

# STAT 516 Lec 09

Randomized complete block designs

*experimental  
design*

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# Nitrogen fertilization data from Kuehl (2000)

Six nitrogen timing schedules in each of four plots in a field of wheat. Response is the nitrate content in stem tissue samples.

**Display 8.1 Arrangement of Experimental Plots for the Wheat Experiment in a Randomized Complete Block Design**

Block	Plot 1	Plot 2	Plot 3	Plot 4	Plot 5	Plot 6
Block 1	1	2	3	4	5	6
	40.89	37.99	37.18	34.98	34.89	42.07
Block 2	1	3	4	6	5	2
	41.22	49.42	45.85	50.15	41.99	46.69
Block 3	6	3	5	1	2	4
	44.57	52.68	37.61	36.94	46.65	40.23
Block 4	2	4	6	5	3	1
	41.90	39.20	43.29	40.45	42.91	39.97

$$a = 6$$

$$b = 4$$

$$a * b = 24$$

Source: Dr. T. Doerge, Department of Soil and Water Science, University of Arizona.

```
ntr <- c(40.89,37.99,37.18,34.98,34.89,42.07,  
        41.22,49.42,45.85,50.15,41.99,46.69,  
        44.57,52.68,37.61,36.94,46.65,40.23,  
        41.90,39.20,43.29,40.45,42.91,39.97)  
block <- as.factor(c(1,1,1,1,1,1,2,2,2,2,2,2,3,3,3,3,3,3,4,4,4,4,4,4))  
trt <- as.factor(c(2,5,4,1,6,3,1,3,4,6,5,2,6,3,5,1,2,4,2,4,6,5,3,1))
```

# Randomized complete block design (RCBD)

- ▶ EUs belong to blocks—groups of EUs homogeneous in some way.
- ▶ Each EU in a block is randomly assigned to a fixed treatment.
- ▶ All treatments appear exactly once<sup>1</sup> in each block.
- ▶ Purpose is to capture the between-block variability among the EUs.
- ▶ This helps us detect treatment effects with greater power.

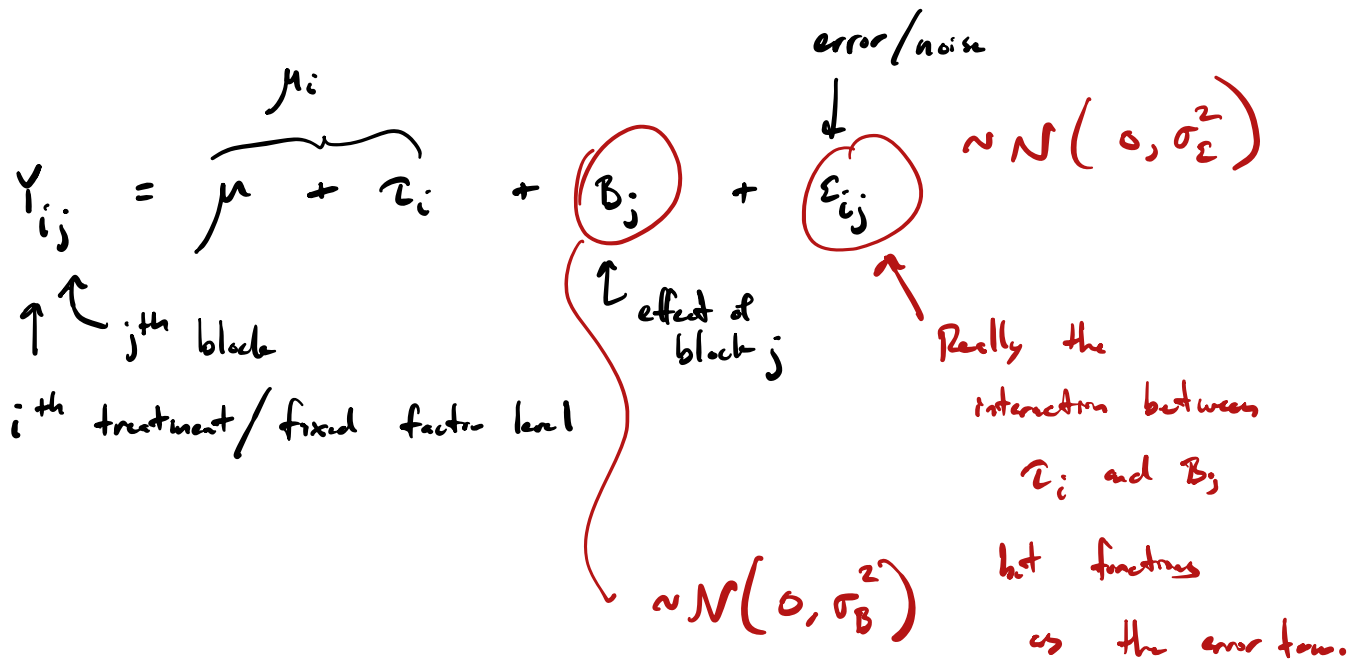
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<sup>1</sup>Can have replication, but many RCBDs do not.

Recall 2-way factorial design

$$Y_{ijk} = \mu + \tau_i + \beta_j + (\tau\beta)_{ij} + \epsilon_{ijk}$$

---



# Treatment effects model for the RCBD

Assume

$n = 1$

there is no little  $n$ , because no replication (yet)

$$Y_{ij} = \mu + \tau_i + B_j + \varepsilon_{ij}, \quad \text{for } i = 1, \dots, a, \quad j = 1, \dots, b,$$

where



- ▶  $Y_{ij}$  is the response of the EU in block  $j$  receiving treatment  $i$ .
- ▶ the  $\tau_i$  are the fixed effects of the treatment.
- ▶ the  $B_j$  are independent  $\text{Normal}(0, \sigma_B^2)$  random block effects.
- ▶ the  $\varepsilon_{ij}$  are independent  $\text{Normal}(0, \sigma_\varepsilon^2)$  error/interaction terms.
- ▶  $\mu$  is an overall or baseline mean.

Define the cell means as

$$\mu_i = \mu + \tau_i, \quad i = 1, \dots, a.$$

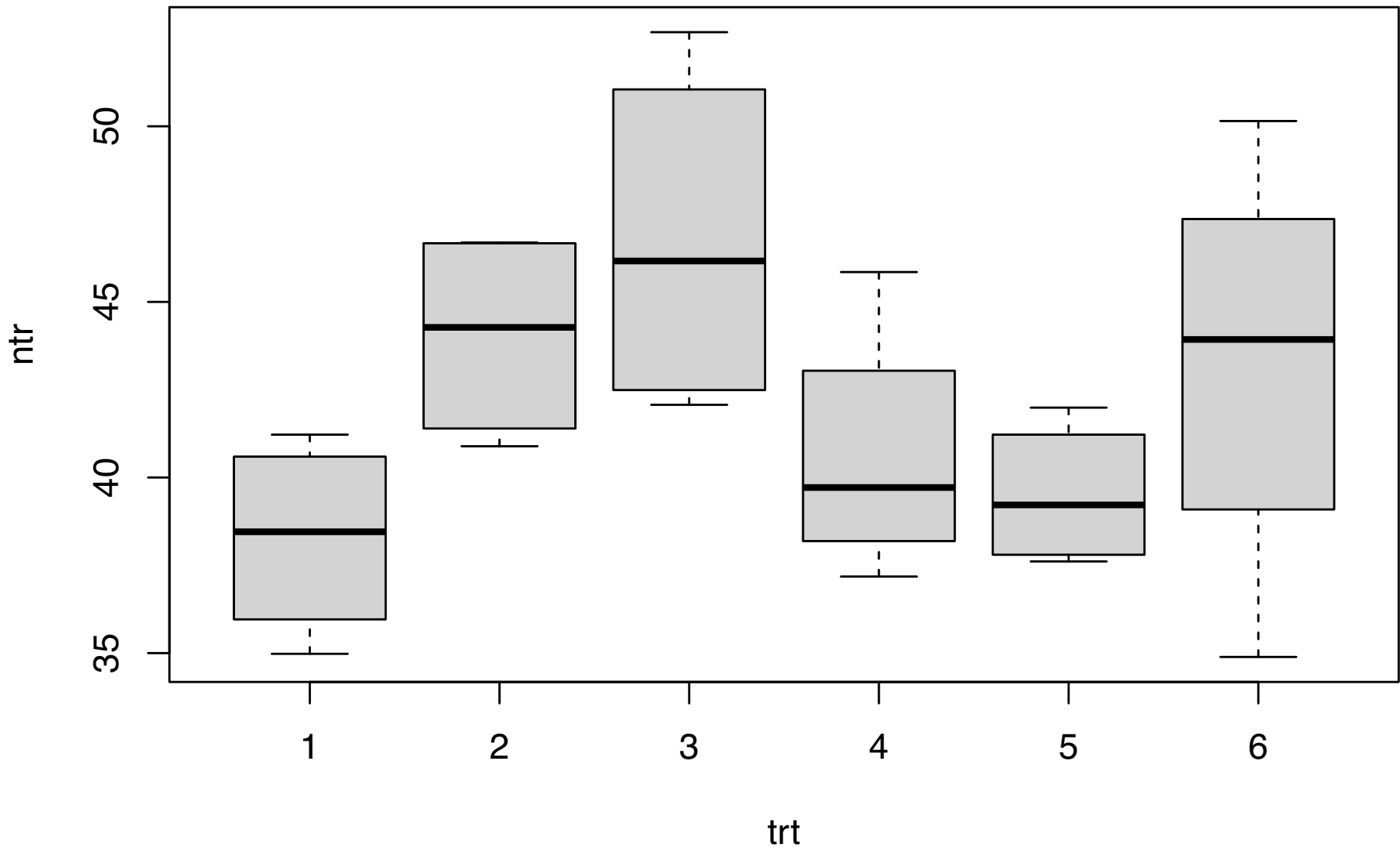
# Goals in the RCBD

In the randomized complete block design we wish to

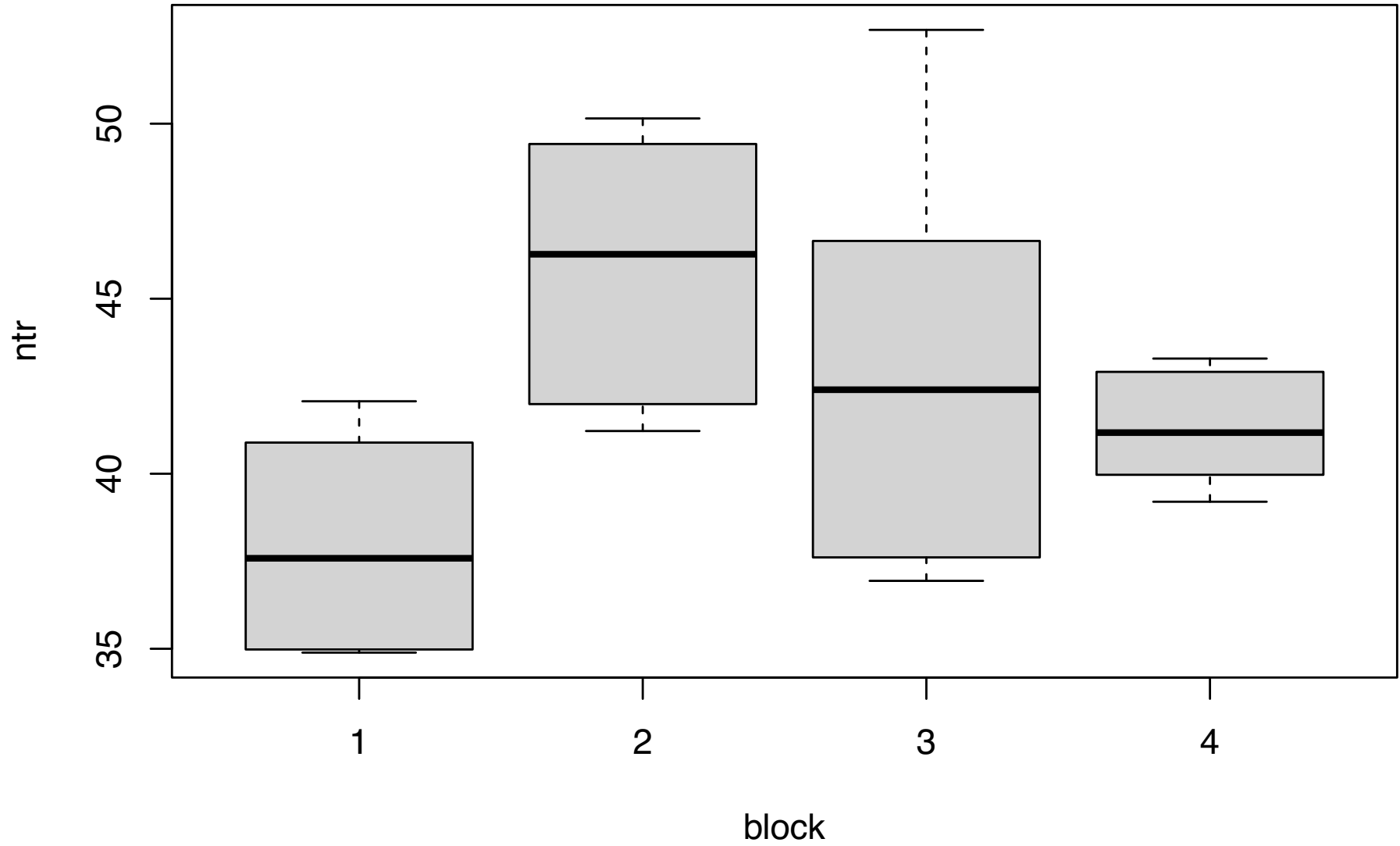
1. Visualize the data.
2. Decompose the variability in the  $Y_{ij}$  into its sources.
3. Estimate the variance components  $\sigma_B^2$  and  $\sigma_\epsilon^2$ .
4. Test whether the treatment has any effect.
5. Make comparisons between treatment means.
6. Check whether the model assumptions are satisfied.

# Nitrogen fertilization data (cont)

```
boxplot(ntr~trt)
```



```
boxplot(ntr~block)
```



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

Estimated mean of combination  $i, j$ .

$$\begin{aligned} \rightarrow \bar{y}_{..} + (\bar{y}_{i.} - \bar{y}_{..}) + (\bar{y}_{.j} - \bar{y}_{..}) \\ = \bar{y}_{i.} + \bar{y}_{.j} - \bar{y}_{..} \end{aligned}$$

SS for RCBD

$$SS_{Tot} = \sum_{i=1}^a \sum_{j=1}^b (y_{ij} - \bar{y}_{..})^2$$

$a = \#$  levels fixed factor

$b = \#$  blocks

$$SS_{Treat} = \sum_{i=1}^a b (\bar{y}_{i.} - \bar{y}_{..})^2$$

$$SS_{blocks} = \sum_{j=1}^b a (\bar{y}_{.j} - \bar{y}_{..})^2$$

$$SS_{Error} = \sum_{i=1}^a \sum_{j=1}^b \left( y_{ij} - (\bar{y}_{i.} + \bar{y}_{.j} - \bar{y}_{..}) \right)^2$$

estimated mean for  $i, j$  group

# Sums of squares for the RCBD

Sum of squares	Symbol	Formula
Total	$SS_{\text{Tot}}$	$\sum_{i=1}^a \sum_{j=1}^b (Y_{ij} - \bar{Y}_{..})^2$
Treatment	$SS_A$	$b \sum_{i=1}^a (\bar{Y}_{i.} - \bar{Y}_{..})^2$
Block	$SS_B$	$a \sum_{j=1}^b (\bar{Y}_{.j} - \bar{Y}_{..})^2$
Error	$SS_{\text{Error}}$	$\sum_{i=1}^a \sum_{j=1}^b (Y_{ij} - (\bar{Y}_{i.} + \bar{Y}_{.j} - \bar{Y}_{..}))^2$

- ▶ We can make the decomposition  $SS_{\text{Tot}} = SS_A + SS_B + SS_{\text{Error}}$ .
- ▶ The  $SS_{\text{Error}}$  is really the interaction sum of squares  $SS_{AB}$ .
- ▶ But without replication, we cannot estimate an interaction.
- ▶ So the interaction serves as the error term.

# ANOVA table for RCBD

$$a-1 + b-1 + (a-1)(b-1) = \cancel{a-1} + \cancel{b-1} + ab - \cancel{a} - \cancel{b} + 1 = ab - 1.$$

add up to total

Really the interaction AB

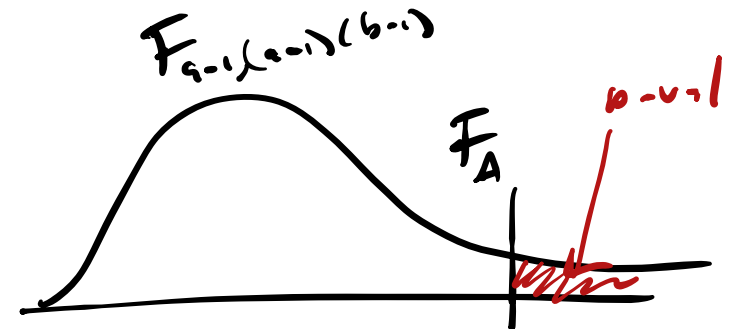
Source	Df	SS	MS	F value
A	$a-1$	$SS_A$	$MS_A$	$F_A = MS_A / MS_{Error}$
B	$b-1$	$SS_B$	$MS_B$	$F_B = MS_B / MS_{Error}$
Error	$(a-1)(b-1)$	$SS_{Error}$	$MS_{Error}$	
Total	$ab-1$	$SS_{Tot}$		

# obs.

1. Reject  $H_0: \mu_1 = \dots = \mu_a$  if  $F_A > F_{a-1, (a-1)(b-1), \alpha}$
2. Reject  $H_0: \sigma_B^2 = 0$  if  $F_B > F_{b-1, (a-1)(b-1), \alpha}$

pv. =  $1 - pf(F_A, a-1, (a-1)(b-1))$

$$Y_{ij} = \underbrace{\mu + \tau_i + \beta_j}_{\mu_i} + \epsilon_{ij}$$



# Nitrogen fertilization data (cont)

```
lm_out <- lm(ntr ~ trt + block) # do not include the interaction
anova_out <- anova(lm_out)
anova_out
```

## Analysis of Variance Table

Response: ntr

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
trt	5	201.32	40.263	5.5917	0.004191	**
block	3	197.00	65.668	9.1198	0.001116	**
Residuals	15	108.01	7.201			

---

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

```

y <- ntr
y.. <- predict(lm(ntr ~ 1))
yi. <- predict(lm(ntr ~ trt))
y.j <- predict(lm(ntr ~ block))

SSA <- sum((yi. - y..)^2)
SSB <- sum((y.j - y..)^2)
SSE <- sum((y - (yi. + y.j - y..))^2)

a <- 6
b <- 4

MSA <- SSA/(a-1)
MSB <- SSB/(b-1)
MSE <- SSE/((a-1)*(b-1))

FA <- MSA / MSE
FB <- MSB / MSE

pA <- 1 - pf(FA, a-1, (a-1)*(b-1))
pB <- 1 - pf(FB, b-1, (a-1)*(b-1))

```

# Expected mean squares in the RCBD

Estimate  $\sigma_{\epsilon}^2$ ,  $\sigma_B^2$

Source	Df	Expected mean square
A	$a - 1$	$\mathbb{E}MS_A = b\theta_A^2 + \sigma_{\epsilon}^2$
B	$b - 1$	$\mathbb{E}MS_B = a\sigma_B^2 + \sigma_{\epsilon}^2$
Error	$(a - 1)(b - 1)$	$\sigma_{\epsilon}^2 = \mathbb{E}MS_{Error}$

In the above  $\theta_A^2 = (a - 1)^{-1} \sum_{i=1}^a (\mu_i - \bar{\mu})^2$ .

$$\hat{\sigma}_B^2 = \frac{1}{a} [MS_B - MS_{Error}]$$

$$\hat{\sigma}_{\epsilon}^2 = MS_{Error}$$

# Method of moments for variance components in RCBD

Equating  $MS_B$  and  $MS_{\text{Error}}$  with their expectations gives

$$\blacktriangleright \hat{\sigma}_\varepsilon^2 = MS_{\text{Error}}.$$

$$\blacktriangleright \hat{\sigma}_B^2 = \frac{MS_B - MS_{\text{Error}}}{a}$$

May obtain  $\hat{\sigma}_B^2 < 0$ , so one should use REML estimation.

# Nitrogen fertilization data (cont)

Obtain REML estimators of  $\sigma_B^2$  and  $\sigma_\varepsilon^2$  on the fertilization data.

```
library(lmerTest) # first time run install.packages("lmerTest")
lmer_out <- lmer(ntr ~ trt + (1|block))
lmer_out
```

Linear mixed model fit by REML ['lmerModLmerTest']

Formula: ntr ~ trt + (1 | block)

REML criterion at convergence: 101.5658

Random effects:

Groups	Name	Std.Dev.
block	(Intercept)	3.122
	Residual	2.683

Number of obs: 24, groups: block, 4

Fixed Effects:

(Intercept)	trt2	trt3	trt4	trt5	trt6
38.278	5.755	8.492	2.337	1.232	4.947

Obtain MoMs estimators for  $\sigma_B^2$  and  $\sigma_\varepsilon^2$  on the fertilization data.

```
sg_B <- sqrt((MSB - MSE)/a)
sg_e <- sqrt(MSE)
```

We have  $\hat{\sigma}_B = 3.122$  and  $\hat{\sigma}_\varepsilon = 2.683$ .

# Variances of some means and difference in means

Contrast	Variance	MoM variance estimator
$\bar{Y}_i.$	$\frac{1}{b}(\sigma_B^2 + \sigma_\varepsilon^2)$	$\frac{2}{ab}[\text{MS}_B + (b - 1) \text{MS}_{\text{Error}}]$
$\bar{Y}_i. - \bar{Y}_{i'}. $	$\frac{2}{b}\sigma_\varepsilon^2$	$\frac{2}{b} \text{MS}_{\text{Error}}$

# Some (unadjusted) CIs in RCB split plot design

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Target	$(1 - \alpha)100\%$ confidence interval
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$$\mu_i \quad \bar{Y}_{i.} \pm t_{\nu^*, \alpha/2} \sqrt{MS_B + (b - 1) MS_{\text{Error}}} \sqrt{\frac{2}{ab}}$$

$$\mu_i - \mu_{i'} \quad \bar{Y}_{i.} - \bar{Y}_{i'.} \pm t_{(a-1)(b-1), \alpha/2} \sqrt{MS_{\text{Error}}} \sqrt{\frac{2}{b}}$$

---

In the above  $\nu^* = \frac{MS_B + (b - 1) MS_{\text{Error}}}{\frac{MS_B^2}{(b-1)} + \frac{(b-1)^2 MS_{\text{Error}}^2}{(a-1)(b-1)}}$  à la Satterthwaite<sup>2</sup>.

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<sup>2</sup>a degrees of freedom approximation when one has not exactly a t-distribution.

# Nitrogen fertilization data (cont)

$$\begin{aligned} & \textcircled{a=6} \\ & \binom{6}{2} = \frac{6!}{2! \cdot 4!} \\ & = \frac{6 \cdot 5}{2} \\ & = 15 \end{aligned}$$

Unadjusted CIs with `ls_means()` from R package `lmerTest`

```
ls_means(lmer_out)
```

Least Squares Means table:

	Estimate	Std. Error	df	t value	lower	upper	Pr(> t )	
trt1	38.2775	2.0582	6.8	18.597	33.3789	43.1761	4.505e-07	***
trt2	44.0325	2.0582	6.8	21.393	39.1339	48.9311	1.767e-07	***
trt3	46.7700	2.0582	6.8	22.724	41.8714	51.6686	1.179e-07	***
trt4	40.6150	2.0582	6.8	19.733	35.7164	45.5136	3.033e-07	***
trt5	39.5100	2.0582	6.8	19.196	34.6114	44.4086	3.646e-07	***
trt6	43.2250	2.0582	6.8	21.001	38.3264	48.1236	2.000e-07	***

---

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

Confidence level: 95%

Degrees of freedom method: Satterthwaite

```
ls_means(lmer_out, pairwise = TRUE)
```

Least Squares Means table:

	Estimate	Std. Error	df	t value	lower	upper	Pr(> t )	
trt1 - trt2	-5.7550	1.8974	15	-3.0330	-9.7993	-1.7107	0.0083887	**
trt1 - trt3	-8.4925	1.8974	15	-4.4758	-12.5368	-4.4482	0.0004443	***
trt1 - trt4	-2.3375	1.8974	15	-1.2319	-6.3818	1.7068	0.2369413	
trt1 - trt5	-1.2325	1.8974	15	-0.6496	-5.2768	2.8118	0.5257996	
trt1 - trt6	-4.9475	1.8974	15	-2.6075	-8.9918	-0.9032	0.0198026	*
trt2 - trt3	-2.7375	1.8974	15	-1.4427	-6.7818	1.3068	0.1696521	
trt2 - trt4	3.4175	1.8974	15	1.8011	-0.6268	7.4618	0.0918200	.
trt2 - trt5	4.5225	1.8974	15	2.3835	0.4782	8.5668	0.0308031	*
trt2 - trt6	0.8075	1.8974	15	0.4256	-3.2368	4.8518	0.6764617	
trt3 - trt4	6.1550	1.8974	15	3.2438	2.1107	10.1993	0.0054516	**
trt3 - trt5	7.2600	1.8974	15	3.8262	3.2157	11.3043	0.0016523	**
trt3 - trt6	3.5450	1.8974	15	1.8683	-0.4993	7.5893	0.0813776	.
trt4 - trt5	1.1050	1.8974	15	0.5824	-2.9393	5.1493	0.5689729	
trt4 - trt6	-2.6100	1.8974	15	-1.3755	-6.6543	1.4343	0.1891597	
trt5 - trt6	-3.7150	1.8974	15	-1.9579	-7.7593	0.3293	0.0691104	.

---

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

Confidence level: 95%

Degrees of freedom method: Satterthwaite

# Multiple comparisons of treatment means in the RCBD

- ▶ Tukey's for comparing all pairs of means among  $\mu_1, \dots, \mu_a$ :

$$\bar{Y}_{i.} - \bar{Y}_{i' .} \pm q_{a, (a-1)(b-1), \alpha} \sqrt{MS_{\text{Error}}} \sqrt{\frac{1}{b}}, \quad 1 \leq i < i' \leq a.$$

- ▶ Dunnett's for comparing  $\mu_2, \dots, \mu_a$  to a baseline  $\mu_1$ :

$$\bar{Y}_{i.} - \bar{Y}_{1.} \pm d_{a, (a-1)(b-1), \alpha} \sqrt{MS_{\text{Error}}} \sqrt{\frac{2}{b}}, \quad i = 2, \dots, a.$$

# Nitrogen fertilization data (cont)

Compare all pairs of fertilizers with Tukey's CIs for mean differences.

```
alpha <- 0.05
a <- 6
b <- 4
MSE <- anova_out$`Mean Sq`[3]
se <- sqrt(MSE) * sqrt(2/b)
me <- qtkey(1-alpha,a,(a-1)*(b-1)) / sqrt(2) * se
ntr_means <- aggregate(ntr, by = list(trt), mean)$x

CIs <- matrix(NA,choose(a,2),2)
comp <- numeric(choose(a,2))

k <- 1
for(i in 1:(a-1))
  for(j in (i+1):a){ # double loop takes us through all pairs

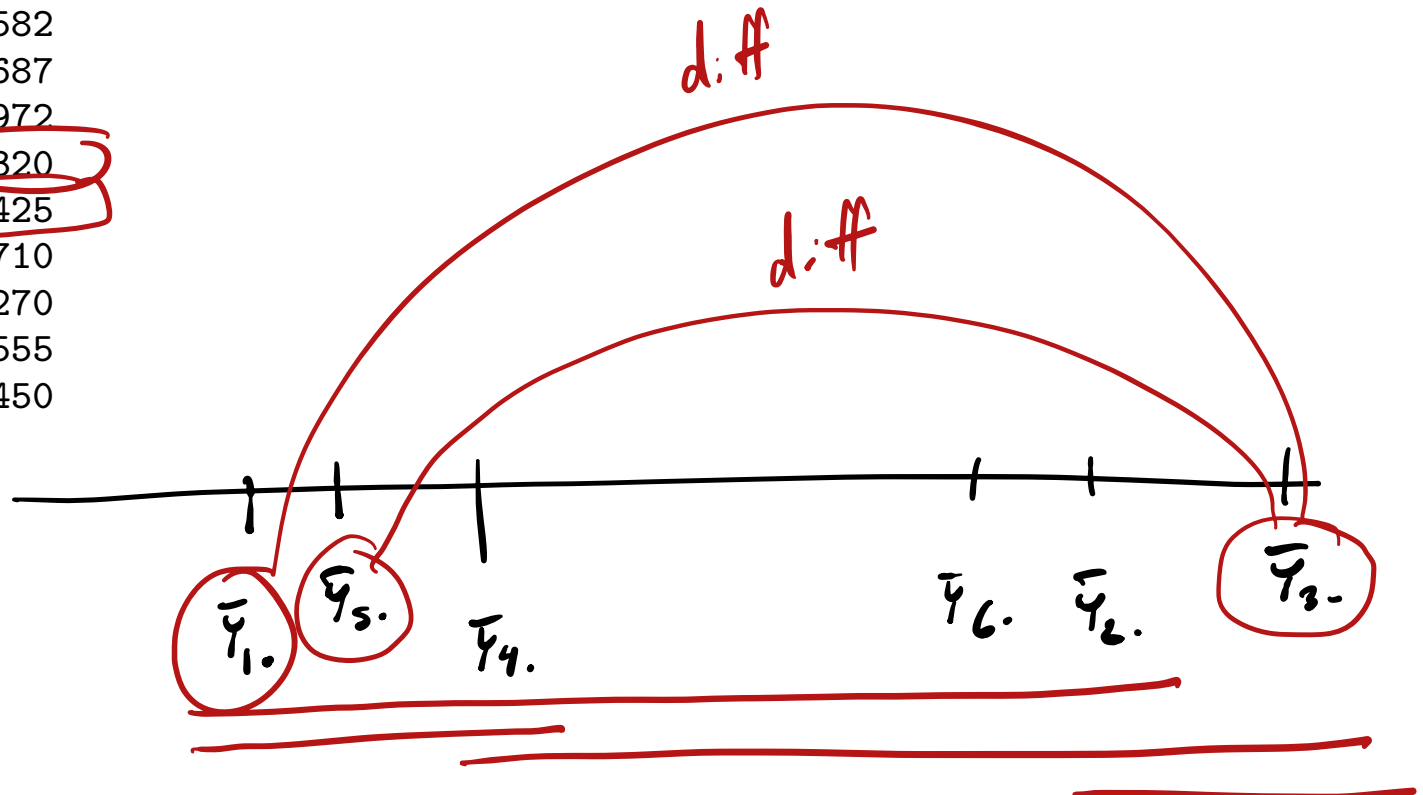
    dij <- ntr_means[i] - ntr_means[j]
    CIs[k,] <- c(dij - me, dij + me)
    comp[k] <- paste(i,"-",j)
    k <- k + 1

  }

colnames(CIs) <- c("lower","upper")
rownames(CIs) <- comp
```

`round(CIs,3)`

	lower	upper
1 - 2	-11.920	0.410
1 - 3	-14.657	-2.328
1 - 4	-8.502	3.827
1 - 5	-7.397	4.932
1 - 6	-11.112	1.217
2 - 3	-8.902	3.427
2 - 4	-2.747	9.582
2 - 5	-1.642	10.687
2 - 6	-5.357	6.972
3 - 4	-0.010	12.320
3 - 5	1.095	13.425
3 - 6	-2.620	9.710
4 - 5	-5.060	7.270
4 - 6	-8.775	3.555
5 - 6	-9.880	2.450



Compare all fertilizers to fertilizer 1 using Dunnett's method.

```
alpha <- 0.05
a <- 6
b <- 4
MSE <- anova_out$`Mean Sq`[3]
me <- 2.82 * sqrt(MSE) * sqrt(2/b) # value 2.82 from Dunnett's table
ntr_means <- aggregate(ntr, by = list(trt), mean)$x

CIs <- matrix(NA, a-1, 2)

k <- 1
for(i in 2:a){

  di <- ntr_means[i] - ntr_means[1]
  CIs[k,] <- c(di - me, di + me)
  k <- k + 1

}

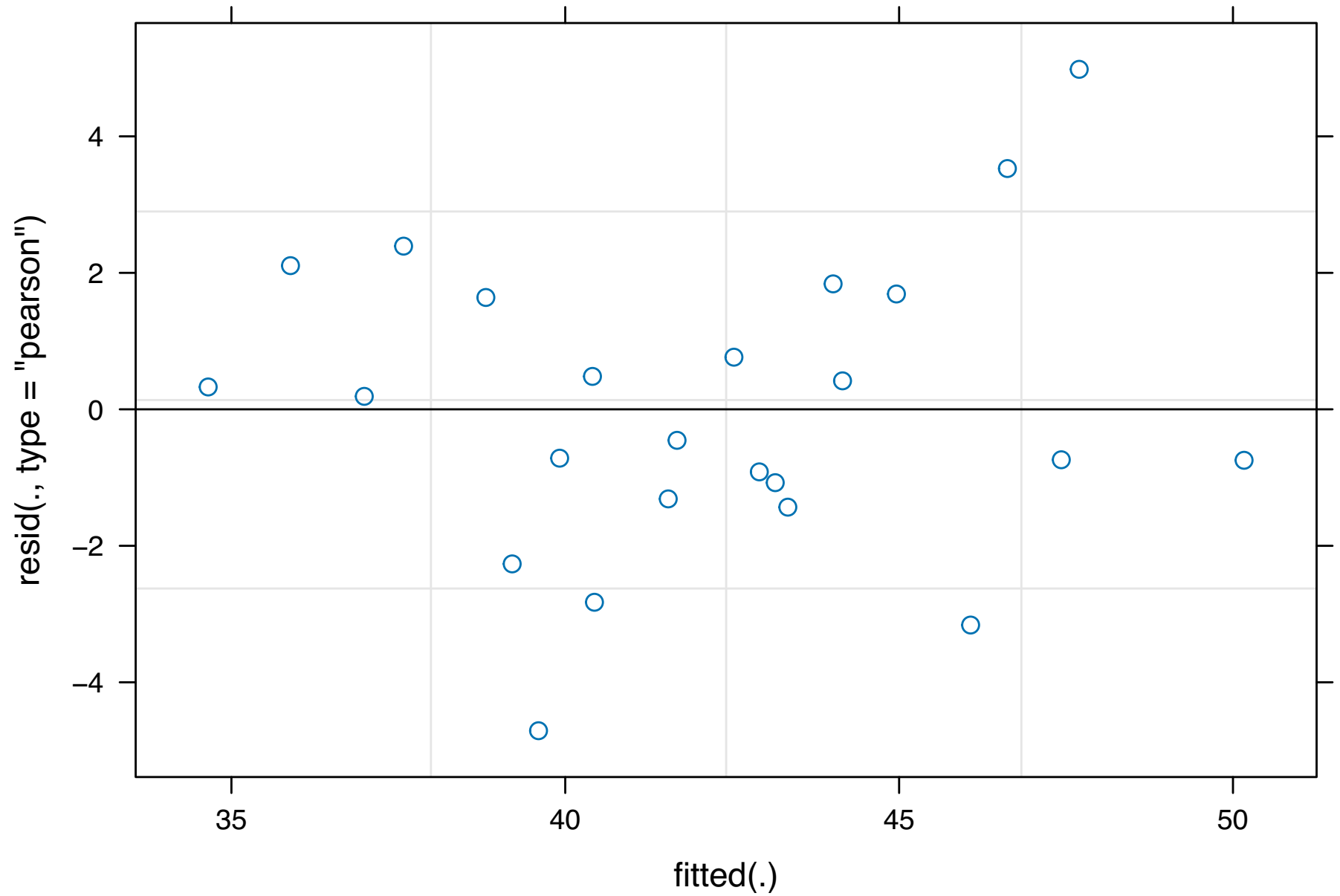
colnames(CIs) <- c("lower", "upper")
rownames(CIs) <- paste(2:a, "- 1")
```

`round(CIs,3)`

		lower	upper
2	- 1	0.404	11.106
3	- 1	3.142	13.843
4	- 1	-3.013	7.688
5	- 1	-4.118	6.583
6	- 1	-0.403	10.298

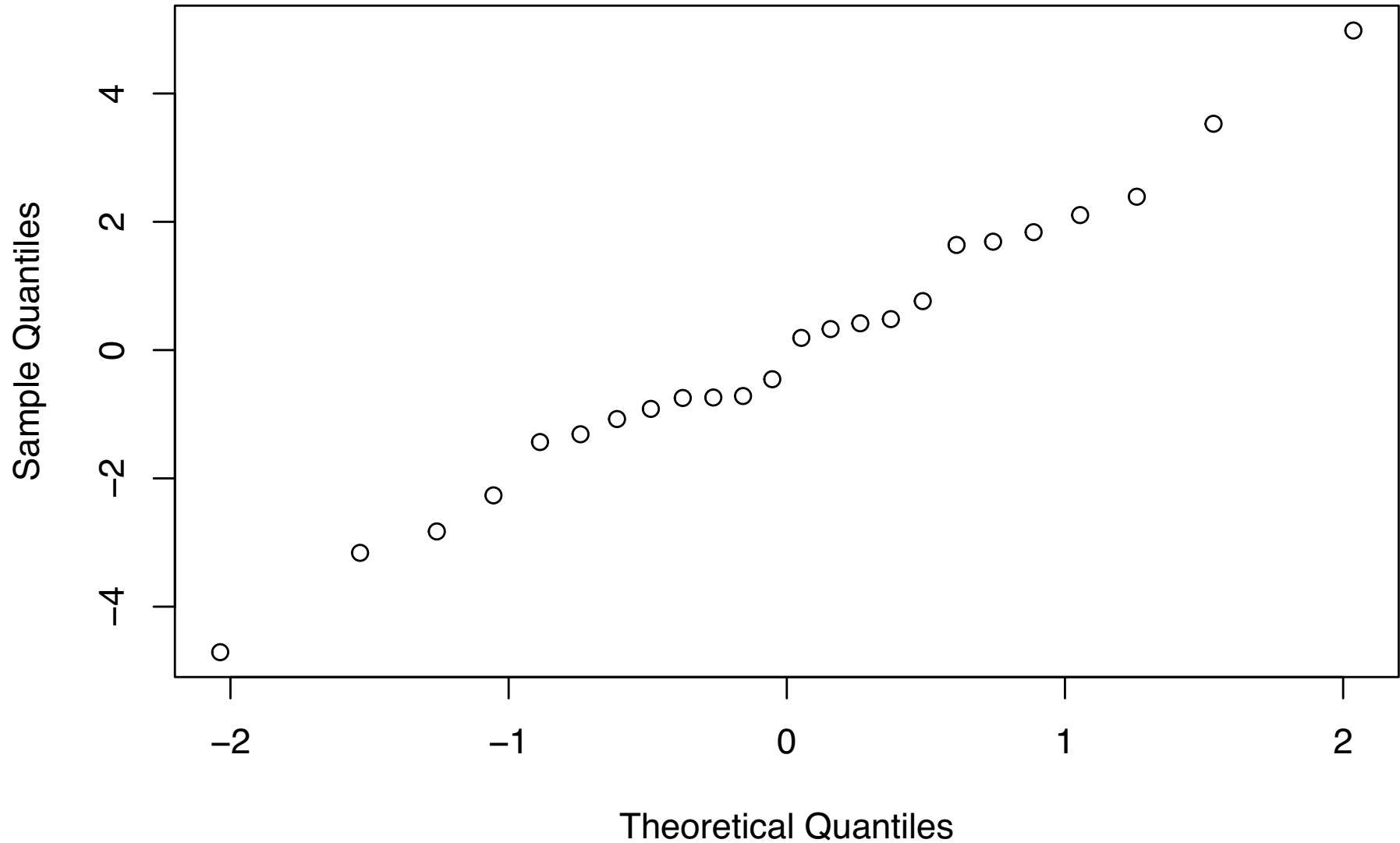


```
plot(lmer_out)
```



```
yhat <- predict(lmer_out)
ehat <- ntr - yhat
qqnorm(ehat)
```

## Normal Q-Q Plot



# Ignoring the blocks in the nitrogen data

If we ignore the blocks, the design looks like a one-way ANOVA:

Treatment	1	2	3	4	5	6
	34.98	40.89	42.07	37.18	37.99	34.89
	41.22	46.69	49.42	45.85	41.99	50.15
	36.94	46.65	52.68	40.23	37.61	44.57
	39.97	41.90	43.29	39.20	40.45	43.29

Suppose we fit  $Y_{ij} = \mu + \tau_i + \varepsilon_{ij}$ ,  $i = 1, 2, 3, 4, 5, 6$ ,  $j = 1, 2, 3, 4$ .

We lose power to detect a treatment effect!

```
anova(lm(ntr ~ trt))
```

↑ ignore the block?

$H_0: \mu_1 = \dots = \mu_a$

Analysis of Variance Table

Response: ntr

	Df	Sum Sq	Mean Sq	F value	Pr(>F)
trt	5	201.32	40.263	2.3761	0.08024 .
Residuals	18	305.01	16.945		

---

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

# Cabbage counts versus nitrogen, Kuehl (2000)

Heads of cabbage in subplots of two field plots.

$a = 5$

Nitrogen	Block 1		Block 2	
0	104	114	109	124
50	134	130	154	164
100	146	142	152	156
150	147	160	160	163
200	133	146	156	161

Source: Dr. W. B. Grew, Department of Plant Sciences, University of Arizona.

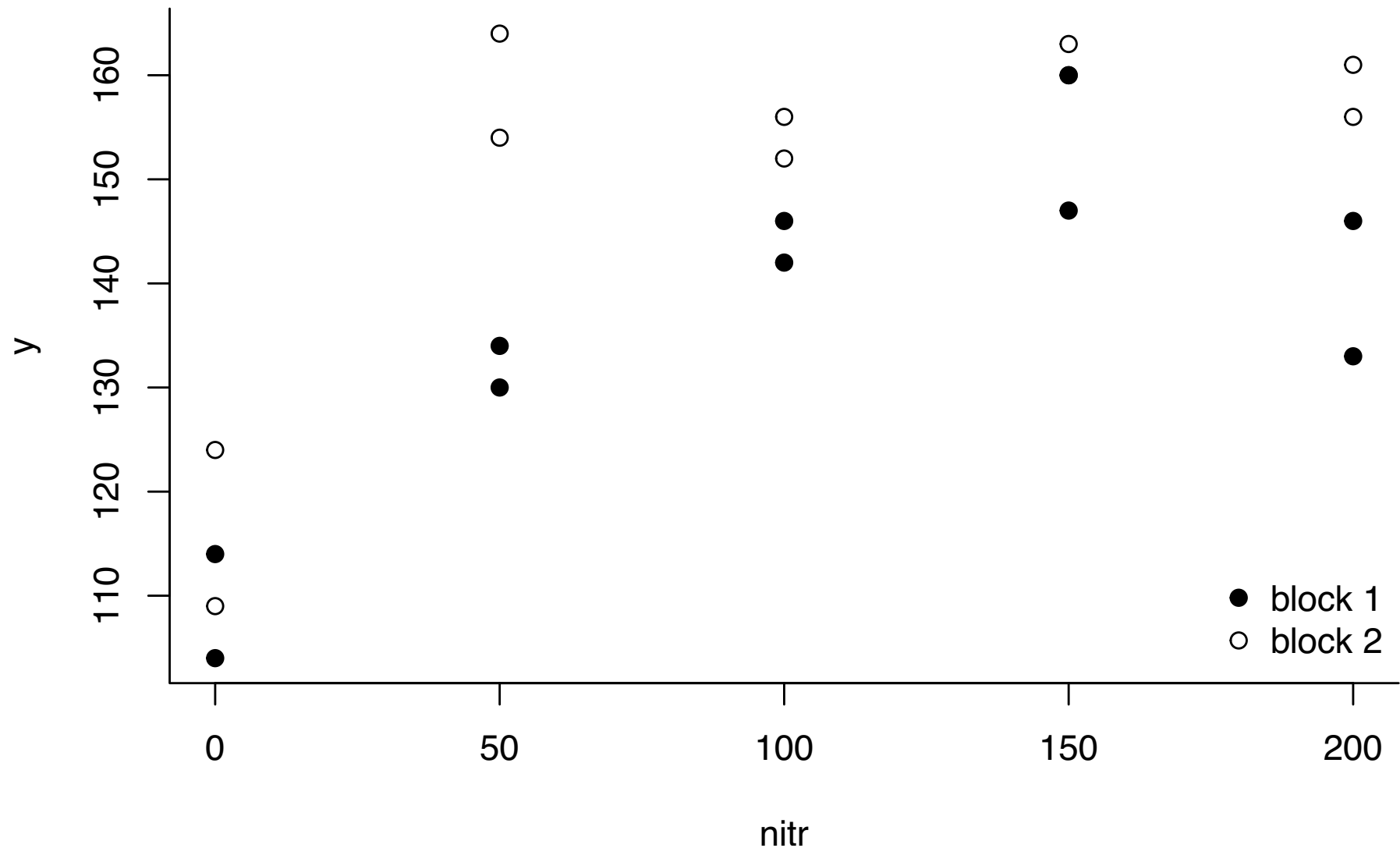
Subplots in each plot randomly assigned to nitrogen levels.

$$Y_{ijk} = \mu + \tau_i + \beta_j + (\tau\beta)_{ij} + \epsilon_{ijk} \quad \begin{array}{l} i = 1, \dots, 5 \\ j = 1, 2 \\ k = 1, 2 \end{array}$$

Block  $\times$  Nit interaction

```
y <- c(104,114,109,124,  
      134,130,154,164,  
      146,142,152,156,  
      147,160,160,163,  
      133,146,156,161)  
  
nitr <- as.factor(c(0,0,0,0,  
                  50,50,50,50,  
                  100,100,100,100,  
                  150,150,150,150,  
                  200,200,200,200))  
  
blk <- as.factor(c(1,1,2,2,  
                 1,1,2,2,  
                 1,1,2,2,  
                 1,1,2,2,  
                 1,1,2,2))
```

```
ylims <- range(y)
par(bty = "l")
stripchart(y~nitr, subset=blk==1, vertical=TRUE, pch=19, ylim=ylims,xlab="nitr")
stripchart(y~nitr, subset=blk==2, vertical=TRUE, pch=1, add=TRUE)
legend("bottomright",legend = c("block 1", "block 2"),pch = c(19, 1),bty="n")
```



# Treatment effects model for RCBD with replication

Assume

$$Y_{ijk} + \mu + \tau_i + B_j + (\tau B)_{ij} + \varepsilon_{ijk}$$

for  $i = 1, \dots, a$ ,  $j = 1, \dots, b$ , and  $k = 1, \dots, n_{ij}$ , where

- ▶  $Y_{ijk}$  is response of EU  $k$  in block  $j$  receiving treatment  $i$ .
- ▶ the  $\tau_i$  are the fixed effects of the treatment.
- ▶ the  $B_j$  are independent  $\text{Normal}(0, \sigma_B^2)$  random block effects.
- ▶ the  $(\tau B)_{ij}$  are indep.  $\text{Normal}(0, \sigma_{AB}^2)$  random interaction effects.
- ▶ the  $\varepsilon_{ijk}$  are independent  $\text{Normal}(0, \sigma_\varepsilon^2)$  error terms.
- ▶  $\mu$  is an overall or baseline mean.

Define the cell means as  $\mu_i = \mu + \tau_i$ ,  $i = 1, \dots, a$ .

Assume for now a balanced design:  $n_{ij} = n$  for all  $i, j$ .

# Sums of squares for RCBD with replication

Exactly like the 2-way ANOVA for two fixed factors

Sum of squares	Symbol	Formula
Total	$SS_{\text{Tot}}$	$\sum_{i=1}^a \sum_{j=1}^b \sum_{k=1}^n (Y_{ijk} - \bar{Y}_{...})^2$
Treatment	$SS_A$	$nb \sum_{i=1}^a (\bar{Y}_{i..} - \bar{Y}_{...})^2$
Block	$SS_B$	$na \sum_{j=1}^b (\bar{Y}_{.j.} - \bar{Y}_{...})^2$
Block $\times$ Treatment	$SS_{AB}$	$n \sum_{i=1}^a \sum_{j=1}^b (\bar{Y}_{ij.} - (\bar{Y}_{i..} + \bar{Y}_{.j.} - \bar{Y}_{...}))^2$
Error	$SS_{\text{Error}}$	$\sum_{i=1}^a \sum_{j=1}^b \sum_{k=1}^n (Y_{ijk} - \bar{Y}_{ij.})^2$

► Can make the decomposition  $SS_{\text{Tot}} = SS_A + SS_B + SS_{AB} + SS_{\text{Error}}$ .

# ANOVA table for RCBD with replication

Source	Df	SS	MS	F value
A	$a - 1$	$SS_A$	$MS_A$	$F_A = MS_A / MS_{AB}$
B	$b - 1$	$SS_B$	$MS_B$	$F_B = MS_B / MS_{AB}$
AB	$(a - 1)(b - 1)$	$SS_{AB}$	$MS_{AB}$	$F_{AB} = MS_{AB} / MS_{Error}$
Error	$ab(n - 1)$	$SS_{Error}$	$MS_{Error}$	
Total	$abn - 1$	$SS_{Tot}$		

1. Reject  $H_0: \mu_1 = \dots = \mu_a$  if  $F_A > F_{a-1, \underline{(a-1)(b-1)}, \alpha}$ .
2. Reject  $H_0: \sigma_B^2 = 0$  if  $F_B > F_{b-1, \underline{(a-1)(b-1)}, \alpha}$ .
3. Reject  $H_0: \sigma_{AB}^2 = 0$  if  $F_{AB} > F_{(a-1)(b-1), ab(n-1), \alpha}$ .

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

$\downarrow \sim N(0, \sigma_B^2)$ 
 $\downarrow \sim N(0, \sigma_{AB}^2)$ 
 $\downarrow \sim N(0, \sigma_\varepsilon^2)$

# Expected MS in RCBD with replication

$$MS_A = SSA / (a-1)$$

$$MS_B = SSB / (b-1)$$

$$MS_{AB} = \frac{SS_{AB}}{(a-1)(b-1)}$$

Source	Df	Expected mean square
A	$a - 1$	$nb\theta_A^2 + n\sigma_{AB}^2 + \sigma_\epsilon^2$
B	$b - 1$	<del><math>n\theta_B^2 + n\sigma_{AB}^2 + \sigma_\epsilon^2</math></del> ← Under $H_0$
AB	$(a - 1)(b - 1)$	$n\sigma_{AB}^2 + \sigma_\epsilon^2$
Error	$ab(n - 1)$	$\sigma_\epsilon^2$

In the above  $\theta_A^2 = (a - 1)^{-1} \sum_{i=1}^a (\mu_i - \bar{\mu})^2$ .

Check expected MS values under each  $H_0$  on previous slide.

$$\underline{H_0: \sigma_B^2 = 0} \quad \cdot \quad F_{\text{test}} = \frac{MS_B}{MS_{AB}}$$

# ANOVA table for cabbage count data

```
a <- nlevels(nitr)
b <- nlevels(blk)
n <- 2

y... <- predict(lm(y ~ 1))
yi.. <- predict(lm(y ~ nitr))
y.j. <- predict(lm(y ~ blk))
yij. <- predict(lm(y ~ nitr + blk + nitr:blk))

SST <- sum((y - y...)**2)
SSA <- sum((yi.. - y...)**2)
SSB <- sum((y.j. - y...)**2)
SSAB <- sum((yij. - (yi.. + y.j. - y...))**2)
SSE <- sum((y - yij.)**2)

MSA <- SSA / (a-1)
MSB <- SSB / (b-1)
MSAB <- SSAB / ((a-1)*(b-1))
MSE <- SSE / (a*b*(n-1))

FA <- MSA / MSAB
FB <- MSB / MSAB
FAB <- MSAB / MSE

pA <- 1 - pf(FA, a-1, (a-1)*(b-1))
pB <- 1 - pf(FB, b-1, (a-1)*(b-1))
pAB <- 1 - pf(FAB, (a-1)*(b-1), a*b*(n-1))
```

# ANOVA table for cabbage counts data

$a=2$   
 $b=2$

$q=5$   
 $b=2$   
 $n=2$

$H_0: \mu_1 = \dots = \mu_5$

Source	Df	SS	MS	F value	p value
A	4	4813.00	1203.25	16.7234	0.0092
B	1	1022.45	1022.45	14.2106	0.0196
AB	4	287.80	71.95	1.7030	0.2253
Error	10	422.50	42.25		
Total	19	6545.75			

$(a-1)(b-1)$

1

$(s-1)(2-1)$   
 $=4$

5-1

2-1

$abn - 1 = 5 \cdot 2 \cdot 2 - 1 = 19.$

$ab(n-1) = 5 \cdot 2 \cdot (2-1) = 10$

$H_0: \sigma_B^2 = 0$

$H_0: \sigma_{AB}^2 = 0$

# ANOVA table for cabbage count data

Note that `anova()` on `lm()` output does not give desired output:

```
anova(lm(y ~ nitr + blk + nitr:blk))
```

Analysis of Variance Table

Response: y

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
nitr	4	4813.0	1203.25	28.479	1.915e-05	***
blk	1	1022.5	1022.45	24.200	0.0006054	***
nitr:blk	4	287.8	71.95	1.703	0.2253096	
Residuals	10	422.5	42.25			

---

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

*Incorrect when  
blk is a random  
effect.*

# Method of moments for estimating $\sigma_B^2$ , $\sigma_{AB}^2$ , and $\sigma_\varepsilon^2$

Set the mean squares equal to their expectations and solve:

$$\begin{aligned} \text{MS}_B &\stackrel{\text{set}}{=} na\sigma_B^2 + n\sigma_{AB}^2 + \sigma_\varepsilon^2 \\ \text{MS}_{AB} &\stackrel{\text{set}}{=} n\sigma_{AB}^2 + \sigma_\varepsilon^2 \\ \text{MS}_{\text{Error}} &\stackrel{\text{set}}{=} \sigma_\varepsilon^2 \end{aligned}$$

Solve for  $\sigma_B^2$ ,  $\sigma_{AB}^2$ , and  $\sigma_\varepsilon^2$ :

$$\begin{aligned} \dot{\sigma}_B^2 &= \frac{1}{na} (\text{MS}_B - \text{MS}_{AB}) \\ \dot{\sigma}_{AB}^2 &= \frac{1}{n} (\text{MS}_{AB} - \text{MS}_{\text{Error}}) \\ \dot{\sigma}_\varepsilon^2 &= \text{MS}_{\text{Error}} \end{aligned}$$

Possible to obtain negative values for  $\dot{\sigma}_B^2$  and  $\dot{\sigma}_{AB}^2$ . Then use REML.

# MOM estimates of cabbage data variance components

```
sigma_B_sq <- (MSB - MSAB) / (n*a)
sigma_AB_sq <- (MSAB - MSE) / n
sigma_e_sq <- MSE

sigma_B <- sqrt(sigma_B_sq)
sigma_AB <- sqrt(sigma_AB_sq)
sigma_e <- sqrt(sigma_e_sq)
```

We obtain  $\hat{\sigma}_B = 9.749$ ,  $\hat{\sigma}_{AB} = 3.854$ , and  $\hat{\sigma}_\varepsilon = 6.500$ .

# REML estimates of cabbage data variance components

```
library(lmerTest) # first time run install.packages("lmerTest")
lmer_out <- lmer(y ~ nitr + (1|blk) + (1|nitr:blk))
lmer_out
```

Linear mixed model fit by REML ['lmerModLmerTest']

Formula:  $y \sim \text{nitr} + (1 \mid \text{blk}) + (1 \mid \text{nitr}:\text{blk})$

REML criterion at convergence: 110.9695

Random effects:

Groups	Name	Std.Dev.
nitr:blk	(Intercept)	3.854
blk	(Intercept)	9.749
Residual		6.500

Number of obs: 20, groups: nitr:blk, 10; blk, 2

Fixed Effects:

(Intercept)	nitr50	nitr100	nitr150	nitr200
112.75	32.75	36.25	44.75	36.25

# CIs (unadjusted) for diffs in means in RCBD with rep.

- ▶ A  $(1 - \alpha)100\%$  confidence interval for  $\mu_i - \mu_{i'}$  is

$$\bar{Y}_{i..} - \bar{Y}_{i'..} \pm t_{(a-1)(b-1), \alpha/2} \sqrt{MS_{AB}} \sqrt{\frac{2}{nb}}$$

for  $1 \leq i < i' \leq a$ .

- ▶ If  $(a - 1)(b - 1) = 1$ , this interval will be frightfully wide.
- ▶ In this case one should drop the interaction term from the model!

# CI for difference in means for cabbage count data

Compare the means of the nitrogen level 50 and 0 groups:

```
y50.. <- mean(y[nitr == "50"])
y0.. <- mean(y[nitr == "0"])
alpha <- 0.05
tval <- qt(1 - alpha/2, (a-1)*(b-1))
me <- tval * sqrt(MSAB) * sqrt(2/(n*b))
lo <- y50.. - y0.. - me
up <- y50.. - y0.. + me
c(lo, up)
```

```
[1] 16.09711 49.40289
```

# CI (unadjusted) for parameters in RCBD with replication

Use `confint()` on output of `lmer()`:

```
confint(lmer_out)
```

	2.5 %	97.5 %
.sig01	0.000000	6.253288
.sig02	2.715017	29.859582
.sigma	4.422824	9.062168
(Intercept)	95.101400	130.398323
nitr50	23.245458	42.254543
nitr100	26.745458	45.754543
nitr150	35.245458	54.254543
nitr200	26.745458	45.754543

These CIs are *not* adjusted to achieve a familywise coverage probability!

# RCBD with replication *without* an interaction term

If we omit the interaction term in the RCBD with replication, we have

$$Y_{ijk} + \mu + \tau_i + B_j + \varepsilon_{ijk}$$

for  $i = 1, \dots, a$ ,  $j = 1, \dots, b$ , and  $k = 1, \dots, n_{ij}$ , where

- ▶  $Y_{ijk}$  is response of EU  $k$  in block  $j$  receiving treatment  $i$ .
- ▶ the  $\tau_i$  are the fixed effects of the treatment.
- ▶ the  $B_j$  are independent  $\text{Normal}(0, \sigma_B^2)$  random block effects.
- ▶ the  $\varepsilon_{ijk}$  are independent  $\text{Normal}(0, \sigma_\varepsilon^2)$  error terms.
- ▶  $\mu$  is an overall or baseline mean.

Assume for now a balanced design:  $n_{ij} = n$  for all  $i, j$ .

Use this model if  $a = b = 2$ , under which  $(a - 1)(b - 1) \notin 1$ .

HW  
8

$$\mu_{ij} = \mu + \tau_i + \delta_j + (\tau\delta)_{ij}$$

$$i=1,2, \quad j=1,2.$$

$$\begin{aligned} \tau_1 &= 0 \\ \delta_1 &= 0 \end{aligned}$$

$\mu_{11}$	$\mu_{12}$	=	$\mu$	$\mu + \delta_2$
$\mu_{21}$	$\mu_{22}$		$\mu + \tau_2$	$\mu + \tau_2 + \delta_2 + (\tau\delta)_{22}$

# Sums of squares for RCBD with rep., no interaction

SS	Symbol	Formula
Total	$SS_{\text{Tot}}$	$\sum_{i=1}^a \sum_{j=1}^b \sum_{k=1}^n (Y_{ijk} - \bar{Y}_{...})^2$
Treat- ment	$SS_A$	$nb \sum_{i=1}^a (\bar{Y}_{i..} - \bar{Y}_{...})^2$
Block	$SS_B$	$na \sum_{j=1}^b (\bar{Y}_{.j.} - \bar{Y}_{...})^2$
Error	$SS_{\text{Error}}$	$\sum_{i=1}^a \sum_{j=1}^b \sum_{k=1}^n (Y_{ijk} - (\bar{Y}_{i..} + \bar{Y}_{.j.} - \bar{Y}_{...}))^2$

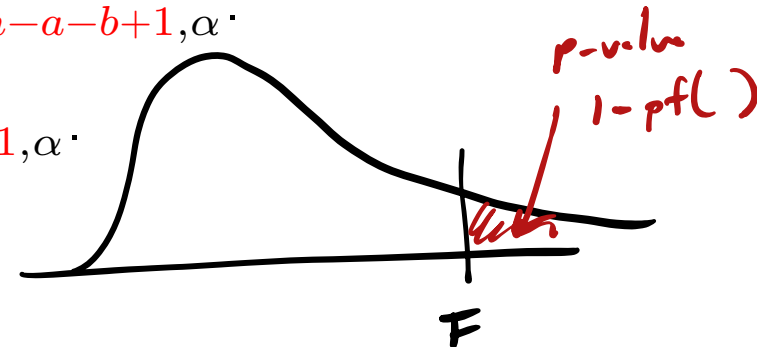
► Can make the decomposition  $SS_{\text{Tot}} = SS_A + SS_B + SS_{\text{Error}}$ .

# ANOVA table for RCBD with rep., omitting interaction

Source	Df	SS	MS	F value
A	$a - 1$	$SS_A$	$MS_A$	$F_A = MS_A / MS_{Error}$
B	$b - 1$	$SS_B$	$MS_B$	$F_B = MS_B / MS_{Error}$
Error	$abn - a - b + 1$	$SS_{Error}$	$MS_{Error}$	$\hat{\sigma}_e^2$
Total	$abn - 1$	$SS_{Tot}$		

1. Reject  $H_0: \mu_1 = \dots = \mu_a$  if  $F_A > F_{a-1, abn-a-b+1, \alpha}$ .

2. Reject  $H_0: \sigma_B^2 = 0$  if  $F_B > F_{b-1, abn-a-b+1, \alpha}$ .



Previously:

A	$a - 1$
B	$b - 1$
AB	$(a - 1)(b - 1)$
error	$ab(n - 1)$
total	$abn - 1$

combine these

$$(a - 1)(b - 1) + ab(n - 1) = ab - a - b + 1 + abn - ab = abn - a - b + 1$$

# Expected MS in RCBD with rep., omitting interaction

Source	Df	Expected mean square
A	$a - 1$	$nb\theta_A^2 + \sigma_\varepsilon^2$
B	$b - 1$	$na\sigma_B^2 + \sigma_\varepsilon^2$
Error	$abn - a - b + 1$	$\sigma_\varepsilon^2$

In the above  $\theta_A^2 = (a - 1)^{-1} \sum_{i=1}^a (\mu_i - \bar{\mu}_.)^2$ .

# ANOVA table for cabbage counts, omitting interaction

```
a <- nlevels(nitr)
b <- nlevels(blk)
n <- 2

y... <- predict(lm(y ~ 1))
yi.. <- predict(lm(y ~ nitr))
y.j. <- predict(lm(y ~ blk))

SST <- sum((y - y...)**2)
SSA <- sum((yi.. - y...)**2)
SSB <- sum((y.j. - y...)**2)
SSE <- sum((y - (yi.. + y.j. - y...))**2)

MSA <- SSA / (a-1)
MSB <- SSB / (b-1)
MSE <- SSE / (a*b*n - a - b + 1)

FA <- MSA / MSE
FB <- MSB / MSE

pA <- 1 - pf(FA, a-1, a*b*n - a - b + 1)
pB <- 1 - pf(FB, b-1, a*b*n - a - b + 1)
```

# ANOVA table for cabbage counts, omitting interaction

Source	Df	SS	MS	F value	p value
A	4	4813.00	1203.25	23.7160	0.0000
B	1	1022.45	1022.45	20.1525	0.0005
Error	14	710.30	50.74		
Total	19	6545.75			

$$H_0: \mu_1 = \dots = \mu_a$$

$$H_0: \sigma_B^2 = 0$$

# ANOVA table for cabbage counts, omitting interaction

With the interaction term omitting, `anova()` on `lm()` output is correct:

```
anova(lm(y ~ nitr + blk))
```

## Analysis of Variance Table

Response: y

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
nitr	4	4813.0	1203.25	23.716	4.13e-06	***
blk	1	1022.5	1022.45	20.152	0.0005097	***
Residuals	14	710.3	50.74			

---

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

# M.o.M. for $\sigma_B^2$ and $\sigma_\varepsilon^2$ when omitting interaction

Set the mean squares equal to their expectations and solve:

$$MS_B \stackrel{\text{set}}{=} na\sigma_B^2 + \sigma_\varepsilon^2$$

$$MS_{\text{Error}} \stackrel{\text{set}}{=} \sigma_\varepsilon^2$$

Solve for  $\sigma_B^2$  and  $\sigma_\varepsilon^2$ :

$$\dot{\sigma}_B^2 = \frac{1}{na} (MS_B - MS_{\text{Error}})$$

$$\dot{\sigma}_\varepsilon^2 = MS_{\text{Error}}$$

Possible to obtain a negative value for  $\dot{\sigma}_B^2$ . In this case use REML.

# M.o.M.s for cabbage data when omitting interaction

```
sigma_B_sq <- (MSB - MSE) / (n*a)
sigma_e_sq <- MSE

sigma_B <- sqrt(sigma_B_sq)
sigma_e <- sqrt(sigma_e_sq)
```

We obtain  $\hat{\sigma}_B = 9.858$  and  $\hat{\sigma}_\varepsilon = 7.123$ .

# REML for cabbage data when omitting interaction

```
library(lmerTest) # first time run install.packages("lmerTest")
lmer_out <- lmer(y ~ nitr + (1|blk))
lmer_out
```

Linear mixed model fit by REML ['lmerModLmerTest']

Formula: y ~ nitr + (1 | blk)

REML criterion at convergence: 111.4024

Random effects:

Groups	Name	Std.Dev.	
blk	(Intercept)	9.858	$\hat{\sigma}_B$
	Residual	7.123	$\hat{\sigma}_\epsilon$

Number of obs: 20, groups: blk, 2

Fixed Effects:

(Intercept)	nitr50	nitr100	nitr150	nitr200
112.75	32.75	36.25	44.75	36.25

# CIs (unadjusted) for diffs in means in RCBD with rep. without interaction

- ▶ A  $(1 - \alpha)100\%$  confidence interval for  $\underline{\mu}_i - \underline{\mu}_{i'}$  is

$$\bar{Y}_{i..} - \bar{Y}_{i'..} \pm t_{abn-a-b+1, \alpha/2} \sqrt{MS_{\text{Error}}} \sqrt{\frac{2}{nb}}$$

for  $1 \leq i < i' \leq a$ .

# CI for difference in means for cabbage count data when omitting interaction

Compare the means of the nitrogen level 50 and 0 groups:

```
y50.. <- mean(y[nitr == "50"])
y0.. <- mean(y[nitr == "0"])
alpha <- 0.05
tval <- qt(1 - alpha/2, a*b*n - a - b + 1)
me <- tval * sqrt(MSE) * sqrt(2/(n*b))
lo <- y50.. - y0.. - me
up <- y50.. - y0.. + me
c(lo, up)
```

$t_{abn - a - b + 1, \alpha/2}$

```
[1] 21.94746 43.55254
```

# CI for difference in means for cabbage count data when omitting interaction

Compare above to `confint()` on output from `lmer()`.

```
confint(lmer_out)
```

	2.5 %	97.5 %
<i>nitr 0</i> .sig01	2.715987	29.862871
.sigma	4.678640	9.062171
(Intercept)	95.101978	130.398015
nitr50	23.558217	41.941783
nitr100	27.058217	45.441783
nitr150	35.558217	53.941783
nitr200	27.058217	45.441783

*CI for  $\sigma_B^2$*   
*CI for  $\sigma_\epsilon^2$*   
*CI for  $\mu$ , mean under nitro 0*  
*CI for  $\tau_{50} = \text{mean nitro 50} - \text{mean nitro 0}$*

$$Y_{ijk} = \underbrace{\mu + \tau_i}_{\mu_i} + \beta_j + \epsilon_{ijk}$$

$\epsilon_{ijk} \sim N(0, \sigma_\epsilon^2)$

# Skin response data, Mohr, Wilson, and Freund (2021)

Galvanic skin responses of five subjects under shock and noise stimuli.

		SUBJECT				
Noise	Shock	1	2	3	4	5
40	0.25	3	7	9	4	1
40	0.50	5	11	13	8	3
40	0.75	9	12	14	11	5
40	1.00	6	11	12	7	4
80	0.25	5	10	10	6	3
80	0.50	6	12	15	9	5
80	0.75	18	18	15	13	9
80	1.00	7	15	14	9	7

```
skin <- data.frame(resp = c(3,7,9,4,1,5,11,13,8,3,9,12,14,11,5,6,11,12,7,4,5,  
                            10,10,6,3,6,12,15,9,5,18,18,15,13,9,7,15,14,9,7),  
                  noise = as.factor(c(rep(40,20),rep(80,20))),  
                  shock = as.factor(rep(c(rep(.25,5),rep(.5,5),  
                                          rep(.75,5),rep(1,5)),2)),  
                  subj = as.factor(rep(1:5,8)))
```

```
head(skin,n=20)
```

	resp	noise	shock	subj
1	3	40	0.25	1
2	7	40	0.25	2
3	9	40	0.25	3
4	4	40	0.25	4
5	1	40	0.25	5
6	5	40	0.5	1
7	11	40	0.5	2
8	13	40	0.5	3
9	8	40	0.5	4
10	3	40	0.5	5
11	9	40	0.75	1
12	12	40	0.75	2
13	14	40	0.75	3
14	11	40	0.75	4
15	5	40	0.75	5
16	6	40	1	1
17	11	40	1	2
18	12	40	1	3
19	7	40	1	4
20	4	40	1	5

# Treatment effects model for two-way factorial RCBD

Assume

$$Y_{ijk} = \mu + \tau_i + \gamma_j + (\tau\gamma)_{ij} + C_k + \varepsilon_{ijk},$$

for  $i = 1, \dots, a$ ,  $j = 1, \dots, b$ , and  $k = 1, \dots, c$ , where

- ▶  $Y_{ijk}$  is the response in block  $k$  under treatment combination  $i \times j$ .
- ▶  $\mu$  is an overall or baseline mean.
- ▶ the  $\tau_i$  are fixed effects for factor A.
- ▶ the  $\gamma_j$  are fixed effects for factor B.
- ▶ the  $(\tau\gamma)_{ij}$  are effects for the A×B interaction.
- ▶ the  $C_k$  are independent  $\text{Normal}(0, \sigma_C^2)$  block effects.
- ▶ the  $\varepsilon_{ijk}$  are independent  $\text{Normal}(0, \sigma_\varepsilon^2)$  error terms.

Define the cell means as

$$\mu_{ij} = \mu + \tau_i + \gamma_j + (\tau\gamma)_{ij}, \quad i = 1, \dots, a, \quad j = 1, \dots, b.$$

# Sums of squares for the two-way factorial RCBD

Sum of squares	Symbol	Formula
Total	$SS_{\text{Tot}}$	$\sum_{i=1}^a \sum_{j=1}^b \sum_{k=1}^c (Y_{ijk} - \bar{Y}_{...})^2$
A	$SS_A$	$bc \sum_{i=1}^a (\bar{Y}_{i..} - \bar{Y}_{...})^2$
B	$SS_B$	$ac \sum_{j=1}^b (\bar{Y}_{.j.} - \bar{Y}_{...})^2$
AB	$SS_{AB}$	$c \sum_{i=1}^a \sum_{j=1}^b (Y_{ij.} - (\bar{Y}_{i..} + \bar{Y}_{.j.} - \bar{Y}_{...}))^2$
C	$SS_C$	$ab \sum_{k=1}^c (\bar{Y}_{..k} - \bar{Y}_{...})^2$
Error	$SS_{\text{Error}}$	$SS_{\text{Tot}} - (SS_A + SS_B + SS_{AB} + SS_C)$

- ▶ Then we have  $SS_{\text{Tot}} = SS_A + SS_B + SS_{AB} + SS_C + SS_{\text{Error}}$ .
- ▶ The error  $SS_{\text{Error}}$  is really the interaction sum of squares  $SS_{\text{Trt} \times C}$ .
- ▶ Again, without replication, we cannot estimate this interaction.
- ▶ So the interaction serves as the error term.

# ANOVA table for two-way factorial RCBD

Source	Df	SS	MS	F value
A	$a - 1$	$SS_A$	$MS_A$	$F_A = MS_A / MS_{\text{Error}}$
B	$b - 1$	$SS_B$	$MS_B$	$F_B = MS_B / MS_{\text{Error}}$
AB	$(a - 1)(b - 1)$	$SS_{AB}$	$MS_{AB}$	$F_{AB} = MS_{AB} / MS_{\text{Error}}$
C	$c - 1$	$SS_C$	$MS_C$	$F_C = MS_C / MS_{\text{Error}}$
Error	$(ab - 1)(c - 1)$	$SS_{\text{Error}}$	$MS_{\text{Error}}$	
Total	$abc - 1$	$SS_{\text{Tot}}$		

1. Reject  $H_0: \mu_{1.} = \dots = \mu_{a.}$  if  $F_A > F_{a-1, (ab-1)(c-1), \alpha}$ .
2. Reject  $H_0: \mu_{.1} = \dots = \mu_{.b}$  if  $F_B > F_{b-1, (ab-1)(c-1), \alpha}$ .
3. R.  $H_0: \mu_{ij} = \bar{\mu}_{i.} + \bar{\mu}_{.j} - \bar{\mu}_{..} \quad \forall ij$  if  $F_{AB} > F_{(a-1)(b-1), (ab-1)(c-1), \alpha}$ .
4. Reject  $H_0: \sigma_C^2 = 0$  if  $F_C > F_{c-1, (ab-1)(c-1), \alpha}$ .

# Skin response data (cont)

```
lm_out <- lm(resp ~ noise + shock + noise:shock + subj, data = skin)
anova(lm_out)
```

## Analysis of Variance Table

Response: resp

	Df	Sum Sq	Mean Sq	F value	Pr(>F)	
noise	1	65.02	65.025	28.3819	1.134e-05	***
shock	3	219.27	73.092	31.9028	3.564e-09	***
subj	4	361.85	90.463	39.4848	3.975e-11	***
noise:shock	3	12.67	4.225	1.8441	0.1621	
Residuals	28	64.15	2.291			

---

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

```

y <- skin$resp
y... <- predict(lm(resp ~ 1,data = skin))
yi.. <- predict(lm(resp ~ noise,data=skin))
y.j. <- predict(lm(resp ~ shock,data=skin))
y..k <- predict(lm(resp ~ subj,data=skin))
yij. <- predict(lm(resp ~ noise + shock + noise:shock,data=skin))

SSA <- sum((yi.. - y...) ^2)
SSB <- sum((y.j. - y...) ^2)
SSC <- sum((y..k - y...) ^2)
SSAB <- sum((yij. - (yi.. + y.j. - y...)) ^2)
SST <- sum((y - y...) ^2)
SSE <- SST - (SSA + SSB + SSC + SSAB)

a <- 2
b <- 4
c <- 5

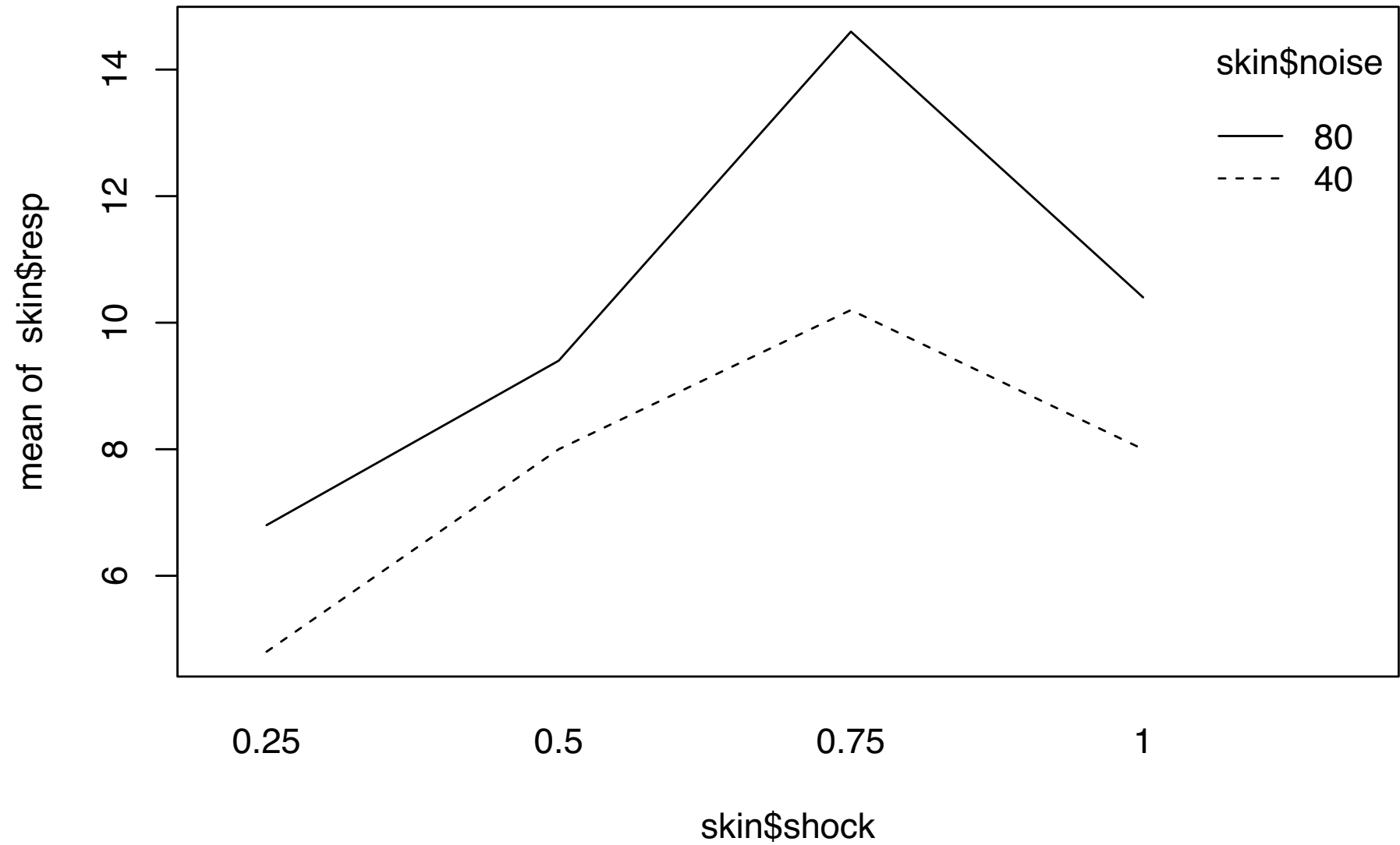
MSA <- SSA/(a-1)
MSB <- SSB/(b-1)
MSC <- SSC/(c-1)
MSAB <- SSAB/((a-1)*(b-1))
MSE <- SSE/((a*b-1)*(c-1))

FA <- MSA / MSE
FB <- MSB / MSE
FAB <- MSAB / MSE
FC <- MSC / MSE

pA <- 1 - pf(FA,a-1,(a*b-1)*(c-1))
pB <- 1 - pf(FB,b-1,(a*b-1)*(c-1))
pAB <- 1 - pf(FAB,(a-1)*(b-1),(a*b-1)*(c-1))
pC <- 1 - pf(FC,c-1,(a*b-1)*(c-1))

```

```
interaction.plot(skin$shock,skin$noise,skin$resp)
```



# Expected mean squares in factorial RCBD

Source	Df	Expected mean square
A	$a - 1$	$bc\theta_A^2 + \sigma_\varepsilon^2$
B	$b - 1$	$ac\theta_B^2 + \sigma_\varepsilon^2$
AB	$(a - 1)(b - 1)$	$c\theta_{AB}^2 + \sigma_\varepsilon^2$
C	$c - 1$	$ab\sigma_C^2 + \sigma_\varepsilon^2$
Error	$(ab - 1)(c - 1)$	$\sigma_\varepsilon^2$

In the above

$$\blacktriangleright \theta_A^2 = (a - 1)^{-1} \sum_{i=1}^a (\bar{\mu}_{i.} - \bar{\mu}_{..})^2$$

$$\blacktriangleright \theta_B^2 = (b - 1)^{-1} \sum_{j=1}^b (\bar{\mu}_{.j} - \bar{\mu}_{..})^2$$

$$\blacktriangleright \theta_{AB}^2 = [(a - 1)(b - 1)]^{-1} \sum_{i=1}^a \sum_{j=1}^b (\mu_{ij} - (\bar{\mu}_{i.} + \bar{\mu}_{.j} - \bar{\mu}_{..}))^2$$

# MoMs for variance components in two-way factorial RCBD

- ▶ Equating  $MS_C$  and  $MS_{\text{Error}}$  with their expectations gives

$$\dot{\sigma}_C^2 = \frac{MS_C - MS_{\text{Error}}}{ab} \quad \text{and} \quad \dot{\sigma}_\varepsilon^2 = MS_{\text{Error}}.$$

- ▶ May obtain  $\dot{\sigma}_C^2 < 0$ , so one should use REML estimation.

# Skin response data (cont)

Obtain REML estimators of  $\sigma_C^2$  and  $\sigma_\varepsilon^2$  on the skin response data.

```
lmer_out <- lmer(resp ~ noise + shock + noise:shock + (1|subj), data = skin)
lmer_out
```

Linear mixed model fit by REML ['lmerModLmerTest']

Formula: resp ~ noise + shock + noise:shock + (1 | subj)

Data: skin

REML criterion at convergence: 144.9199

Random effects:

Groups	Name	Std.Dev.
subj	(Intercept)	3.320
Residual		1.514

Number of obs: 40, groups: subj, 5

Fixed Effects:

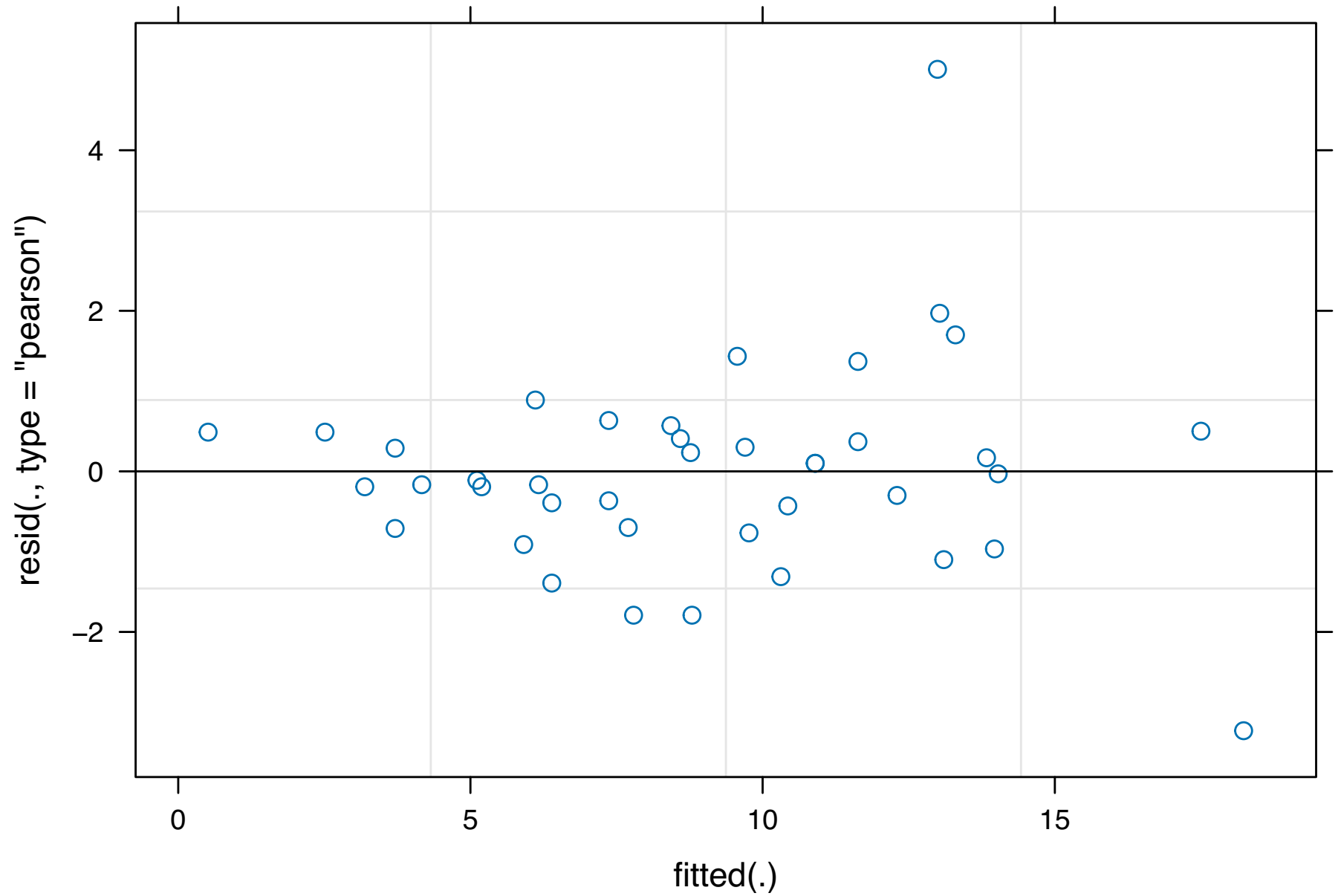
(Intercept)	noise80	shock0.5	shock0.75
4.8	2.0	3.2	5.4
shock1	noise80:shock0.5	noise80:shock0.75	noise80:shock1
3.2	-0.6	2.4	0.4

Obtain MoMs estimators for  $\sigma_C^2$  and  $\sigma_\varepsilon^2$  on the skin response data.

```
sg_C <- sqrt((MSC - MSE)/(a*b))  
sg_e <- sqrt(MSE)
```

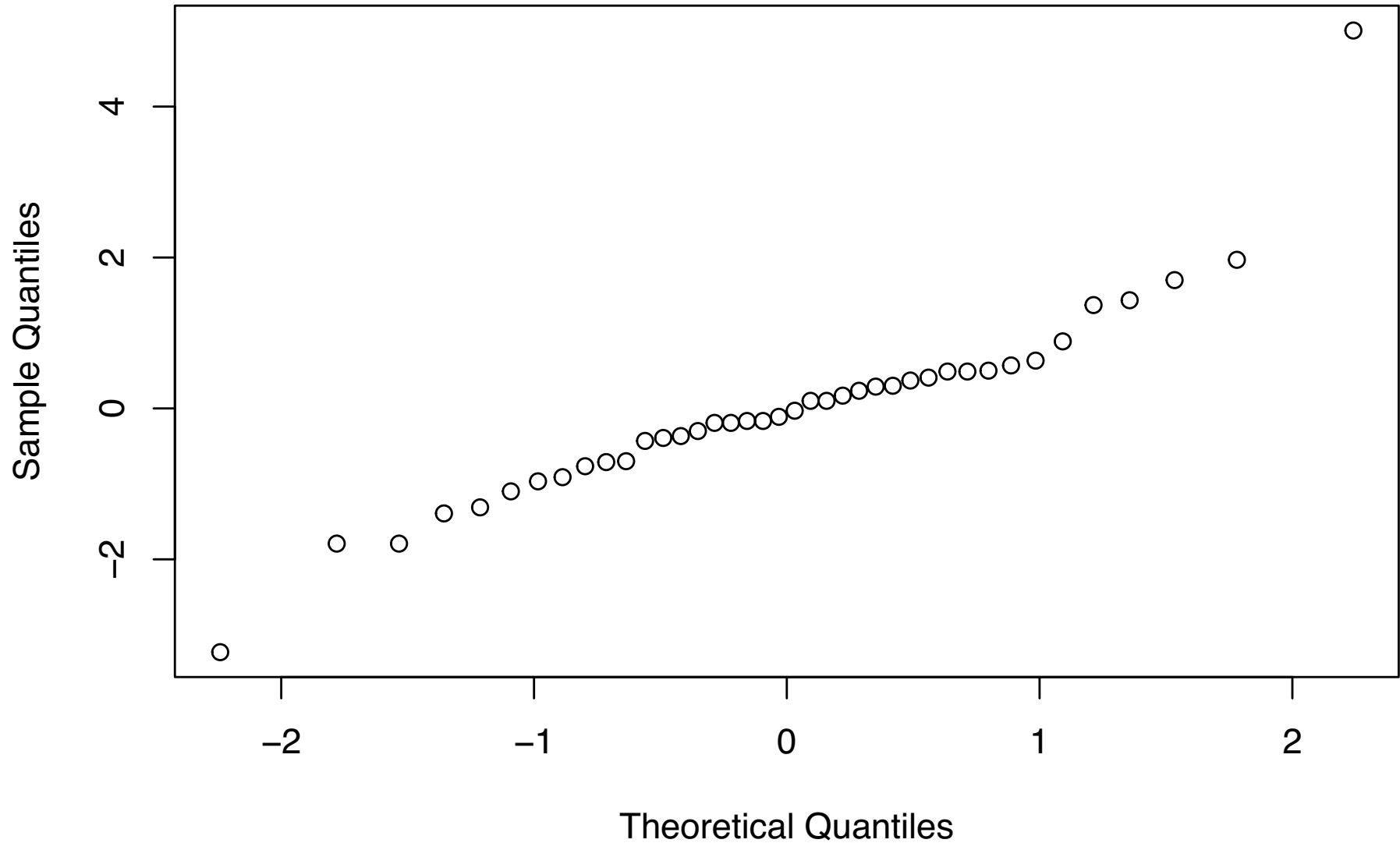
We have  $\hat{\sigma}_C = 3.320$  and  $\hat{\sigma}_\varepsilon = 1.514$ .

```
plot(lmer_out)
```



```
yhat <- predict(lmer_out)
ehat <- skin$resp - yhat
qqnorm(ehat)
```

## Normal Q-Q Plot



# References

- Kuehl, R. O. 2000. *Design of Experiments: Statistical Principles of Research Design and Analysis*. Duxbury/Thomson Learning.
- Mohr, Donna L, William J Wilson, and Rudolf J Freund. 2021. *Statistical Methods*. Academic Press.