warmXtrophic Project: CN Analyses

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Load in packages & data

```
# Clear all existing data
rm(list = ls())

# Load packages
library(tidyverse)
library(bbmle)
library(lmerTest)

# Set working directory to Google Drive
setwd("/Volumes/GoogleDrive/Shared drives/SpaCE_Lab_warmXtrophic/data/")

# Read in data
cn <- read.csv("L1/final_CN_L1.csv")

# create dataframes for kbs and umbs
cn_kbs <- subset(cn, site == "kbs")
cn_umbs <- subset(cn, site == "umbs")</pre>
```

Check what species are measured at each site

```
unique(cn_kbs$species)

## [1] "Soca"
unique(cn_umbs$species)

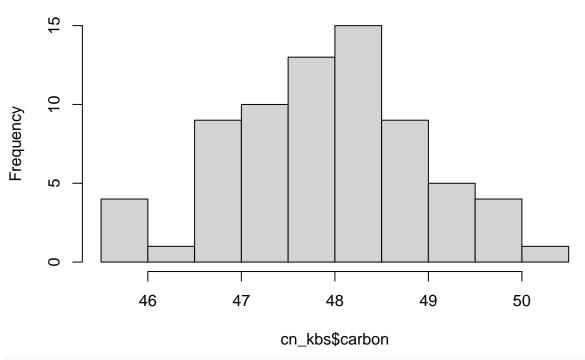
## [1] "Cest" "Popr"

# separate dataframes for each species
cn_cest_umbs <- subset(cn_umbs, species == "Cest")
cn_popr_umbs <- subset(cn_umbs, species == "Popr")</pre>
```

Data exploration: KBS Soca

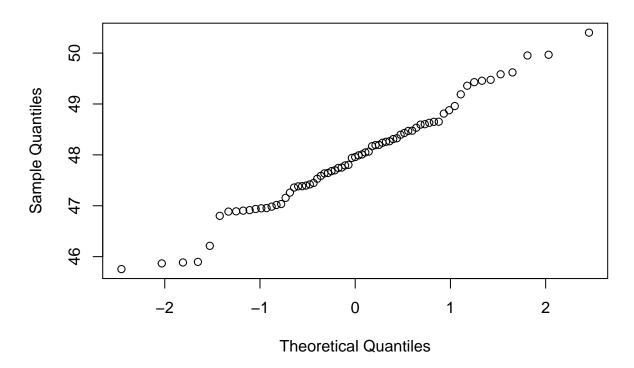
hist(cn_kbs\$carbon)

Histogram of cn_kbs\$carbon



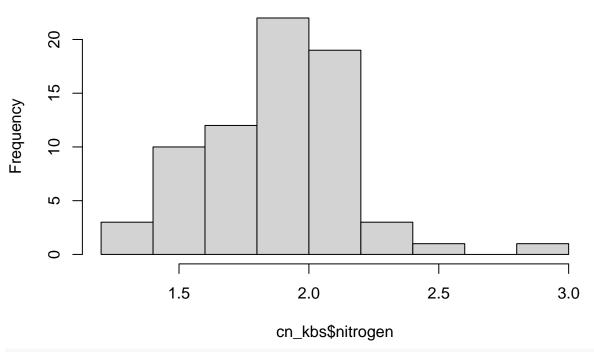
qqnorm(cn_kbs\$carbon)

Normal Q-Q Plot



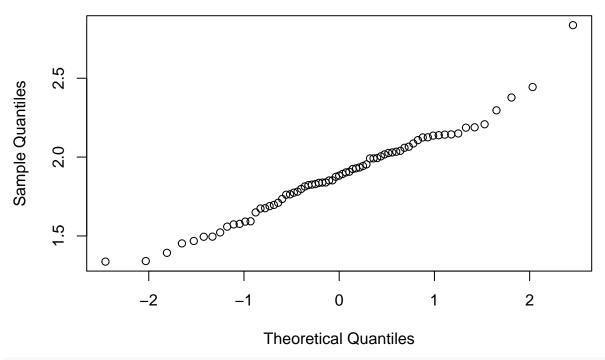
shapiro.test(cn_kbs\$carbon) ## ## Shapiro-Wilk normality test ## ## data: cn_kbs\$carbon ## W = 0.98512, p-value = 0.5658 hist(cn_kbs\$nitrogen)

Histogram of cn_kbs\$nitrogen



qqnorm(cn_kbs\$nitrogen)

Normal Q-Q Plot



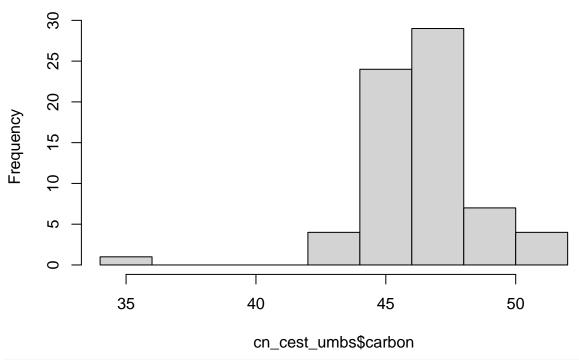
shapiro.test(cn_kbs\$nitrogen)

```
##
## Shapiro-Wilk normality test
##
## data: cn_kbs$nitrogen
## W = 0.97562, p-value = 0.1819
```

Data exploration: UMBS Cest

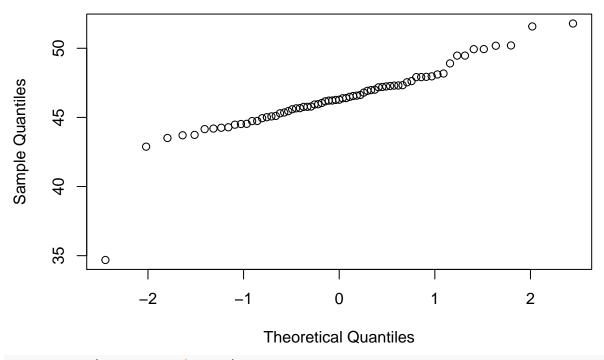
hist(cn_cest_umbs\$carbon)

Histogram of cn_cest_umbs\$carbon



qqnorm(cn_cest_umbs\$carbon)

Normal Q-Q Plot

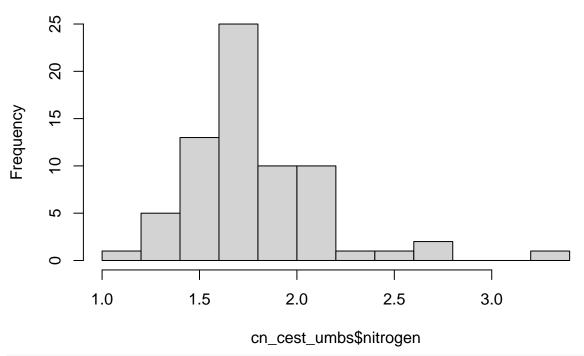


shapiro.test(cn_cest_umbs\$carbon)

##

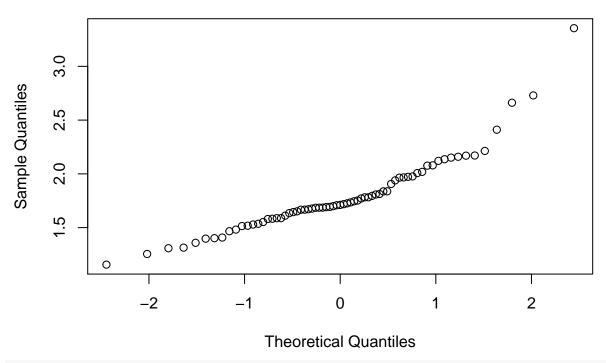
```
## Shapiro-Wilk normality test
##
## data: cn_cest_umbs$carbon
## W = 0.87921, p-value = 7.329e-06
hist(cn_cest_umbs$nitrogen)
```

Histogram of cn_cest_umbs\$nitrogen



qqnorm(cn_cest_umbs\$nitrogen)

Normal Q-Q Plot



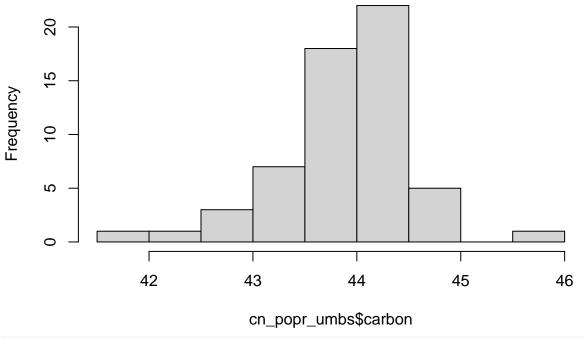
```
shapiro.test(cn_cest_umbs$nitrogen)
```

```
##
## Shapiro-Wilk normality test
##
## data: cn_cest_umbs$nitrogen
## W = 0.88786, p-value = 1.507e-05
```

Data exploration: UMBS Popr

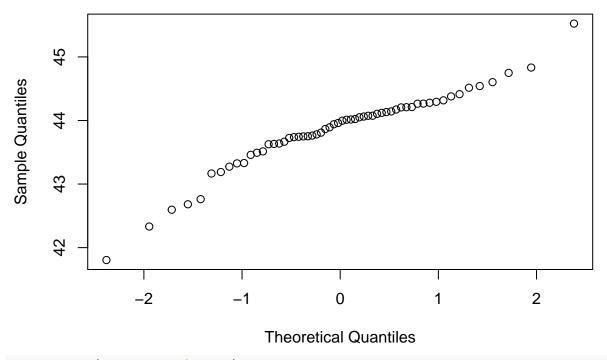
hist(cn_popr_umbs\$carbon)

Histogram of cn_popr_umbs\$carbon



qqnorm(cn_popr_umbs\$carbon)

Normal Q-Q Plot



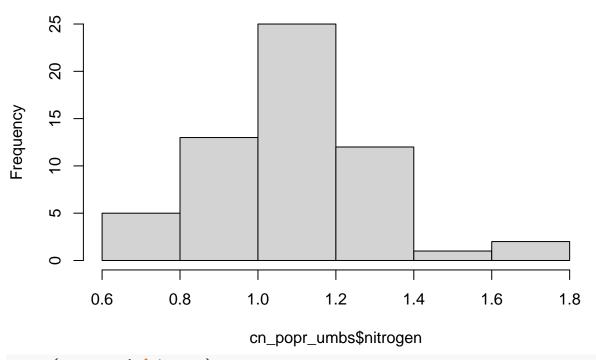
shapiro.test(cn_popr_umbs\$carbon)

##

```
## Shapiro-Wilk normality test
##
## data: cn_popr_umbs$carbon
## W = 0.93934, p-value = 0.006065
```

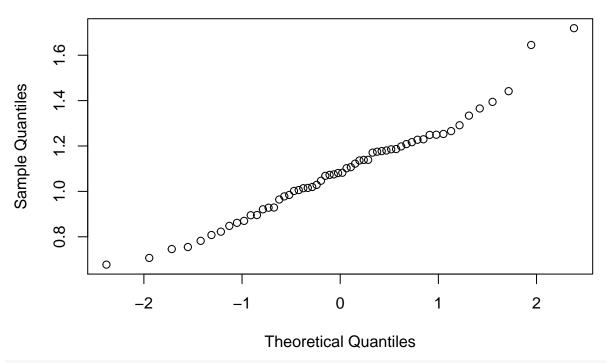
hist(cn_popr_umbs\$nitrogen)

Histogram of cn_popr_umbs\$nitrogen



qqnorm(cn_popr_umbs\$nitrogen)

Normal Q-Q Plot



shapiro.test(cn_popr_umbs\$nitrogen)

```
##
## Shapiro-Wilk normality test
##
## data: cn_popr_umbs$nitrogen
## W = 0.97376, p-value = 0.2403
```

Model comparison: KBS Soca

```
# carbon
m1a <- lm(carbon ~ state, data = cn_kbs)</pre>
m1b <- lm(carbon ~ state + insecticide, data = cn_kbs)</pre>
m1c <- lmer(carbon ~ state + (1 | plot), data = cn kbs)
m1d <- lmer(carbon ~ state + insecticide + (1 | plot), data = cn_kbs)
AICctab(m1a, m1b, m1c, m1d, weights = T)
##
       dAICc df weight
## m1a 0.0
           3 0.7098
           4 0.2332
## m1b 2.2
           4 0.0473
## m1c 5.4
## m1d 8.6
           5 0.0097
summary(m1a)
##
## Call:
## lm(formula = carbon ~ state, data = cn_kbs)
## Residuals:
##
       Min
                 1Q
                     Median
                                   3Q
## -2.26732 -0.72064 0.04298 0.55648 2.24858
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 47.7409 0.1673 285.28
                                            <2e-16 ***
## statewarmed 0.4125
                           0.2383
                                             0.088 .
                                     1.73
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 1.004 on 69 degrees of freedom
## (2 observations deleted due to missingness)
## Multiple R-squared: 0.04159,
                                   Adjusted R-squared: 0.0277
## F-statistic: 2.995 on 1 and 69 DF, p-value: 0.08802
summary(m1c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: carbon ~ state + (1 | plot)
     Data: cn_kbs
##
##
## REML criterion at convergence: 203.2
## Scaled residuals:
       \mathtt{Min}
              1Q
                     Median
                                   3Q
## -2.14660 -0.60048 -0.01159 0.56066 2.33512
## Random effects:
## Groups
                        Variance Std.Dev.
           Name
## plot
            (Intercept) 0.07219 0.2687
## Residual
                        0.94047 0.9698
## Number of obs: 71, groups: plot, 24
```

```
##
## Fixed effects:
              Estimate Std. Error
                                       df t value Pr(>|t|)
## (Intercept) 47.7386
                        0.1795 20.2987 265.909 <2e-16 ***
## statewarmed 0.4129
                           0.2553 20.8739
                                            1.617
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Correlation of Fixed Effects:
##
               (Intr)
## statewarmed -0.703
# nitrogen
m2a <- lm(nitrogen ~ state, data = cn_kbs)</pre>
m2b <- lm(nitrogen ~ state + insecticide, data = cn_kbs)</pre>
m2c <- lmer(nitrogen ~ state + (1 | plot), data = cn_kbs)</pre>
m2d <- lmer(nitrogen ~ state + insecticide + (1 | plot), data = cn_kbs)
AICctab(m2a, m2b, m2c, m2d, weights = T)
      dAICc df weight
## m2b 0.0
           4 0.5138
## m2a 0.2
           3 0.4574
           4 0.0251
## m2c 6.0
## m2d 9.9
           5 0.0037
summary(m2a)
##
## Call:
## lm(formula = nitrogen ~ state, data = cn_kbs)
##
## Residuals:
       Min
                 1Q Median
## -0.47449 -0.17566 -0.02393 0.16347 1.02461
##
## Coefficients:
##
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 1.95273 0.04392 44.458
                                            <2e-16 ***
## statewarmed -0.14094
                          0.06256 - 2.253
                                            0.0274 *
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 0.2635 on 69 degrees of freedom
     (2 observations deleted due to missingness)
## Multiple R-squared: 0.06852,
                                  Adjusted R-squared: 0.05502
## F-statistic: 5.076 on 1 and 69 DF, p-value: 0.02745
summary(m2c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: nitrogen ~ state + (1 | plot)
##
      Data: cn_kbs
##
## REML criterion at convergence: 13.7
##
```

```
## Scaled residuals:
##
      Min 1Q Median
                           3Q
                                    Max
## -1.8561 -0.5037 -0.1174 0.5903 3.1426
##
## Random effects:
## Groups Name
                       Variance Std.Dev.
            (Intercept) 0.02190 0.148
## plot
## Residual
                       0.04928 0.222
## Number of obs: 71, groups: plot, 24
##
## Fixed effects:
              Estimate Std. Error
                                      df t value Pr(>|t|)
##
## (Intercept) 1.95576 0.05665 20.85402 34.523 <2e-16 ***
## statewarmed -0.14870
                         0.08032 21.09931 -1.851
                                                   0.0782 .
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Correlation of Fixed Effects:
##
              (Intr)
## statewarmed -0.705
```

Model comparison: UMBS Cest

```
# carbon
m3a <- lm(carbon ~ state, data = cn_cest_umbs)
m3b <- lm(carbon ~ state + insecticide, data = cn cest umbs)
m3c <- lmer(carbon ~ state + (1 | plot), data = cn_cest_umbs)</pre>
m3d <- lmer(carbon ~ state + insecticide + (1 | plot), data = cn_cest_umbs)
AICctab(m3a, m3b, m3c, m3d, weights = T)
##
      dAICc df weight
           3 0.48
## m3a 0.0
## m3c 1.4
            4 0.23
## m3b 2.1
            4 0.17
## m3d 2.7
           5 0.12
summary(m3a)
##
## Call:
## lm(formula = carbon ~ state, data = cn_cest_umbs)
##
## Residuals:
       Min
                 1Q Median
                                    3Q
                                            Max
## -11.9371 -1.1730 0.0917 1.1498
                                       5.1632
##
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 46.1235
                           0.4067 113.408
                                             <2e-16 ***
## statewarmed 0.5035
                           0.5710
                                    0.882
                                              0.381
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 2.371 on 67 degrees of freedom
```

```
## Multiple R-squared: 0.01147, Adjusted R-squared: -0.003283
## F-statistic: 0.7775 on 1 and 67 DF, p-value: 0.3811
summary(m3c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: carbon ~ state + (1 | plot)
     Data: cn_cest_umbs
##
##
## REML criterion at convergence: 312.1
## Scaled residuals:
##
      Min
              1Q Median
                               ЗQ
                                      Max
## -5.0954 -0.4926 0.0019 0.4734 2.4473
##
## Random effects:
## Groups Name
                        Variance Std.Dev.
            (Intercept) 0.6428 0.8018
## plot
## Residual
                        5.0133
                                 2.2390
## Number of obs: 69, groups: plot, 24
##
## Fixed effects:
              Estimate Std. Error
                                       df t value Pr(>|t|)
## (Intercept) 46.1270
                        0.4492 23.0887 102.695
                                                    <2e-16 ***
                           0.6316 22.6923 0.786
## statewarmed 0.4966
                                                      0.44
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Correlation of Fixed Effects:
##
              (Intr)
## statewarmed -0.711
# nitrogen
m4a <- lm(nitrogen ~ state, data = cn_cest_umbs)
m4b <- lm(nitrogen ~ state + insecticide, data = cn_cest_umbs)
m4c <- lmer(nitrogen ~ state + (1 | plot), data = cn_cest_umbs)</pre>
m4d <- lmer(nitrogen ~ state + insecticide + (1 | plot), data = cn_cest_umbs)
AICctab(m4a, m4b, m4c, m4d, weights = T)
      dAICc df weight
## m4a 0.0 3 0.7383
## m4b 2.1 4 0.2528
## m4c 9.0 4 0.0082
## m4d 14.1 5 <0.001
summary(m4a)
##
## Call:
## lm(formula = nitrogen ~ state, data = cn_cest_umbs)
##
## Residuals:
##
      Min
               1Q Median
                               3Q
                                      Max
## -0.5740 -0.1970 -0.1005 0.1940 1.4726
##
```

```
## Coefficients:
##
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 1.69764
                         0.05935 28.605
## statewarmed 0.18473
                          0.08333
                                             0.03 *
                                   2.217
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 0.346 on 67 degrees of freedom
## Multiple R-squared: 0.06834, Adjusted R-squared: 0.05444
## F-statistic: 4.915 on 1 and 67 DF, p-value: 0.03003
summary(m4c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: nitrogen ~ state + (1 | plot)
     Data: cn_cest_umbs
## REML criterion at convergence: 54.1
##
## Scaled residuals:
      Min
           1Q Median
                              3Q
                                     Max
## -1.5491 -0.5223 -0.2124 0.4340 3.9535
##
## Random effects:
## Groups Name
                       Variance Std.Dev.
## plot
            (Intercept) 0.01481 0.1217
## Residual
                        0.10569 0.3251
## Number of obs: 69, groups: plot, 24
## Fixed effects:
              Estimate Std. Error
                                       df t value Pr(>|t|)
## (Intercept) 1.69816
                         0.06602 22.98884 25.721
                                                    <2e-16 ***
## statewarmed 0.18338
                         0.09285 22.59847
                                           1.975
                                                    0.0606 .
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Correlation of Fixed Effects:
              (Intr)
## statewarmed -0.711
```

Model comparison: UMBS Popr

```
# carbon
m5a <- lm(carbon ~ state, data = cn_popr_umbs)
m5b <- lm(carbon ~ state + insecticide, data = cn_popr_umbs)
m5c <- lmer(carbon ~ state + (1 | plot), data = cn_popr_umbs)
m5d <- lmer(carbon ~ state + insecticide + (1 | plot), data = cn_popr_umbs)
AICctab(m5a, m5b, m5c, m5d, weights = T)

## dAICc df weight
## m5a 0.0 3 0.7260
## m5b 2.2 4 0.2365
## m5c 6.2 4 0.0328</pre>
```

```
## m5d 10.1 5 0.0047
summary(m5a)
##
## lm(formula = carbon ~ state, data = cn_popr_umbs)
## Residuals:
      Min
              1Q Median
                               3Q
                                     Max
## -1.9645 -0.2155 0.0791 0.3955 1.5634
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
##
## (Intercept) 43.7693
                        0.1105 395.993 <2e-16 ***
## statewarmed 0.1915
                           0.1620 1.182
                                            0.242
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 0.6154 on 56 degrees of freedom
## Multiple R-squared: 0.02435,
                                  Adjusted R-squared: 0.006928
## F-statistic: 1.398 on 1 and 56 DF, p-value: 0.2421
summary(m5c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: carbon ~ state + (1 | plot)
     Data: cn_popr_umbs
##
##
## REML criterion at convergence: 110.1
## Scaled residuals:
           1Q Median
##
      Min
                               3Q
## -3.0725 -0.3584 0.1327 0.5730 2.1594
##
## Random effects:
## Groups Name
                        Variance Std.Dev.
## plot
            (Intercept) 0.05839 0.2416
## Residual
                        0.32389 0.5691
## Number of obs: 58, groups: plot, 21
##
## Fixed effects:
              Estimate Std. Error
##
                                      df t value Pr(>|t|)
## (Intercept) 43.7747 0.1258 18.7118 347.948
                                                   <2e-16 ***
## statewarmed 0.1778
                          0.1842 18.9535
                                          0.965
                                                    0.346
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Correlation of Fixed Effects:
##
              (Intr)
## statewarmed -0.683
# nitrogen
m6a <- lm(nitrogen ~ state, data = cn_popr_umbs)</pre>
m6b <- lm(nitrogen ~ state + insecticide, data = cn_popr_umbs)</pre>
```

```
m6c <- lmer(nitrogen ~ state + (1 | plot), data = cn_popr_umbs)</pre>
m6d <- lmer(nitrogen ~ state + insecticide + (1 | plot), data = cn_popr_umbs)
AICctab(m6a, m6b, m6c, m6d, weights = T)
      dAICc df weight
## m6a 0.0 3 0.7542
## m6b 2.3 4 0.2376
## m6c 9.1 4 0.0079
## m6d 15.4 5 <0.001
summary(m6a)
##
## Call:
## lm(formula = nitrogen ~ state, data = cn_popr_umbs)
## Residuals:
       Min
                 1Q
                     Median
                                  3Q
                                         Max
## -0.36742 -0.11669 -0.00407 0.07251 0.52918
##
## Coefficients:
              Estimate Std. Error t value Pr(>|t|)
## (Intercept) 1.19002
                         0.03227
                                  36.88 < 2e-16 ***
## statewarmed -0.22326
                         0.04730
                                  -4.72 1.62e-05 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 0.1797 on 56 degrees of freedom
## Multiple R-squared: 0.2846, Adjusted R-squared: 0.2719
## F-statistic: 22.28 on 1 and 56 DF, p-value: 1.615e-05
summary(m6c) # plot as a random effect
## Linear mixed model fit by REML. t-tests use Satterthwaite's method [
## lmerModLmerTest]
## Formula: nitrogen ~ state + (1 | plot)
     Data: cn_popr_umbs
## REML criterion at convergence: -29.8
## Scaled residuals:
      Min
               1Q Median
                              3Q
                                     Max
## -1.9564 -0.6459 -0.0134 0.5663 2.5154
## Random effects:
## Groups Name
                       Variance Std.Dev.
## plot
            (Intercept) 0.008205 0.09058
## Residual
                       0.024490 0.15649
## Number of obs: 58, groups: plot, 21
##
## Fixed effects:
                                      df t value Pr(>|t|)
              Estimate Std. Error
## (Intercept) 1.18916 0.03929 19.10091 30.26 < 2e-16 ***
## ---
```

```
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Correlation of Fixed Effects:
## (Intr)
## statewarmed -0.684
```