Displays for Statistics 5303

Lecture 30

November 13, 2002

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Class Web Page

http://www.stat.umn.edu/~kb/classes/5303

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Nested Random Effects Designs

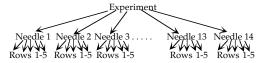
We have looked at the one-factor random effect design as a particular case of random effect factorial designs.

But it is also a particular case of a socalled nested design:

Example in the sample exam

Cmd> data <- read("","stomata")
stomata 70 2
) Data on the number of stomata in 5 rows randomly selected
) on each of 14 randomly selected evergreen needles
) Col. 1: Needle number (1 - 14)
) Col. 2: stomata per cm.
) Source of the data is unknown
Read from file "TP1:Stat5303:Data:stomata.dat"
Cmd> makecols(data,needle,stomata)
Cmd> needle <- factor(needle)

Here needles were first randomly selected. Then, within each needle, 5 rows were randomly selected. It's a sort of tree-like structure



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You can define a factor for row, that is nested within each needle:

The model we have have previously used for this has been

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

An equivalent model would be

$$y_{ij} = \mu + \alpha_i + \beta_{i(i)} + \widetilde{\epsilon}_{ij}$$

where $\beta_{j(i)}$ is the random effect of row j within needle i and $\beta_{j(i)}$ + $\widetilde{\epsilon}_{ij}$ = $\epsilon_{ij}.$

The notation j(i) is intended to convey that j has a different meaning for each i, that is for each needle.

Here how you would analyze it with the nested model.

 Cmd> anova("stomata=needle+row.needle")

 Model used is stomata=needle+row.needle

 DF
 SS
 MS

 CONSTANT
 1
 1.1762e+06
 1.1762e+06

 needle
 13
 2111.1
 162.4

 row.needle
 56
 2667.2
 47.629

 ERROR1
 0
 0
 undefined

Since there is only 1 measurement per row, there are no error d.f.

row.needle does not signify an interaction here but a nesting of row Within needle.

Q. How can you tell from that row.needle doesn't indicate interaction?

A. From the absence of a line for row.

The degrees of freedom for needle is $DF_{\Delta} = a - 1 = 14 - 1 = 13$.

The degrees of freedom for row.needle (row nested in needle) is

$$DF_{R(A)} = a(b-1) = 14(5-1) = 56.$$

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If the experimenter made n=3 quick counts for each row of each needle, so there were $14\times5\times3=210$ values, then an appropriate model would be

$$y_{ijk} = \mu + \alpha_i + \beta_{j(i)} + \epsilon_{k(ij)}$$

where again, the notation k(ij) is meant to indicate that the level k is specific to the particular row i within needle j.

When you have an experiment that consists of randomly selecting

- a entities of type A (needles say)
- b entities of type B (rows, say) within each type A entity
- c entities of type C (random places in a row, say) within each type B entity
- Making n measurements y_{ijk 1} on each type C entity

the nested model would be

$$y_{ijl} = \mu + \alpha_i + \beta_{j(i)} + \delta_{k(ij)} + \epsilon_{k(ijk)}$$

$$A \quad B(A) C(AB) \quad Error(ABC)$$

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Here is an example. An experiment was designed to study the sources of variability in measurements of the fat content of dried whole eggs.

All material to be analyzed came from a single well mixed can.

- 24 samples from the can were packaged for sending to labs.
- 4 samples were sent to each of a = 6 labs (A) which can be considered a random sample of labs.
- At each lab, each of b = 2 analysts (B) on the staff were given c = two samples (C) to analyze.
- Each analyst made n = 2 determination of the fat content of the sample.

Note there are no symbols containing two or more letters. This is characteristic of fully nested designs.

The α_i , $\beta_{j(i)}$, $\delta_{k(ij)}$ and $\epsilon_{\ell(ijk)}$ are assumed to be random variables with

- Zero means $(\mu_{x} = \mu_{y} = \mu_{y} = 0)$
- Variances $\sigma_{\alpha}^{\ 2}$, $\sigma_{\beta}^{\ 2}$, $\sigma_{\sigma}^{\ 2}$, and σ^{2} are constant

For tests and confidence intervals you assume

All random variables are normal

The parameters are μ and the variance components σ_{χ}^2 , σ_{κ}^2 , σ_{χ}^2 , and σ^2

The variance of a single observation is
$$V(y_{ijk0}) = \sigma_{\alpha}^2 + \sigma_{\beta}^2 + \sigma_{\alpha}^2 + \sigma^2$$

The variance of the grand mean \overline{y} ... is $V(\overline{y}$...) = $\sigma_{\pi}^2/a + \sigma_{\pi}^2/ab + \sigma_{\pi}^2/abc + \sigma^2/abcn$

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Model used is y = lab+ analyst.lab+sample.analyst.lab 7.2075 7.2075 1001.62131 4.3928e-21 CONSTANT 1 0.088605 0.44302 5.4864e-06 12.31338 lab lab.analyst 0.24748 0.041246 5.73191 0.00081653 lab.analyst. 0.1599 0.013325 1.85177 0.096155 0.0071958 0.1727

Each SS is computed from the means at that level.

Example:

$$SS_{B(A)} = nc \sum_{1 < i < a} \sum_{1 < j < b} (\overline{y_{ij \bullet \bullet}} - \overline{y_{i \bullet \bullet \bullet}})^2$$

nc = number of values averaged to compute $\overline{y_{ij+}}$.

Numerical check

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The skeleton ANOVA is

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	Source	DF	EMS		
	А	a-1	$\sigma^2 + n\sigma_{g}^2 + nc\sigma_{g}^2 + nbc\sigma_{\alpha}^2$		
	B(A)	a(b-1)	$\sigma^2 + n\sigma_{\sigma}^2 + nc\sigma_{\beta}^2$		
	C(AB)	ab(c-1)	$\sigma^2 + n\sigma_{\sigma}^2$		
	Error	abc(n-1)	σ^2		

In this case

Cmd>
$$vector(a-1,a^*(b-1),a^*b^*(c-1),a^*b^*c^*(n-1))$$
 (1) 5 6 12 24

Source	DF	EMS
А	5	$\sigma^2 + 2\sigma_{g}^2 + 4\sigma_{g}^2 + 8\sigma_{\alpha}^2$
B(A)	6	$\sigma^2 + 2\sigma_{g}^2 + 4\sigma_{g}^2$
C(AB)	12	$\sigma^2 + 2\sigma_{g}^2$
Error	24	σ^2

From this estimates of the o2's are $\hat{\sigma}_{\alpha}^{2} = (MS_{\Delta} - MS_{B(\Delta)})/nbc,$ $\hat{\sigma}_{\beta}^{2}$ = (MS_{B(A)} - MS_{C(AB)})/nc, etc.

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You can use this output to compute approximate confidence intervals using χ^2 (assuming normality of effects).

```
Cmd> estimate <- vcomp[1,1]; estimate</pre>
       Estimate
lab
Cmd> eps <- .025; chisqpts <- invchi(vector(1-eps/2,eps/2),df)
Cmd> vector(df*estimate/chisqpts) # 95% confidence interval
(1) 0.0028102
                    0.14089
```

Cmd> df <- vcomp[1,3]; df

3.5755

lab

ems() can compute these formulas:

```
 \begin{array}{ll} {\it Cmd>\ ems("y=lab+analyst.lab+sample.analyst.lab", \backslash \\ & vector("lab","analyst","sample"))} \\ {\it EMS(CONSTANT)=V(ERROR1)+2V(lab.exper.sample)+4V(lab.exper)+8V(lab)+48Q(CONSTANT)} \end{array} 
EMS(lab) = V(ERROR1) + 2V(lab.exper.sample) + 4V(lab.exper) +
8V(lab)
EMS(lab.exper) = V(ERROR1) + 2V(lab.exper.sample) +
4V(lab.exper)
EMS(lab.exper.sample) = V(ERROR1) + 2V(lab.exper.sample)
EMS(ERROR1) = V(ERROR1)
```

As before, v stands for the variance of a random effect and o stands for a contribution from one or more fixed parameters. Only μ is fixed here and $O(CONSTANT) = U^2$.

```
\label{eq:cmd} $$\operatorname{Cmd}> \operatorname{sigmasqA\_hat} <- (MS[2] - MS[3])/(n*b*c); \ \operatorname{sigmasqA\_hat} $$
(1) 0.0059199
Cmd> sigmasgB hat <- (MS[3] - MS[4])/(n*b); sigmasgB hat
      0.0069802
Cmd> sigmasqC_hat <- (MS[4] - MS[5])/n; sigmasqC_hat
(1) 0.0030646
Cmd> sigmasq_hat <- MS[5]; sigmasq_hat
      ERROR1
   0.0071958
Cmd> vcomp <- varcomp("y=exper + lab + lab.exper +
    lab.exper.sample", vector("lab", "sample"))
Cmd> vcomp
                       Estimate
                                                     3.5755
                                   0.0070378
lab
                       0.00941
                     0.0088221
                                   0.0078058
exper.lab
exper.lab.sample
                     0.0030646
                                   0.0029115
                                                     2.2158
                      0.0071958
                                   0.0020773
```

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Suppose the two experimenters are selected so that one is inexperienced (< 2 years in the lab) and the other is experienced (≥ 2 years).

Experience is a factor that is crossed with lab and sample is nested within combinations of lab and experience.

```
Cmd> exper <- analyst # experience factor
Cmd> anova("y=exper + lab + lab.exper + lab.exper.sample", \
Model used is y=exper + lab + lab.exper + lab.exper.sample
                                   MS F P-value
7.2075 1001.62131 4.3928e-21
                           SS
                      7.2075
CONSTANT
               1
                   0.0044083
                                0.0044083
                                               0.61262
                                                           0.44146
exper
                      0.44303
                                                        5.4864e-06
lab
                                 0.088605
                                              12.31338
exper.lab
                                                       0.00046361
               5
                                 0.048613
                                               6.75576
                     0.24307
                                 0.013325
 sample
              12
                      0.1599
                                               1.85177
                                                          0.096155
                                0.0071958
                       0.1727
```

The model

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y=exper+lab+lab.exper+lab.exper.sample specifies that exper and lab are crossed and not nested so that lab.exper is a random interaction term.

sample is nested within lab.exper.

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The mathematical model is

```
\begin{array}{l} y_{ijkl} = \mu + \alpha_i + \beta_j + \alpha \beta_{ij} + \mathcal{T}_{k(ij)} = \epsilon_{ijkl} \\ \text{where } \alpha_i \ \alpha \beta_{ij}, \ \mathcal{T}_{k(ij)} \ \text{and} \ \epsilon_{ijkl} \ \text{are random} \\ \text{variables with zero means and variances} \\ \sigma_{\alpha}^{\ 2}, \ \sigma_{\alpha\beta}^{\ 2}, \ \sigma_{\sigma}^{\ 2} \ \text{and} \ \sigma^2. \end{array}
```

Note that, because exper is a fixed factor, Q(exper) and not V(exper) is part of EMS(exper).

There is no line for exper.

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