tau}_4 - \hat{\tau}_1$

The t value in the table for timeB gives a t-test for comparing the first two groups (tread A vs. B, p-value = 0.0294), the next row in the table has a t-test for comparing groups 1 and 3 (tread A vs C, p-value=0.0009) and finally tireD compares tread A vs D, with p-value 0.011.

However, there is not a test for comparing groups 2 and 3, or 2 and 4 in the table.

One could redifine the factor time to include some of the missing comparison but then some others will dissapear.

This is not a convenient way to get all the tests, particularly if there are many levels for the treatment.

There exists a function for comparing all groups called pairwise.t.test.

However, one must correct for multiple testing.

The problem is that if we perform many tests, the probability of finding one of them to be significant by chance alone increases.

Consider one hundred statistical tests at the 5% level and assume all null hypothesis are true. We expect to reject 5 of them by chance alone. This is the expected value of Type I errors.

If the tests are independent, the probability of rejecting at least one hull hypothesis can be calculated using the binomial distribution:

```
1 - dbinom(0,100,0.05); 1 - dbinom(0,100,0.01)
```

```
## [1] 0.9940795
```

```
## [1] 0.6339677
```

One simple and frequently used method for correction is based on the Bonferroni inequalities and is known as the Bonferroni correction:

$$P(\cup_1^n B_i) \leq \sum_1^n P(B_i)$$

Thus, by dividing the significance level by the number of tests we get a test with true significance level smaller or equal to the nominal significance level.

This is equivalent to multiplying the p-values by the number of tests.

This test is conservative (tends to produce p values that are larger than they really are).

The Bonferroni correction should be applied at the planning stage of the experiment.

```
with(Tire, pairwise.t.test(StopDist, tire,
               p.adjust.method = 'bonferroni'))
##
   Pairwise comparisons using t tests with pooled SD
##
##
## data: StopDist and tire
##
## A B C
## B 0.1770 -
## C 0.0058 0.8696 -
## D 0.0636 1.0000 1.0000
##
## P value adjustment method: bonferroni
```

We have seen that the Bonferroni inequality can be used in the case of preplanned comparisons.

We will consider here briefly another method proposed by Tukey for comparisons of the form $H_0: \tau_s = \tau_u$ for $s \neq u$ in favor of the alternative $H_1: \tau_s \neq \tau_u$.

The procedure considers simultaneously all pairs of effects and adjusts the critical region by using the studentized range statistic instead of student's *t*-distribution.

The test is

reject
$$H_0$$
 if $|\hat{\tau}_u - \hat{\tau}_s| > \sqrt{2}q_{I,n-k,1-\alpha/2}\hat{\sigma}(\bar{y}_{sullet} - \bar{y}_{uullet})$

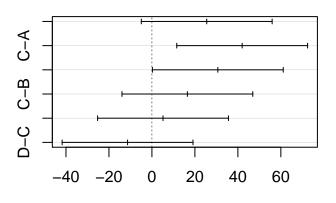
where $q_{I,n-k,1-\alpha}$ is the $1-\alpha$ percentile of the studentized range and $\hat{\sigma}(\bar{y}_{s\bullet}-\bar{y}_{u\bullet})$ is the estimated standard error for the difference between the averages.

If $X_1, ..., X_I$ are independent random variables with a $N(\mu, \sigma^2)$ distribution and $R = \max_i X_i - \min_i X_i$ then $R/\hat{\sigma}$ follows the studentized range distribution.

This is only approximate when the sample sizes are unequal.

```
(mod1.tky <-TukeyHSD(mod1))</pre>
##
    Tukey multiple comparisons of means
##
      95% family-wise confidence level
##
## Fit: aov(formula = StopDist ~ tire, data = Tire)
##
## $tire
##
            diff
                         lwr
                                upr padj
## B-A 25.500000 -4.9446409 55.94464 0.1213153
## C-A 42.000000 11.5553591 72.44464 0.0049515
## D-A 30.666667 0.2220258 61.11131 0.0479540
## C-B 16.500000 -13.9446409 46.94464 0.4464584
## D-B 5.166667 -25.2779742 35.61131 0.9637307
## D-C -11.333333 -41.7779742 19.11131 0.7273681
```

95% family-wise confidence level



Differences in mean levels of tire