

Analyses of species distributions in peatlands

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Load the R packages that you will use

If you do not have the R packages installed, you need to install them.

```
library(tidyverse)
library(readxl)
library(knitr)
library(ggeffects)
library(car)
library(glmTMB)
library(ggplot2)
library(BiodiversityR)
library(vegan)
```

Data preparation

Read data from Excel file

Note that you need to change the path to the folder where you have the Excel file

```
data_peat<-read_excel("data/raw/Modelling_SDM_species_data_Updated.xlsx",
                      sheet="Sheet1")
data_peat<-data_peat[1:29] # This is to remove some columns that we do not need!
```

Have a look at the data

This shows the first rows of your data file in “tibble” format. You can also see the variable type for each variable (double or character).

data_peat

```
## # A tibble: 115 x 29
##   n_samples depth depth_corrected fen tot_Sphagnum Erio Carex Erica
##   <dbl> <chr> <dbl> <chr> <dbl> <dbl> <dbl> <dbl>
## 1 1 0 0 N 100 0 0 0
## 2 2 14-15 14 N 87 3 2 8
## 3 3 20-21 20 N 74 6 6 14
## 4 4 30-31 30 N 87 5 4 4
## 5 5 40-41 40 N 90 2 3 5
## 6 6 50-51 50 N 41 46 9 4
## 7 7 55-56 55 N 74 19 6 1
## 8 8 60-61 60 N 94 4 1 1
## 9 9 65-66 65 N 95 3 1 1
## 10 10 70-71 70 N 90 3 4 3
##   other_veg Balticum Medium Cuspidata Austinii Fuscum Rubellum Acutifolia
##   <dbl> <dbl> <dbl> <dbl> <dbl> <dbl> <dbl> <dbl>
## 1 0 0.32 0.68 0 0 0 0 0
## 2 0 0.724 0.253 0.0230 0 0 0 0
## 3 0 0.324 0.649 0 0 0 0 0.0270
## 4 0 0.264 0.276 0 0 0 0.368 0.0920
## 5 0 0.189 0.0556 0.0333 0 0 0 0.722
## 6 0 0.171 0.415 0.220 0 0 0.195 0
## 7 0 0.568 0.216 0.216 0 0 0 0
## 8 0 0.489 0.383 0.128 0 0 0 0
## 9 0 0.589 0.263 0.147 0 0 0 0
## 10 0 0.544 0.111 0.344 0 0 0 0
##   'Diseased Acutifolia' Angustifolium Tenellum Papillosum Fallax Stems age
##   <dbl> <dbl> <dbl> <dbl> <dbl> <dbl> <dbl> <dbl>
## 1 0 0 0 0 0 0 0 0
## 2 0 0 0 0 0 0 0 75
## 3 0 0 0 0 0 0 0 136
## 4 0 0 0 0 0 0 0 244
## 5 0 0 0 0 0 0 0 352
## 6 0 0 0 0 0 0 0 452
## 7 0 0 0 0 0 0 0 505
## 8 0 0 0 0 0 0 0 555
## 9 0 0 0 0 0 0 0 606
## 10 0 0 0 0 0 0 0 643
##   temp imp_temp moist nutrient fire dry
##   <dbl> <dbl> <chr> <dbl> <dbl> <dbl>
## 1 6.35 0 NA 0 0 0
## 2 6.92 1 NA 0 0 0
## 3 7.39 1 NA 0 0 0
## 4 8.21 0 NA 0 0 0
## 5 8.00 1 -0.38 1 0 0
## 6 7.81 1 -0.3 1 0 0
## 7 7.71 1 0.01 1 0 0
## 8 7.62 1 -0.26 1 0 0
```

```
## 9 7.52      1 0.22      1      0      0
## 10 7.45     1 -0.25     1      0      0
## # i 105 more rows
```

Convert some variables to factors

It is better to convert some variables (those that are Y/N or 0/1) to factors.

```
data_peat<-data_peat%>%
  mutate(fen=as.factor(fen),imp_temp=as.factor(imp_temp),
         nutrient=as.factor(nutrient),fire=as.factor(fire),dry=as.factor(dry))
# with mutate you create new variables that are equal to the old variables
# but are coded as factors
```

Convert moist to numeric

For some reason, moist appears as a character variable. It should be numeric, so we convert it.

```
data_peat<-data_peat%>%
  mutate(moist=as.numeric(moist))
```

Ordinations (vegan package)

Suggested reading: <https://www.davidzeleny.net/anadat-r/doku.php/en:ordination>

(lots of info on this webpage!)

Chapter 10 in this pdf: <https://apps.worldagroforestry.org/downloads/Publications/PDFS/b13695.pdf>

Using the vegan package.

I performed some ordinations with Sphagnum species. I tried different methods, but if I would need to choose one, I would do a constrained ordination, specifically a Distance-based redundancy analysis (db-RDA) with Bray-Curtis distance. You can read about all types in the webpage above if you feel like it.

Data for ordination:

```
data_ordi2<-data_peat %>%
  filter_at(vars(Balticum:Fallax),
            all_vars(!is.na(.)))%>% # Remove rows with all NAs
  filter_at(vars(Balticum:Fallax),
            any_vars(>0))%>% # Remove rows with all zeros - WHY?
  filter(!is.na(age)&!is.na(temp)&!is.na(moist)&!is.na(nutrient)&!is.na(fire)&
         !is.na(dry))%>%
  rename(Diseased_Acutifolia=`Diseased Acutifolia`) #Rename to avoid problems
```

Distance-based redundancy analysis (db-RDA) with Bray-Curtis distance.

See <https://www.davidzeleny.net/anadat-r/doku.php/en:similarity> for info on distances.

Calculate ordination:

```
ordi6<-capscale(data_ordi2[10:21]~ # species data matrix
                age+temp+moist+nutrient+fire+dry, # Environmental variables
                data = data_ordi2, distance="bray") # Bray-Curtis distance
```

Result of the ordination:

```
ordi6
```

```
## Call: capscale(formula = data_ordi2[10:21] ~ age + temp + moist +
## nutrient + fire + dry, data = data_ordi2, distance = "bray")
##
##              Inertia Proportion Rank
## Total          31.0246      1.0000
## Constrained     8.3892      0.2704    6
## Unconstrained  26.3948      0.8508   34
## Imaginary      -3.7594     -0.1212   38
## Inertia is squared Bray distance
## Species scores projected from '[' 'data_ordi2' '10:21'
##
## Eigenvalues for constrained axes:
## CAP1 CAP2 CAP3 CAP4 CAP5 CAP6
## 4.886 2.239 0.760 0.276 0.138 0.090
##
## Eigenvalues for unconstrained axes:
## MDS1 MDS2 MDS3 MDS4 MDS5 MDS6 MDS7 MDS8
## 7.438 4.231 3.328 2.484 2.345 2.009 1.230 0.598
## (Showing 8 of 34 unconstrained eigenvalues)
```

“Intertia” is the total variance - your environmental variables explain 0.2704 of this variance (“constrained” part).

Proportion explained by each ordination axis. CAP1-CAP6 are the “constrained” axes, explained by your environmental variables. MDS1-MDS34 are the “unconstrained” axes.

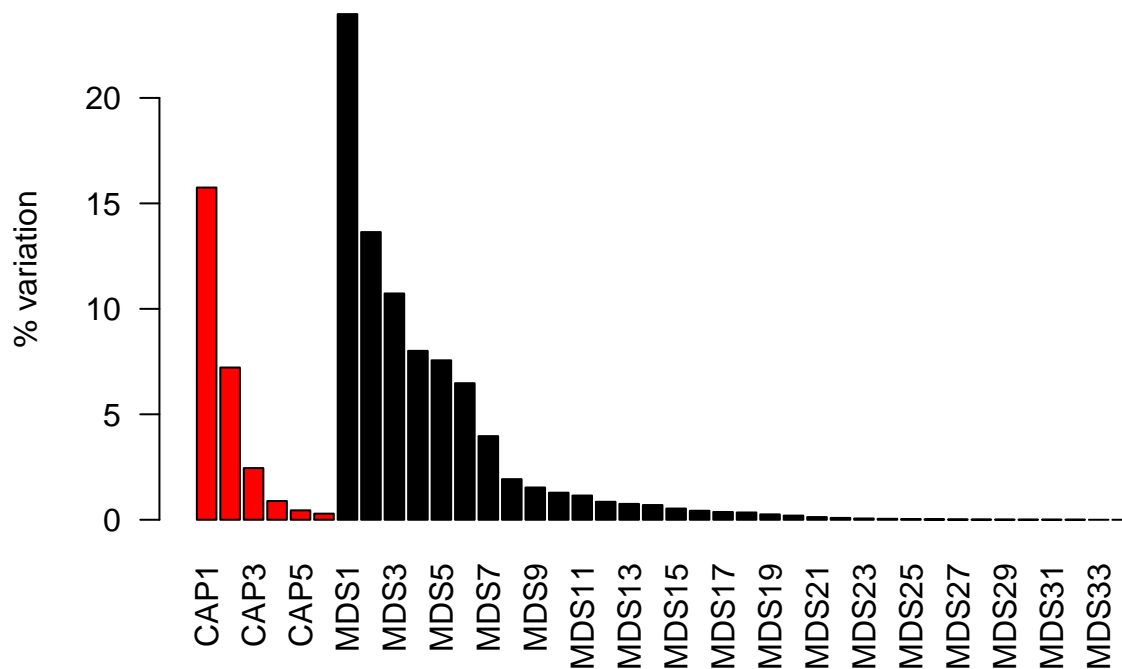
```
eigenvals(ordi6) %>%
summary()
```

```
## Importance of components:
##              CAP1      CAP2      CAP3      CAP4      CAP5      CAP6      MDS1
## Eigenvalue      4.8863 2.23908 0.76030 0.275526 0.138265 0.089651 7.4376
## Proportion Explained 0.1405 0.06437 0.02186 0.007921 0.003975 0.002577 0.2138
## Cumulative Proportion 0.1405 0.20485 0.22671 0.234627 0.238602 0.241179 0.4550
##              MDS2      MDS3      MDS4      MDS5      MDS6      MDS7      MDS8
## Eigenvalue      4.2314 3.32835 2.48447 2.34511 2.00855 1.23007 0.59751
## Proportion Explained 0.1216 0.09569 0.07143 0.06742 0.05774 0.03536 0.01718
## Cumulative Proportion 0.5766 0.67234 0.74376 0.81118 0.86892 0.90429 0.92147
##              MDS9      MDS10      MDS11      MDS12      MDS13      MDS14
## Eigenvalue      0.47443 0.39821 0.35601 0.264468 0.233343 0.216337
## Proportion Explained 0.01364 0.01145 0.01023 0.007603 0.006708 0.006219
## Cumulative Proportion 0.93510 0.94655 0.95679 0.964391 0.971099 0.977319
##              MDS15      MDS16      MDS17      MDS18      MDS19      MDS20
## Eigenvalue      0.165074 0.132727 0.114648 0.10713 0.079384 0.061005
```

```
## Proportion Explained 0.004746 0.003816 0.003296 0.00308 0.002282 0.001754
## Cumulative Proportion 0.982065 0.985880 0.989176 0.99226 0.994538 0.996292
##                      MDS21      MDS22      MDS23      MDS24      MDS25
## Eigenvalue          0.039707 0.0272079 0.0186386 0.0140238 0.0100911
## Proportion Explained 0.001142 0.0007822 0.0005358 0.0004032 0.0002901
## Cumulative Proportion 0.997434 0.9982159 0.9987517 0.9991549 0.9994450
##                      MDS26      MDS27      MDS28      MDS29      MDS30
## Eigenvalue          0.0072677 0.0045643 3.108e-03 1.815e-03 1.313e-03
## Proportion Explained 0.0002089 0.0001312 8.935e-05 5.217e-05 3.774e-05
## Cumulative Proportion 0.9996539 0.9997851 9.999e-01 9.999e-01 1.000e+00
##                      MDS31      MDS32      MDS33      MDS34
## Eigenvalue          7.435e-04 4.026e-04 9.116e-05 9.039e-07
## Proportion Explained 2.138e-05 1.157e-05 2.621e-06 2.599e-08
## Cumulative Proportion 1.000e+00 1.000e+00 1.000e+00 1.000e+00
```

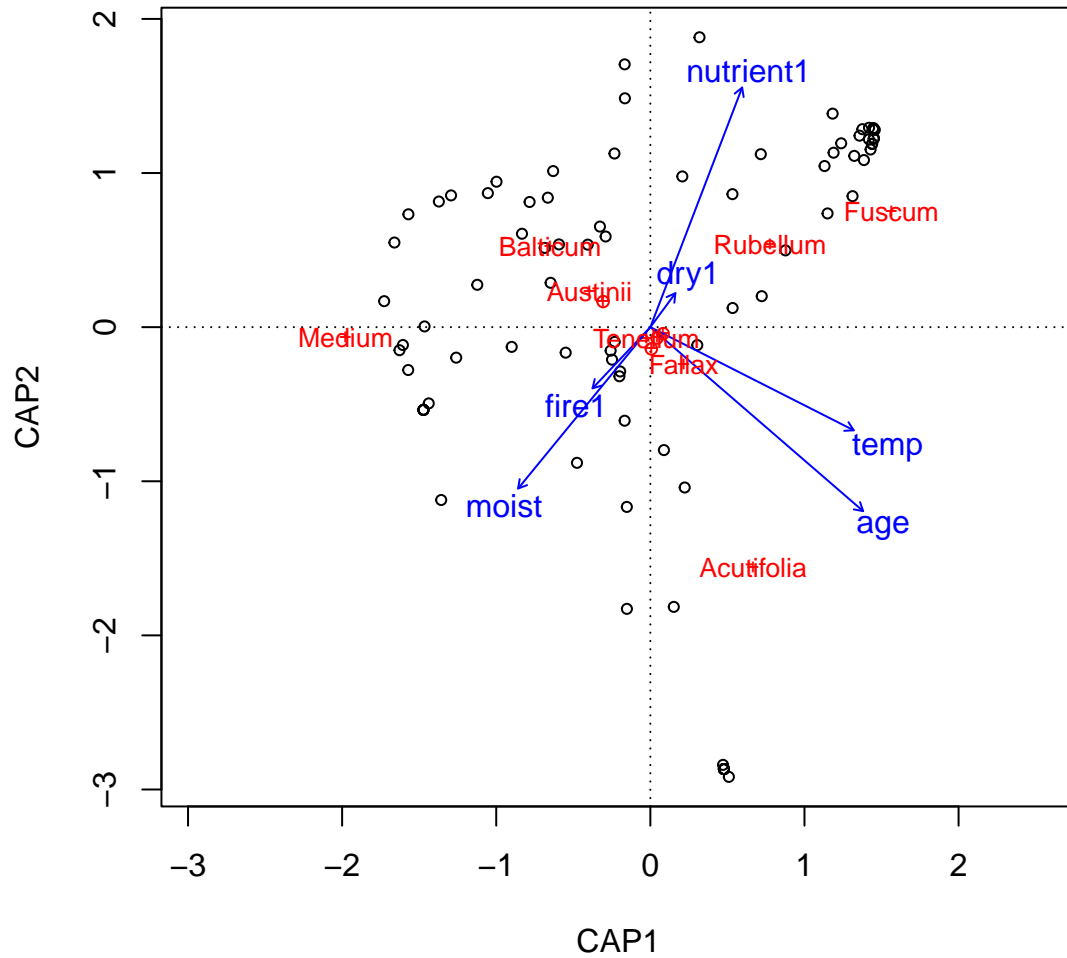
Barplot of percentage variance explained by individual axes

```
expl_var_ordi6 <- c(ordi6$CCA$eig/ordi6$tot.chi*100,
                    ordi6$CA$eig/ordi6$tot.chi*100)
barplot (expl_var_ordi6, col = c(rep ('red',
                                     length (ordi6$CCA$eig/ordi6$tot.chi*100)),
                                rep ('black',
                                     length (ordi6$CA$eig/ordi6$tot.chi*100))),
        las = 2, ylab = '% variation')
```



Plot of the ordination (species in red and sites-samples in black):

```
vegan::ordiplot(ordi6, display = c('species', 'sites', 'bp'))
orditorp(ordi6, display="species", cex=0.8, col="red")
```



This shows the two first constrained axes of the ordination. You can see how the sites and species distribute along these axes.

Test significance of the ordination with Monte Carlo permutation test.

For the whole model:

```
anova (ordi6, permutations = 999)
```

```
## Permutation test for capscale under reduced model
## Permutation: free
## Number of permutations: 999
##
## Model: capscale(formula = data_ordi2[10:21] ~ age + temp + moist + nutrient + fire + dry, data = data_ordi2)
```

```
##           Df SumOfSqs      F Pr(>F)
## Model      6   8.3892 4.2378 0.001 ***
## Residual 80  26.3948
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

The model is significant.

For each explanatory variable (with all the others used as covariables, independently from their order in the model):

```
anova (ordi6, by = 'margin', permutations = 999)
```

```
## Permutation test for capscale under reduced model
## Marginal effects of terms
## Permutation: free
## Number of permutations: 999
##
## Model: capscale(formula = data_ordi2[10:21] ~ age + temp + moist + nutrient + fire + dry, data = dat
##           Df SumOfSqs      F Pr(>F)
## age        1   1.2989 3.9369 0.002 **
## temp        1   0.3226 0.9778 0.411
## moist       1   0.8304 2.5168 0.024 *
## nutrient    1   1.1292 3.4226 0.005 **
## fire        1   0.1146 0.3472 0.948
## dry         1   0.8159 2.4728 0.025 *
## Residual 80  26.3948
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

Age, moist, nutrient and dry show significant effects.

For each axis:

```
anova (ordi6, by = 'axis', permutations = 999)
```

```
## Permutation test for capscale under reduced model
## Forward tests for axes
## Permutation: free
## Number of permutations: 999
##
## Model: capscale(formula = data_ordi2[10:21] ~ age + temp + moist + nutrient + fire + dry, data = dat
##           Df SumOfSqs      F Pr(>F)
## CAP1       1   4.8863 14.8100 0.001 ***
## CAP2       1   2.2391  6.7864 0.002 **
## CAP3       1   0.7603  2.3044 0.315
## CAP4       1   0.2755  0.8351 0.972
## CAP5       1   0.1383  0.4191 0.997
## CAP6       1   0.0897  0.2717 0.989
## Residual 80  26.3948
## ---
## Signif. codes:  0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
```

Axis 1 and 2 are significant.

Ordination plot with ggplot2.

Install ggord package (you only need to do this once):

```
# Enable the r-universe repo
options(repos = c(
  fawda123 = 'https://fawda123.r-universe.dev',
  CRAN = 'https://cloud.r-project.org'))

# Install ggord
install.packages('ggord')

## package 'ggord' successfully unpacked and MD5 sums checked
##
## The downloaded binary packages are in
## C:\Users\alici\AppData\Local\Temp\RtmpMTfBdL\downloaded_packages
```

Load ggord package:

```
library(ggord)
```

```
ggord(ordi6,ptslab=T,repel=T,labcol="red",veccol="red",size=NA,addsize=3,
      xlims=c(-1.1,1.1),ylim=c(-1.1,1.1))+
  geom_point(size=3,shape=20,color="black",alpha=0.2)
```

