



# Factorial Treatment Structure: Part II

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# Individual Analyses: Example (Quinn & Keough, 2000)

- Example: Fecundity of limpets.
- Response: Average number of eggs per snail.
- Factors:
  - A: Season (2 levels: Spring / Summer)
  - B: Population density (3 levels: 6 / 12 / 24 limpets under wire mesh guard)
- Per treatment combination: 3 observations



	6	12	24
<i>Spring</i>	1.17	1.50	0.67
	0.50	0.83	0.67
	1.67	1.00	0.75
<i>Summer</i>	4.00	3.33	2.54
	3.83	2.58	1.83
	3.83	2.75	1.63

# Individual Analyses: Example (Quinn & Keough, 2000)

## ■ Output of full model

```
> fit <- aov(y ~ season * density, data = snails)
> summary(fit)
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
season	1	17.131	17.131	119.373	1.36e-07	***
density	2	4.001	2.001	13.940	0.000742	***
season:density	2	1.689	0.845	5.885	0.016552	*
Residuals	12	1.722	0.144			

Need individual models per season

## ■ Output of individual models

```
> fit.spring <- aov(y ~ density, data = subset(snails, season == "Spring"))
> fit.summer <- aov(y ~ density, data = subset(snails, season == "Summer"))
>
> summary(fit.spring)
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
density	2	0.3445	0.1722	1.104	0.391
Residuals	6	0.9361	0.1560		

```
> summary(fit.summer)
```

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
density	2	5.346	2.673	20.41	0.00211 **
Residuals	6	0.786	0.131		

# Individual Analyses: Example (Quinn & Keough, 2000)

- Can improve tests by “re-using”  $MS_E$  with corresponding df’s of **full model** (a better, more precise estimate of  $\sigma$ ).
- The more df’s we have for  $MS_E$ , the **more powerful** the test will be because quantiles of the  $F$ -distribution are (much) **smaller**, as seen in previous lecture.
- E.g., for the spring data-set, the “better”  $F$ -test is:

$$\frac{MS_{\text{density}}}{MS_E} = \frac{0.1722}{0.144} = 1.2$$

- We have to use an  $F_{2,12}$ -distribution, hence the  $p$ -value is

```
> pf(1.2, 2, 12, lower.tail = FALSE)
[1] 0.334898
```

(which is only a slight improvement here).

# Individual Analyses

- Similarly, for testing contrasts etc. we can make use of the global  $\sigma$  estimate given by  $MS_E$  of the full model.
- This means: whenever we have an  $MS_E$  in our formulas for the **individual** models, we can “plug in” the **global** estimate with the corresponding degrees of freedom.
- This is especially useful if the error df's of the individual model are small (say below 10).

# Single Replicates

# Single Replicates

- If we have a factorial experiment with only **one** observation per factor-level combination we **cannot fit a full model** anymore.
- Reason: **Perfect fit**, all residuals are **zero** (or # parameters = # observations).
- Think of two-way ANOVA situation with no replicates. If we have no replicates we have no index  $k$ :

$$Y_{ij} = \mu + \alpha_i + \beta_j + \underbrace{(\alpha\beta)_{ij} + \epsilon_{ij}}$$

Cannot distinguish between these two terms in  $n = 1$  situation.

- Also: Remember factor  $n - 1$  for error in ANOVA table last time (i.e., we implicitly assumed  $n > 1$ ).

# Single Replicates

- We can of course still fit a model without interaction term, i.e. a main-effects model only (= additive effects).
- If there is an underlying interaction term, we get an error estimate that is biased upwards (because it contains the error **and** the interaction term).
- Tests will be **conservative** ( $p$ -values will be too large).
- See also R-File.



# Single Replicates

- Parameter estimates for main-effects model are as for previous model.
- ANOVA table now looks as follows:

Source	df	Sum of squares (SS)
A	$a - 1$	$\sum_{i=1}^a b \cdot \hat{\alpha}_i^2$
B	$b - 1$	$\sum_{j=1}^b a \cdot \hat{\beta}_j^2$
Error	$(a - 1) \cdot (b - 1)$	$\sum_{i=1}^a \sum_{j=1}^b (y_{ij} - \underbrace{(\hat{\mu} + \hat{\alpha}_i + \hat{\beta}_j)}_{\text{fitted value}})^2$
Total	$ab - 1$	$\sum_{i=1}^a \sum_{j=1}^b (y_{ij} - \bar{y}_{..})^2$

Annotations:

- A box labeled "#observations - 1" has an arrow pointing to the total df value  $ab - 1$ .
- A box labeled "df of total - sum(df above)" has an arrow pointing to the error df value  $(a - 1) \cdot (b - 1)$ .
- A box labeled "fitted value" has an arrow pointing to the term  $(\hat{\mu} + \hat{\alpha}_i + \hat{\beta}_j)$  in the error SS formula.

# Single Replicates

- If we have no replicates and **more than two factors** we would typically remove some of the higher-order interaction terms.
- This means: We put them into the error term (the df's of the error term will therefore increase).
- Often: Transformations of the response help getting rid of interactions, see blackboard.
- Alternative: Tukey one-degree-of freedom model for interaction (see next slide).

# Tukey One-Degree of Freedom Interaction

- Tukey's idea was to use only **one** additional parameter for the interaction term.
- For the two-factor model it is

$$Y_{ij} = \mu + \alpha_i + \beta_j + \underbrace{\lambda\alpha_i\beta_j}_{\text{interaction}} + \epsilon_{ij}$$

Here, interaction actually **is** the product of the main-effects!

- This is a **very special form of interaction**.
- Some add-on packages can test  $H_0: \lambda = 0$  (we will not discuss this any further).

# Contrasts for Factorial Data

# Contrasts for Factorial Data

- As we have already seen in the one-way ANOVA situation, **contrasts** can be used to get a more precise picture or to answer more specific question.
- Contrasts can of course also be used for factorial data structure.
- E.g., perform a **pairwise comparison between all possible treatment combinations** (think of having one “huge” cell means model with all the treatment combinations as levels).
- We can also construct “**main-effects**” **contrasts**.
- See R-File for examples.

# Unbalanced Data

# Unbalanced Data



- Up to now we assumed that our factorial data is **balanced**, i.e. all the “cells” contained the **same number of replicates** ( $= n$ ).
- This assumption is **crucial** for the **decomposition** of the total variability into the different sources.
- With balanced data we can estimate the effects of a factor by **ignoring** the remaining factors.
- Unbalanced data **destroy** these properties.
- Calculations are more involved. **No** independent estimates anymore and **no** unique decomposition either.



# Unbalanced Data: Toy Example

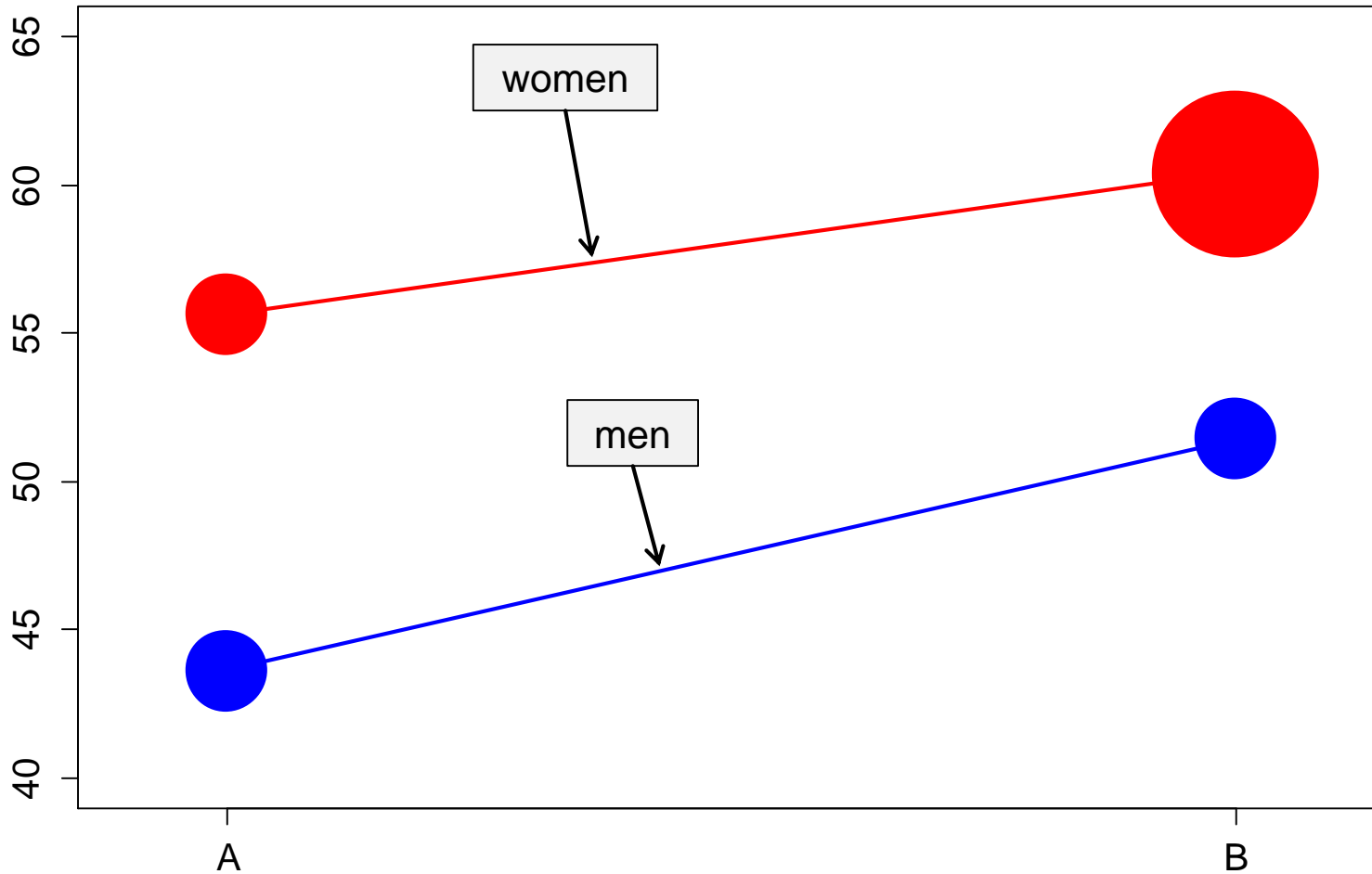
Consider the following data-set for a sports experiment (running time in seconds for a specific track, unpaired data). In red: number of observations.

	<i>Energy Drink A</i>	<i>Energy Drink B</i>
<b>Men</b>	40.6, 49.7, 42.1, 42.2, 39.0, 44.2, 44.1, 43.1, 44.7, 46.3	49.7, 48.1, 49.7, 52.0, 51.5, 49.9, 55.6, 53.0, 53.5, 51.1
<b>Women</b>	55.7, 61.0, 58.0, 54.1, 51.9, 54.2, 54.4, 55.4, 55.4, 56.1	62.0, 60.3, 59.9, 61.2, 66.2, 56.5, 59.7, 63.0, 58.4, 61.7, 61.4, 62.6, 56.8, 55.2, 66.1, 60.6, 58.9, 59.1, 56.8, 62.5, 58.5, 61.3, 62.2, 62.5, 60.8, 57.1, 61.6, 65.9, 58.6, 60.6, 56.1, 53.6, 62.4, 62.2, 59.2, 62.9, 57.0, 58.5, 60.9, 63.4,



# Unbalanced Data: Toy Example

Interaction plot. Circle area proportional to sample size



# Unbalanced Data: Toy Example

- If we simply ignore the “gender” structure and estimate the drink effect we are estimating the wrong effect.
- Why? We have **more women** with **Energy Drink *B***.
- Gender and Energy Drink are **correlated** in this example.
- When moving from Drink *A* to Drink *B* we are also moving to a different gender structure.
- In the **balanced** case this would **not** happen.

# Unbalanced Data: Toy Example

- Hence, we can't estimate the parameters one at a time anymore.
- Parameters have to be estimated **simultaneously** using the principle of least squares (no problem for the computer).
- Parameter estimates estimate the “right thing”.
- Similarly: Sum of squares **cannot** be partitioned into different sources anymore.
- Note: There are unbalanced situations that are still “nice”.

# Unbalanced Data

- “Solution” to decomposition of sum of squares: use **model comparison** approach.
- Remember:  $SS_B$  (and the other sum of squares) can be thought of as the **reduction of residual sum of squares** when adding the factor  $B$  to our model.
- In the balanced case, it does **not** matter whether we have  $A$  (or  $AB$ ) in our model or not:  $SS_B$  is always the same.
- For unbalanced data,  $SS_B$  depends on the remaining terms in the model.  $\mu$
- Notation:  $SS(B \mid 1, A)$  is **reduction of error sum of squares** when comparing the models  $(1, A, B)$  with  $(1, A)$ .

# Unbalanced Data

- **Model terms** in two-way ANOVA situation:  $(1, A, B, AB)$
- Could have a look at
  - $SS(A | 1)$
  - $SS(B | 1, A)$
  - $SS(A | 1, B)$
  - $SS(AB | 1, A, B)$
  - $SS(A | 1, B, AB)$
  - ...
- $SS_E$  or  $MS_E$  are typically taken from the **full** model (including **all** terms).



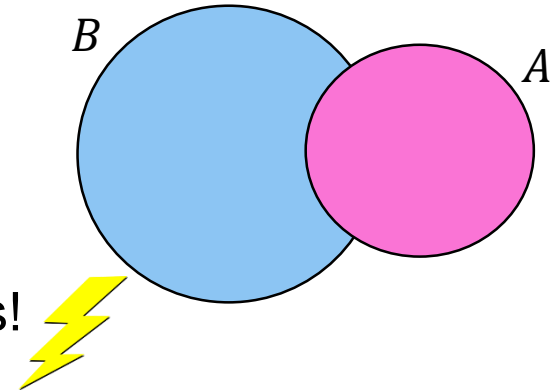
# Different Types of Sum of Squares

Terminology from SAS

## ■ Type I: Sequential sum of squares

Sequentially build up model

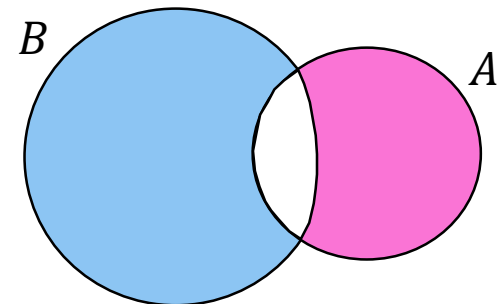
- $SS(A | 1)$
- $SS(B | 1, A)$
- $SS(AB | 1, A, B)$
- Hence: Depends on **ordering** of factors!
- R: `aov`



## ■ Type II: Hierarchical / partially sequential approach

Control for the influence of the largest hierarchical model not including the term of interest.

- $SS(A | 1, B)$
- $SS(B | 1, A)$
- $SS(AB | 1, A, B)$
- R: Function `Anova` in package `car`.



# Different Types of Sum of Squares

- **Type III: Fully adjusted / marginal** approach

Control for **all** other terms

- $SS(A \mid 1, B, AB)$  (meaningful?)
- $SS(B \mid 1, A, AB)$  (meaningful?)
- $SS(AB \mid 1, A, B)$
- R: `drop1`
- Typically the preferred type.

# In R...

or aov



## R FAQ 7.18: Why does the output from `anova()` depend on the order of factors in the model?

- In a model such as `~A+B+A:B`, R will report the difference in sums of squares between the models `~1`, `~A`, `~A+B` and `~A+B+A:B`. If the model were `~B+A+A:B`, R would report differences between `~1`, `~B`, `~A+B`, and `~A+B+A:B`. In the first case the sum of squares for A is comparing `~1` and `~A`, in the second case it is comparing `~B` and `~B+A`. In a non-orthogonal design (i.e., most unbalanced designs) these comparisons are (conceptually and numerically) different.
- Some packages report instead the sums of squares based on comparing the full model to the models with each factor removed one at a time (the famous ‘Type III sums of squares’ from SAS, for example). These do not depend on the order of factors in the model. **The question of which set of sums of squares is the Right Thing provokes low-level holy wars on R-help from time to time.**
- **There is no need to be agitated about the particular sums of squares that R reports. You can compute your favorite sums of squares quite easily. Any two models can be compared with `anova(model1, model2)`, and `drop1(model1)` will show the sums of squares resulting from dropping single terms.**



# Recommendations / Comments

- With **balanced data** we always get the **same result**, no matter what type we use.
- For main-effects only models, type II and type III coincide.
- Type I is useful if there is some intrinsic ordering of the terms.
- If there is a significant interaction, tests of the corresponding main-effects are typically difficult to interpret.