# mixed\_models\_guide

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8/3/22

## Table of contents

Pr	reface	3
1	Introduction	4
2	Basic Models	5
3	Generalized Linear Mixed Models 3.1 Hurdle model	<b>6</b>
4	Special Conditions 4.1 Split plot with repeated measures	<b>17</b> 17
5	Summary	23
Re	eferences	24

## **Preface**

This is a defaulf file accompanying a quarto book.

To learn more about Quarto books visit https://quarto.org/docs/books.

### 1 Introduction

This guide is focused on frequentist implementations of mixed models in R. If someone wants to write a Bayesian guide, please go for it! I'm not experience sufficiently in Bayesian to do this.

Each section contains the minimum to run a model, with more detail found at the later chapters. Unless I decide it makes more sense to include early materials.

A Tidymodels framework is used whenever possible because that is a promising avenue for making the syntax easier to write across packages.

## 2 Basic Models

### 3 Generalized Linear Mixed Models

```
The following objects are masked from 'package:stats':

filter, lag

The following objects are masked from 'package:base':

intersect, setdiff, setequal, union

library(ggplot2)
library(glmmTMB)
library(DHARMa)

This is DHARMa 0.4.6. For overview type '?DHARMa'. For recent changes, type news(package = 'Package')
```

### 3.1 Hurdle model

library(dplyr)

Attaching package: 'dplyr'

```
insect_exp <- read.csv("data/insect_count_data_glmm.csv")
plot: a unique number referring to each experimental unit
treatment: pesticidal treatment (6 different products)
row: plot position for row
col: plot positions for column or range
block: the blocking unit</pre>
```

insect\_counts: response variable

sampling\_date: dates when each experimental unit were evaluated for insect counts

```
head(insect_exp)
```

	plot	treatment	row	column	block	<pre>insect_counts</pre>	sampling_date
1	101	2	1	1	1	4	6/17/88
2	102	5	1	2	1	1	6/17/88
3	103	1	1	3	1	0	6/17/88
4	104	6	1	4	1	0	6/17/88
5	201	3	2	1	1	0	6/17/88
6	202	4	2	2	1	0	6/17/88

Two new variables created:

treatment: original variable treatment converted to a factor

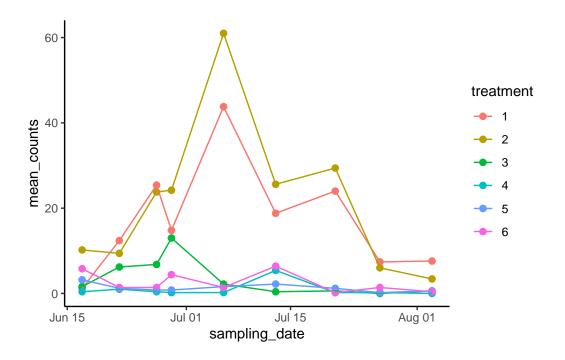
block: original variable block converted to a factor

Date: factor version of sampling\_date

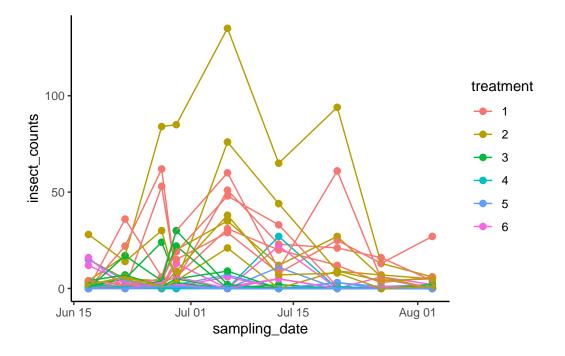
Visualise data

```
library(ggplot2)

insect_exp %>% group_by(sampling_date, treatment) %>%
   summarise(mean_counts = mean(insect_counts)) %>%
   ggplot(., aes(x = sampling_date, y = mean_counts, color = treatment)) +
   geom_point(size = 2) +
   geom_line() +
   theme_classic()
```



```
ggplot(insect_exp, aes(x = sampling_date, y = insect_counts, color = treatment, group = pl
  geom_point(size = 2) +
  geom_line() +
  theme_classic()
```



Model statement {[[[[ FIX THIS - it's still written for alfalfa ]]]]}

$$y_{ijk} = \mu + \alpha_i + \beta_j + \gamma_k + a_l + b_m + c_n + \epsilon$$

where

 $\mu=$  overall mean/intercept  $\alpha_i=$  effect of the  $i^{th}$  pesticide treatment  $\beta_j=$  effect of the  $j^{th}$  block  $\gamma_k=$  effect of the  $k^{th}$  sampling date

To make things easier, the interactions between the fixed effects are not shown.

```
library(glmmTMB)

m1 = glmmTMB(
  insect_counts ~ treatment + Date + ar1(Date + O|plot) + (1|block),
  ziformula = ~ treatment,
  data = insect_exp, na.action = na.exclude,
  family = nbinom2)
```

special correlation structure for correlated error terms ar1() (autoregressive 1).

There are several other specialized covariance structures implemented by glmmTMB. In general, repeated measures syntax follow this convention: (time + 0 | grouping).

We can test other distributions

```
m2 <- update(m1, family = poisson)</pre>
```

Warning in fitTMB(TMBStruc): Model convergence problem; non-positive-definite Hessian matrix. See vignette('troubleshooting')

Warning in fitTMB(TMBStruc): Model convergence problem; false convergence (8). See vignette('troubleshooting')

```
m3 <- update(m1, family = nbinom1)</pre>
```

Warning in (function (start, objective, gradient = NULL, hessian = NULL, : NA/NaN function evaluation

Warning in fitTMB(TMBStruc): Model convergence problem; non-positive-definite Hessian matrix. See vignette('troubleshooting')

Warning in fitTMB(TMBStruc): Model convergence problem; false convergence (8). See vignette('troubleshooting')

Fitting glmm is hard. Basic guidance on model fitting: https://glmmtmb.github.io/glmmTMB/articles/troubles

```
diagnose(m2)
```

Unusually large Z-statistics (|x|>5):

```
treatment3 treatment4 treatment5 Date1988-07-06 Date1988-07-13
-6.616584 -15.016232 -10.313314 9.371860 6.381998
Date1988-08-03 zi~(Intercept) zi~treatment4
-5.415075 -22.156803 59.213969
```

Large Z-statistics (estimate/std err) suggest a \*possible\* failure of the Wald approximation - often also associated with parameters that are at or near the edge of their range (e.g. random-effects standard deviations approaching 0). (Alternately, they may simply represent very well-estimated parameters; intercepts of non-centered models may fall in this category.) While the Wald p-values and standard errors listed in summary() may be unreliable, profile confidence intervals (see ?confint.glmmTMB) and likelihood ratio test p-values derived by

comparing models (e.g. ?drop1) are probably still OK. (Note that the LRT is conservative when the null value is on the boundary, e.g. a variance or zero-inflation value of O (Self and Liang 1987; Stram and Lee 1994; Goldman and Whelan 2000); in simple cases these p-values are approximately twice as large as they should be.)

Non-positive definite (NPD) Hessian

The Hessian matrix represents the curvature of the log-likelihood surface at the maximum likelihood estimate (MLE) of the parameters (its inverse is the estimate of the parameter covariance matrix). A non-positive-definite Hessian means that the likelihood surface is approximately flat (or upward-curving) at the MLE, which means the model is overfitted or poorly posed in some way. NPD Hessians are often associated with extreme parameter estimates.

parameters with non-finite standard deviations: (Intercept), Date1988-06-22, zi~treatment3, zi~treatment6, theta\_Date+0|plot.2, theta\_1|block.1

recomputing Hessian via Richardson extrapolation. If this is too slow, consider setting check

Hessian has complex eigenvalues

We would have used the smallest eigenvalues of the Hessian to determine which components were bad but instead we got complex eigenvalues. (Not really sure what to do with this ...)

```
diagnose(m3)
```

Unusually large coefficients (|x|>10):

```
d~(Intercept)
-26.35181
```

Large negative coefficients in zi (log-odds of zero-inflation), dispersion, or random effects (log-standard deviations) suggest

unnecessary components (converging to zero on the constrained scale); large negative and/or positive components in binomial or Poisson conditional parameters suggest (quasi-)complete separation. Large values of nbinom2 dispersion suggest that you should use a Poisson model instead.

Unusually large Z-statistics (|x|>5):

treatment5 -6.281128

Large Z-statistics (estimate/std err) suggest a \*possible\* failure of the Wald approximation - often also associated with parameters that are at or near the edge of their range (e.g. random-effects standard deviations approaching 0). (Alternately, they may simply represent very well-estimated parameters; intercepts of non-centered models may fall in this category.) While the Wald p-values and standard errors listed in summary() may be unreliable, profile confidence intervals (see ?confint.glmmTMB) and likelihood ratio test p-values derived by comparing models (e.g. ?drop1) are probably still OK. (Note that the LRT is conservative when the null value is on the boundary, e.g. a variance or zero-inflation value of 0 (Self and Liang 1987; Stram and Lee 1994; Goldman and Whelan 2000); in simple cases these p-values are approximately twice as large as they should be.)

Non-positive definite (NPD) Hessian

The Hessian matrix represents the curvature of the log-likelihood surface at the maximum likelihood estimate (MLE) of the parameters (its inverse is the estimate of the parameter covariance matrix). A non-positive-definite Hessian means that the likelihood surface is approximately flat (or upward-curving) at the MLE, which means the model is overfitted or poorly posed in some way. NPD Hessians are often associated with extreme parameter estimates.

parameters with non-finite standard deviations:  $d \sim (Intercept)$ 

recomputing Hessian via Richardson extrapolation. If this is too slow, consider setting check

Hessian has complex eigenvalues

We would have used the smallest eigenvalues of the Hessian to determine which components were bad but instead we got complex eigenvalues. (Not really sure what to do with this ...)

#### Summary info

m1

#### Formula:

insect\_counts ~ treatment + Date + ar1(Date + 0 | plot) + (1 | block)

Zero inflation: ~treatment

Data: insect\_exp

AIC BIC logLik df.resid 1298.7328 1385.0949 -625.3664 246

Random-effects (co)variances:

#### Conditional model:

Groups Name Std.Dev. Corr
plot Date1988-06-17 0.7748 0.49 (ar1)

block (Intercept) 0.3333

Number of obs: 270 / Conditional model: plot, 30; block, 5

Dispersion parameter for nbinom2 family (): 1.76

#### Fixed Effects:

#### Conditional model:

(Intercept)	treatment2	treatment3	treatment4	treatment5
2.39231	-0.04978	-1.53159	-2.75395	-2.50652
treatment6	Date1988-06-22	Date1988-06-27	Date1988-06-29	Date1988-07-06
-1.48975	0.24054	0.26618	0.62692	1.17067
Date1988-07-13	Date1988-07-21	Date1988-07-27	Date1988-08-03	
0.83442	0.19962	-0.96749	-1.11938	

#### Zero-inflation model:

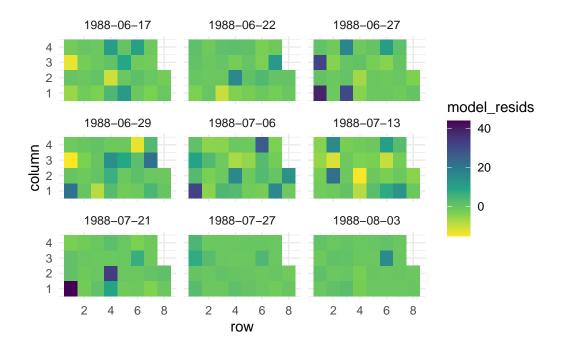
(Intercept)	treatment2	treatment3	treatment4	treatment5	treatment6
-2.608	-1.200	1.568	2.607	1.542	2.134

#### Diagnostics

Look at residuals over space

```
insect_exp$model_resids <- residuals(m1)

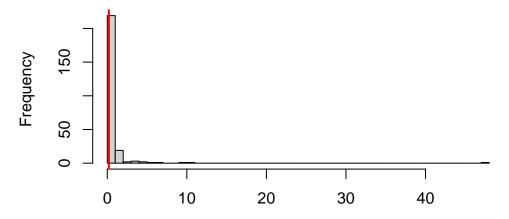
ggplot(insect_exp, aes(x = row, y = column, fill = model_resids)) +
    geom_tile() +
    facet_wrap(facets = vars(Date), nrow = 3, ncol = 3) +
    scale_fill_viridis_c(direction = -1) +
    theme_minimal()</pre>
```



#### use **DHARMa** to conduct residual tests

```
simulated_resids <- simulateResiduals(m1)
testDispersion(simulated_resids)</pre>
```

## DHARMa nonparametric dispersion test via sd of residuals fitted vs. simulated



Simulated values, red line = fitted model. p-value (two.sided) = 0.336

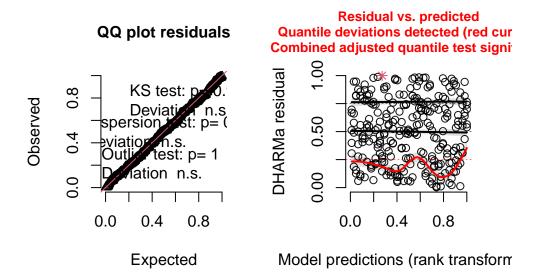
DHARMa nonparametric dispersion test via sd of residuals fitted vs. simulated

data: simulationOutput

dispersion = 0.23324, p-value = 0.336
alternative hypothesis: two.sided

plot(simulated\_resids)

#### DHARMa residual



#### ANOVA

car::Anova(m1)

Analysis of Deviance Table (Type II Wald chisquare tests)

Response: insect\_counts

Chisq Df Pr(>Chisq)

treatment 54.358 5 1.769e-10 \*\*\*
Date 41.652 8 1.574e-06 \*\*\*

---

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

glmmTMB is compatible with emmeans and effects.

### 4 Special Conditions

### 4.1 Split plot with repeated measures

```
Main plot is "irrigation" and split plot is "mix".
   alfalfa_sp <- read.csv("data/alfalfa2021_data.csv")</pre>
  library(dplyr)
Attaching package: 'dplyr'
The following objects are masked from 'package:stats':
    filter, lag
The following objects are masked from 'package:base':
    intersect, setdiff, setequal, union
cut: a cutting (harvest) of alfalfa within a single growing season. This is a temporal unit for
repeated measures analysis. There were three cuttings in total for that year and field. The
dates are not known, but we cannot assume they are evenly spaced apart.
irrigation: irrigation treatment ("Full" or "Deficit")
plot: a unique number referring to each experimental unit
block: the blocking unit
yield: response variable
row: plot position for row
col: plot positions for column or range
  head(alfalfa_sp)
```

```
cut irrigation plot block
                                mix
                                       yield row col
1 First
             Full 1101
                           1 50A+500 221.0418
                                                   1
2 First
             Full 1102
                           1 75A+250 288.7987
                                                   2
3 First
             Full 1103
                           1 50A+50F 466.7924
                                                   3
4 First
                           1 75A+25M 556.9506
                                                   4
             Full 1104
5 First
             Full 1105
                           1 50A+50M 422.9160
                                                   5
             Full 1106 1 75A+25F 289.8350
6 First
                                              2 1
```

Two new variables created:

**rep**: factor version of block (We should treat rep/block as a factor rather than an integer in modelling)

Cut: number version of cut where 1 is the first cutting. This is required by nlme::lme for specialized correlation structures.

```
alfalfa_sp <- alfalfa_sp %>%
  mutate(rep = as.factor(block)) %>%
  mutate(Cut = case_when(
    cut == "First" ~ 1L,
    cut == "Second" ~ 2L,
    cut == "Third" ~ 3L,
    is.na(cut) ~ NA_integer_))
```

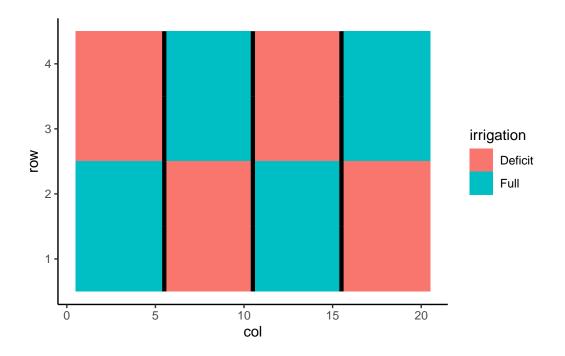
Visualise data

```
library(ggplot2); library(desplot)

alfalfa_sp %>% filter(cut == "First") %>%

ggplot(aes(x = col, y = row)) +
   geom_raster(aes(fill = irrigation)) +
   geom_tileborder(aes(group = 1, grp = rep), lwd = 1.5) +
   theme_classic()
```

Warning: Using `size` aesthetic for lines was deprecated in ggplot2 3.4.0. i Please use `linewidth` instead.



Model statement

$$y_{ijk} = \mu + \alpha_i + \beta_j + \gamma_k + a_l + b_m + c_n + \epsilon$$

where

 $\mu = {\rm overall~mean/intercept}$ 

 $\alpha_i$  = effect of the  $i^{th}$  irrigation treatment  $\beta_j$  = effect of the  $j^{th}$  planting mix treatment  $\gamma_k$  = effect of the  $k^{th}$  cutting [[need all those interactions]]

library(nlme)

Attaching package: 'nlme'

The following object is masked from 'package:dplyr':

collapse

use a special correlation structure for correlated error terms <code>corCompSymm()</code> is for compound symmetry. There are several other options in the <code>nlm</code> machinery (search "cor" for more options and details on the syntax). In general, repeated measures syntax follow this convention: <code>form = ~ time|grouping</code>. You can also use <code>l|group</code> and the observation order for each group will be. The default starting value (<code>value</code>) is zero, and if <code>fixed = FALSE</code> (the current nlme default), this value will be allowed to change during the model fitting process.

It's important that these two terms match after the "|" in the random and form arguments:

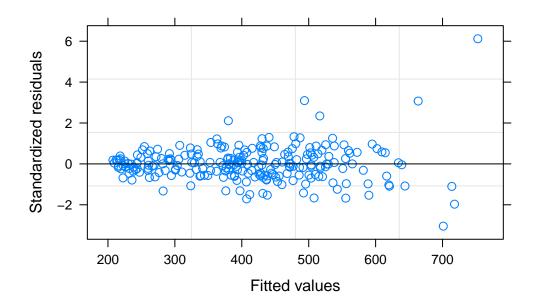
Update the model:

```
m2 <- update(m1, cor = corstr)</pre>
```

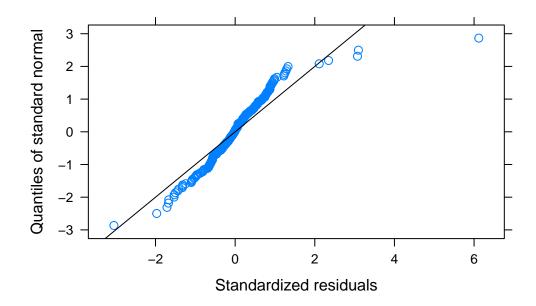
The usual next steps:

check diagnostics

```
plot(m2)
```



$$qqnorm(m2, \sim resid(., type = "p"), abline = c(0, 1))$$



Look at the variance components.

### VarCorr(m2)

	Variance	StdDev
rep =	pdLogChol(1)	
(Intercept)	83.17553	9.120062
irrigation =	pdLogChol(1)	
(Intercept)	280.54819	16.749573
plot =	pdLogChol(1)	
(Intercept)	481.44292	21.941808
Residual	16182.24438	127.209451

### Run ANOVA

### anova(m2)

	${\tt numDF}$	${\tt denDF}$	F-value	p-value
(Intercept)	1	102	1432.6369	<.0001
mix	9	102	13.6932	<.0001
irrigation	1	3	4.8770	0.1143
cut	2	102	6.0434	0.0033
mix:irrigation	9	102	0.5256	0.8530
mix:cut	18	102	0.8029	0.6927
irrigation:cut	2	102	14.2649	<.0001
mix:irrigation:cut	18	102	1.0226	0.4418

always check the degrees of freedom (denominator and numerator)!

# Summary

In summary, mixed models are complicated.

## References