

ANOVA Table with Degrees of Freedom

each line is 1 df and we have a total of r 1-df lines

| Sum of Squares | Degrees of Freedom | DF |
|--|---|---------------|
| $\mathbf{y}^\top (\mathbf{P}_2 - \mathbf{P}_1) \mathbf{y}$ | $\text{rank}(\mathbf{X}_2) - \text{rank}(\mathbf{X}_1)$ | $r_2 - 1$ |
| $\mathbf{y}^\top (\mathbf{P}_3 - \mathbf{P}_2) \mathbf{y}$ | $\text{rank}(\mathbf{X}_3) - \text{rank}(\mathbf{X}_2)$ | $r_3 - r_2$ |
| \vdots | \vdots | \vdots |
| $\mathbf{y}^\top (\mathbf{P}_m - \mathbf{P}_{m-1}) \mathbf{y}$ | $\text{rank}(\mathbf{X}_m) - \text{rank}(\mathbf{X}_{m-1})$ | $r - r_{m-1}$ |
| $\mathbf{y}^\top (\mathbf{P}_{m+1} - \mathbf{P}_m) \mathbf{y}$ | $\text{rank}(\mathbf{X}_{m+1}) - \text{rank}(\mathbf{X}_m)$ | $n - r$ |
| $\mathbf{y}^\top (\mathbf{I} - \mathbf{P}_1) \mathbf{y}$ | $\text{rank}(\mathbf{X}_{m+1}) - \text{rank}(\mathbf{X}_1)$ | $n - 1$ |

Mean Squares

For $j = 1, \dots, m - 1$, define

$$MS(j + 1 \mid j) = \frac{SS(j + 1 \mid j)}{r_{j+1} - r_j} = \frac{\mathbf{y}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{y}}{r_{j+1} - r_j}.$$

These sums of squares divided by their degrees of freedom are known as *mean squares*.

ANOVA Table with Mean Squares

$$MS(j+1|j) / MSE = F$$

| Sum of Squares | Degrees of Freedom | Mean Square |
|-----------------|--------------------|------------------------|
| $SS(2 1)$ | $r_2 - 1$ | $MS(2 1)$ |
| $SS(3 2)$ | $r_3 - r_2$ | $MS(3 2)$ |
| \vdots | \vdots | \vdots |
| $SS(m m - 1)$ | $r - r_{m-1}$ | $MS(m m - 1)$ |
| SSE | $n - r$ | $MSE - \hat{\sigma}^2$ |
| $SSTo$ | $n - 1$ | |

Independence of ANOVA Sums of Squares

Because

$$(P_{j+1} - P_j) (\sigma^2 I) (P_{\ell+1} - P_\ell) = 0$$

for all $j \neq \ell$, any two ANOVA sums of squares (not including SST_o) are independent.

It is also true that the ANOVA sums of squares (not including SST_o) are mutually independent by Cochran's Theorem, but that stronger result is not usually needed.

Stat 6110

ANOVA F Statistics

For $j = 1, \dots, m - 1$ we have

$$\underline{F_j} = \frac{MS(j+1 | j)}{MSE} = \frac{\mathbf{y}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{y} / (r_{j+1} - r_j)}{\mathbf{y}^\top (\mathbf{I} - \mathbf{P}_X) \mathbf{y} / (n - r)}$$

$$\sim F_{\boxed{r_{j+1} - r_j}, n-r} \left(\underbrace{\frac{\beta^\top \mathbf{X}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{X} \beta}{2\sigma^2}}_{ncp} \right).$$

Handwritten notes:
- "den df" with an arrow pointing to the denominator of the fraction inside the F-distribution.
- "n df" with an arrow pointing to the first parameter of the F-distribution, $r_{j+1} - r_j$.
- "ncp" with a bracket underneath the fraction inside the F-distribution.

ANOVA Table with F -Statistics

| Sum of Squares | Degrees of Freedom | Mean Square | F Stat |
|--------------------|--------------------|---------------|-----------|
| $SS(2 \mid 1)$ | $r_2 - 1$ | $MS(2 1)$ | F_1 |
| $SS(3 \mid 2)$ | $r_3 - r_2$ | $MS(3 2)$ | F_2 |
| \vdots | \vdots | \vdots | \vdots |
| $SS(m \mid m - 1)$ | $r - r_{m-1}$ | $MS(m m - 1)$ | F_{m-1} |
| SSE | $n - r$ | MSE | |
| SST_o | $n - 1$ | | |

Relationship with Reduced vs. Full Model F -Statistic

The ANOVA F_j statistic:

$$F_j = \frac{\mathbf{y}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{y} / (r_{j+1} - r_j)}{\mathbf{y}^\top (\mathbf{I} - \mathbf{P}_X) \mathbf{y} / (n - r)} = \frac{MS(j+1 | j)}{MSE}$$

The reduced vs. full model F statistic:

$$F = \frac{\mathbf{y}^\top (\mathbf{P}_X - \mathbf{P}_{X_0}) \mathbf{y} / (r - r_0)}{\mathbf{y}^\top (\mathbf{I} - \mathbf{P}_X) \mathbf{y} / (n - r)}$$

What do ANOVA F -statistics test?

In general, an F -statistic is used to test

$$\delta = \lambda_{PC} = 0$$

H_0 : “The non-centrality parameter of the F -statistic is zero.”

vs.

H_A : “The non-centrality parameter of the F -statistic is not zero.”

$$\delta = \lambda_{PC} \neq 0$$

What do ANOVA F -statistics test?

The ANOVA F -statistic

$$F_j = \frac{\mathbf{y}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{y} / (r_{j+1} - r_j)}{\mathbf{y}^\top (\mathbf{I} - \mathbf{P}_X) \mathbf{y} / (n - r)} = \frac{MS(j+1 | j)}{MSE}$$

has non-centrality parameter

$$mcp = \frac{\beta^\top \mathbf{X}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{X} \beta}{2\sigma^2}.$$

mcp will be zero
as long as
numerator = 0

Thus, F_j can be used to test

$$H_{0j} : \frac{\beta^\top \mathbf{X}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{X} \beta}{2\sigma^2} = 0 \text{ versus}$$

$$H_{Aj} : \frac{\beta^\top \mathbf{X}^\top (\mathbf{P}_{j+1} - \mathbf{P}_j) \mathbf{X} \beta}{2\sigma^2} \neq 0.$$

What do ANOVA F -statistics test? $= a$

$\beta^T X^T (P_{j+1} - P_j) X \beta$ due to symmetry & idempotency

The following are equivalent ways to write the null and alternative hypotheses tested by F_j .

$$H_{0j} : \mu_{cp} = 0$$

$$H_{Aj} : \mu_{cp} \neq 0 \quad \text{which is of form } a^T a$$

follows from previous slide (not considering 25^2 in denom)

$$\beta^T X^T (P_{j+1} - P_j) X \beta = 0$$

$$\beta^T X^T (P_{j+1} - P_j) X \beta \neq 0$$

$$(P_{j+1} - P_j) X \beta = 0$$

$$(P_{j+1} - P_j) X \beta \neq 0$$

which can only be 0 if $a = 0$

$$P_j E(y) = P_{j+1} E(y)$$

$$P_j E(y) \neq P_{j+1} E(y)$$

$$P_{j+1} E(y) \in \mathcal{C}(X_j)$$

$$P_{j+1} E(y) \in \mathcal{C}(X_{j+1}) \setminus \mathcal{C}(X_j)$$

for $E(y)$

$E(y)$ is found in the smaller of the two column spaces, thus it does matter in which column space we look

What do ANOVA F -statistics test?

Recall $C\beta$ is estimable $\Leftrightarrow C = AX$

$$H_{0j} : (P_{j+1} - P_j)X\beta = 0 \quad \text{vs.} \quad H_{Aj} : (P_{j+1} - P_j)\underline{X}\beta \neq 0$$

is of the form

$$H_{0j} : C_j^* \beta = 0 \quad \text{vs.} \quad H_{Aj} : C_j^* \beta \neq 0,$$

$$\text{where } C_j^* = \underbrace{(P_{j+1} - P_j)}_{C=A} \underline{X}.$$

As written, H_{0j} is not a testable hypothesis because C_j^* has n rows but $\text{rank } r_{j+1} - r_j < n$ (homework problem).

We can rewrite H_{0j} as a testable hypothesis by replacing C_j^* with any matrix C_j whose $q = r_{j+1} - r_j$ rows form a basis for the row space of C_j^* .

Example: Multiple Regression

multiple linear regression

intercept only model

adding a covariate assuming linear relationship with $y \Rightarrow$ SLR

$$\begin{aligned} X_1 &= 1 \\ X_2 &= [1, x_1] \\ X_3 &= [1, x_1, x_2] \\ &\vdots \\ X_m &= [1, x_1, \dots, x_{m-1}] \end{aligned}$$

$SS(j+1 | j)$ is the decrease in SSE that results when the explanatory variable x_j is added to a model containing an intercept and explanatory variables x_1, \dots, x_{j-1} .

Example: Polynomial Regression

$$\mathbf{X}_1 = 1$$

$$\mathbf{X}_2 = [1, \mathbf{x}]$$

$$\mathbf{X}_3 = [1, \mathbf{x}, \mathbf{x}^2]$$

$$\vdots$$

$$\mathbf{X}_m = [1, \mathbf{x}, \mathbf{x}^2, \dots, \mathbf{x}^{m-1}]$$

quadratic
model

$SS(j+1 \mid j)$ is the decrease in SSE that results when the explanatory variable x^j is added to a model containing an intercept and explanatory variables x, x^2, \dots, x^{j-1} .

An Example in R

```
> #An example from "Design of Experiments: Statistical  
> #Principles of Research Design and Analysis"  
> #2nd Edition by Robert O. Kuehl  
>  
> d=read.delim("https://.../S510/PlantDensity.txt")
```

The Data

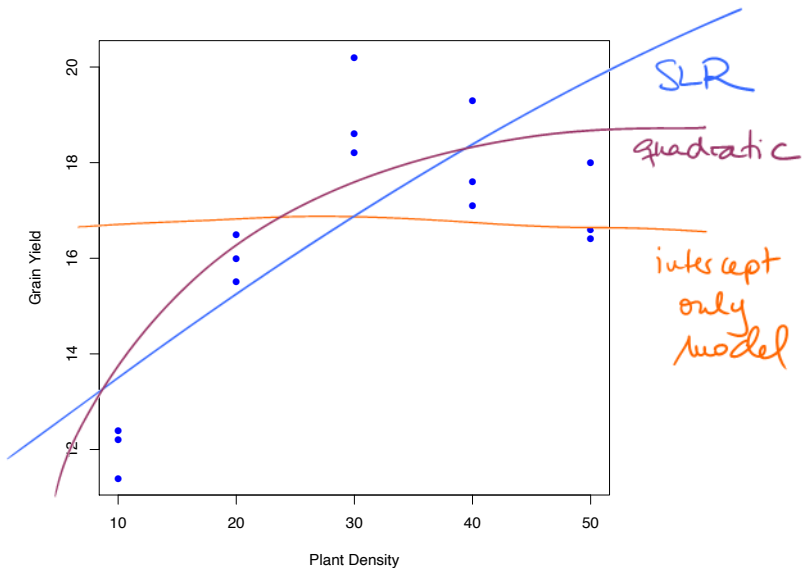
X

```
> d
  PlantDensity GrainYield
1           10      12.2
2           10      11.4
3           10      12.4
4           20      16.0
5           20      15.5
6           20      16.5
7           30      18.6
8           30      20.2
9           30      18.2
10          40      17.6
11          40      19.3
12          40      17.1
13          50      18.0
14          50      16.4
15          50      16.6
```

5 different levels
of treatment
3 replicates per
level

Renaming the Variables and Plotting the Data

```
> names(d)=c("x", "y")
> head(d)
      x      y
1 10 12.2
2 10 11.4
3 10 12.4
4 20 16.0
5 20 15.5
6 20 16.5
>
> plot(d[,1],d[,2],col=4,pch=16,xlab="Plant Density",
+      ylab="Grain Yield")
```

Matrices with Nested Column Spaces

$$\mathbf{X}_1 = \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{bmatrix}, \mathbf{X}_2 = \begin{bmatrix} 1 & 10 \\ 1 & 10 \\ 1 & 10 \\ 1 & 20 \\ 1 & 20 \\ 1 & 20 \\ 1 & 30 \\ 1 & 30 \\ 1 & 30 \\ 1 & 40 \\ 1 & 40 \\ 1 & 40 \\ 1 & 50 \\ 1 & 50 \\ 1 & 50 \end{bmatrix}, \mathbf{X}_3 = \begin{bmatrix} 1 & 10 & 100 \\ 1 & 10 & 100 \\ 1 & 10 & 100 \\ 1 & 20 & 400 \\ 1 & 20 & 400 \\ 1 & 20 & 400 \\ 1 & 30 & 900 \\ 1 & 30 & 900 \\ 1 & 30 & 900 \\ 1 & 40 & 1600 \\ 1 & 40 & 1600 \\ 1 & 40 & 1600 \\ 1 & 50 & 2500 \\ 1 & 50 & 2500 \\ 1 & 50 & 2500 \end{bmatrix},$$

Matrices with Nested Column Spaces

$$\mathbf{X}_4 = \begin{bmatrix} 1 & 10 & 100 & 1000 \\ 1 & 10 & 100 & 1000 \\ 1 & 10 & 100 & 1000 \\ 1 & 20 & 400 & 8000 \\ 1 & 20 & 400 & 8000 \\ 1 & 20 & 400 & 8000 \\ 1 & 30 & 900 & 27000 \\ 1 & 30 & 900 & 27000 \\ 1 & 30 & 900 & 27000 \\ 1 & 40 & 1600 & 64000 \\ 1 & 40 & 1600 & 64000 \\ 1 & 40 & 1600 & 64000 \\ 1 & 50 & 2500 & 125000 \\ 1 & 50 & 2500 & 125000 \\ 1 & 50 & 2500 & 125000 \end{bmatrix}, \mathbf{X}_5 = \begin{bmatrix} 1 & 10 & 100 & 1000 & 10000 \\ 1 & 10 & 100 & 1000 & 10000 \\ 1 & 10 & 100 & 1000 & 10000 \\ 1 & 20 & 400 & 8000 & 160000 \\ 1 & 20 & 400 & 8000 & 160000 \\ 1 & 20 & 400 & 8000 & 160000 \\ 1 & 30 & 900 & 27000 & 810000 \\ 1 & 30 & 900 & 27000 & 810000 \\ 1 & 30 & 900 & 27000 & 810000 \\ 1 & 40 & 1600 & 64000 & 2560000 \\ 1 & 40 & 1600 & 64000 & 2560000 \\ 1 & 40 & 1600 & 64000 & 2560000 \\ 1 & 50 & 2500 & 125000 & 6250000 \\ 1 & 50 & 2500 & 125000 & 6250000 \\ 1 & 50 & 2500 & 125000 & 6250000 \end{bmatrix}$$

Centering and Standardizing for Numerical Stability

It is typically best for numerical stability to center and scale a quantitative explanatory variable prior to computing higher order terms.

In the plant density example, we could replace x by $(x - 30)/10$ and work with the matrices on the next two slides.

Because these matrices have the same column spaces as the original matrices, the ANOVA table entries are mathematically identical for either set of matrices.

Matrices with Centered and Scaled x

[illegible]

Matrices with Centered and Scaled x

$$X_4 = \begin{bmatrix} 1 & -2 & 4 & -8 \\ 1 & -2 & 4 & -8 \\ 1 & -2 & 4 & -8 \\ 1 & -1 & 1 & -1 \\ 1 & -1 & 1 & -1 \\ 1 & -1 & 1 & -1 \\ 1 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 \\ 1 & 2 & 4 & 8 \\ 1 & 2 & 4 & 8 \\ 1 & 2 & 4 & 8 \end{bmatrix}$$

$, X_5 =$

$$\begin{bmatrix} 1 & -2 & 4 & -8 & 16 \\ 1 & -2 & 4 & -8 & 16 \\ 1 & -2 & 4 & -8 & 16 \\ 1 & -1 & 1 & -1 & 1 \\ 1 & -1 & 1 & -1 & 1 \\ 1 & -1 & 1 & -1 & 1 \\ 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 \\ 1 & 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 & 1 \\ 1 & 1 & 1 & 1 & 1 \\ 1 & 2 & 4 & 8 & 16 \\ 1 & 2 & 4 & 8 & 16 \\ 1 & 2 & 4 & 8 & 16 \end{bmatrix}$$

a different mean for each treatment

↓
5 different means

Regardless of whether we center and scale x , the column space of X_5 is the same as the column space of the cell means model matrix

$$\underline{\underline{X}} = \begin{bmatrix} 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 \\ 1 & 0 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix}.$$

ANOVA Table for the Plant Density Data

by how much will the $SE \downarrow$ if we add x to the model that contains only intercept


| Source | Sum of Squares | DF |
|------------------------|---------------------|---------------|
| $x 1$ intercept only | $y^T (P_2 - P_1) y$ | $2 - 1 = 1$ |
| $x^2 1, x$ | $y^T (P_3 - P_2) y$ | $3 - 2 = 1$ |
| $x^3 1, x, x^2$ | $y^T (P_4 - P_3) y$ | $4 - 3 = 1$ |
| $x^4 1, x, x^2, x^3$ | $y^T (P_5 - P_4) y$ | $5 - 4 = 1$ |
| Error | $y^T (I - P_5) y$ | $15 - 5 = 10$ |
| C. Total | $y^T (I - P_1) y$ | $15 - 1 = 14$ |

Creating the Matrices in R

```
> y=d$y
> x=(d$x-mean(d$x))/10
> x
[1] -2 -2 -2 -1 -1 -1  0  0  0  1  1  1  2  2  2
>
> n=nrow(d)
>
> x1=matrix(1,nrow=n,ncol=1)
> x2=cbind(x1,x)
> x3=cbind(x2,x^2)
> x4=cbind(x3,x^3)
> x5=matrix(model.matrix(~0+factor(x)),nrow=n)
> I=diag(rep(1,n))
```

Creating the Projection Matrices in R

```
> library(MASS)
> proj=function(x) {
+   x%*%ginv(t(x)%*%x)%*%t(x)
+ }
>
> p1=proj(x1)
> p2=proj(x2)
> p3=proj(x3)
> p4=proj(x4)
> p5=proj(x5)
```



A handwritten diagram in purple ink is positioned to the right of the R code. It consists of a vertical ellipsis (three dots) between the labels P_1 and P_5 . A diagonal line segment connects the label P_1 to the first line of code (`p1=proj(x1)`). A horizontal line segment connects the label P_5 to the last line of code (`p5=proj(x5)`).

Computing the Sums of Squares in R

```
> t(y) %*% (p2-p1) %*% y  
      [,1]  
[1,] 43.2  
> t(y) %*% (p3-p2) %*% y  
      [,1]  
[1,] 42  
> t(y) %*% (p4-p3) %*% y  
      [,1]  
[1,] 0.3  
> t(y) %*% (p5-p4) %*% y  
      [,1]  
[1,] 2.1  
> t(y) %*% (I-p5) %*% y  
      [,1]  
[1,] 7.48  
> t(y) %*% (I-p1) %*% y  
      [,1]  
[1,] 95.08
```

$y^T (P_2 - P_1) y$
 $y^T (P_3 - P_2) y$

sizable reductions in SSE

reductions are neglig compared to SLR & quadratic regression model

end lecture 14