# Multivariate analysis: from 2D PCA to 3D PTA

Romain Frelat
20 october 2016

# **Objectives:**

At the end of this workshop, you should be able to:

- Run a Principal component analysis (PCA) on a matrix (2D)
- Interpret the Principal Components (PC)
- Run a Principal tensor analysis (PTA) on a array (3D)
- Interpret the Principal Tensor (PT)
- Run a clustering analysis with Hierarchical Clustering
- Understand what is a multivariate analysis, and when it can be useful

## A. Getting ready:

## Get ready:

- 1. Get the zip file Multivariate2D3D.zip (can be downloaded: https://github.com/rfrelat/Multivariate2D3D/)
- 2. Unzip the archive in a new folder. The zip file contain the data (*IBTS\_Tensor.Rdata*), the R-script (*script\_Multivariate2D3D.R*) and the present document as a pdf
- 3. Open the R script script\_Multivariate2D3D.R with your favourite R editor (RStudio is recommended)
- 4. Be sure to set the working directory (Session > Set Working Directory) to the directory where the script and the data are.

### Load the package and needed functions

For this work, we will need to use two package ade4 and PTAk.

```
library(ade4)
library(PTAk)
```

If you have an error message, check if the packages are installed. If not, use the function install.packages(c("ade4", "PTAk").

### Load the dataset

The data is a *Rdata* file, that can be loaded with the function load().

```
load("IBTS_Tensor.Rdata")
dim(IBTS_tensor)
```

[1] 65 31 7

You loaded a variable called IBTS\_tensor which is an array with three dimension: 65 fish species in the first dimension, 31 years in the second dimension, and 7 roundfish areas (RA) in the third dimension. To see the names of the dimension, you can type:

dimnames(IBTS\_tensor)

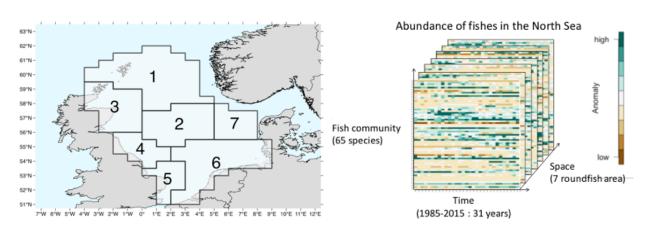


Figure 1:

Abundance data comes from the ICES DAtabase for TRAwl Surveys (DATRAS; http://datras.ices.dk/Home/Default.aspx). The North Sea International Bottom Trawl Survey (NS-IBTS) is an international effort to sampled the demersal fish communities in the North Sea annually and consistently with a standard otter trawl net (chalut Grande Ouverture Verticale, GOV) hauled over the seabed for 30 min. The data is openly available online and the Catch per Unit Effort (CPUE) per length class and per area was downloaded for the first quarter of the period 1985 to 2015 for the roundfish area (RA) 1 to 7. Pre-processing was performed to clean the data, remove the pelagic and the rare species and to transform the CPUE in a three dimensional array, per species, year and roundfish areas (RA).

### Understanding the variables

While loading the data, we saw two different type of variables, quite unusual in R: array and list.

### Array

The object IBTS\_Tensor is an array. Array is generalization of matrix, with more than 2 dimensions. It can only contain numbers. The dimension of the array is given by the function is dim(), and the different elements are accessed with [], similar to a matrix or a data frame:

### dim(IBTS\_tensor)

[1] 65 31 7

```
IBTS_tensor[18,14,6] #Select one element, e.g. Abundance of Cod, in 1998, in RA6
```

### [1] 35.3962

### IBTS\_tensor[18,,6] #Select one vector, e.g. abundance of Cod in 1998

```
1985
               1986
                          1987
                                     1988
                                                1989
                                                           1990
46.253098 91.028164 188.006110 56.636292
                                           44.569500 22.414534
    1991
               1992
                          1993
                                     1994
                                                1995
                                                           1996
66.590250 66.341834 22.999904 31.000150
                                           35.882394 24.778148
    1997
               1998
                          1999
                                     2000
                                                2001
                                                           2002
                                            8.784874 13.173040
42.584734 35.396200 18.648022 12.961334
    2003
               2004
                          2005
                                     2006
                                                2007
                                                           2008
3.980926 10.179750 11.275388
                                6.595208 12.452990 18.515794
    2009
               2010
                          2011
                                     2012
                                                2013
                                                           2014
19.472388 13.342956 14.416228 7.701044
                                            4.403672
                                                      9.886396
     2015
7.990462
```

### IBTS\_tensor[18,,] #Select one matrix, e.g. abundance of Cod

```
    1
    2
    3
    4
    5
    6
    7

    1985
    83.49736
    60.15421
    14.40386
    98.84760
    31.85639
    46.25310
    46.95024

    1986
    68.47722
    48.67987
    31.83467
    100.66903
    32.85298
    91.02816
    93.96932

    1987
    51.54965
    54.80659
    29.67527
    227.12319
    66.39998
    188.00611
    54.00147

    1988
    47.44849
    30.64436
    24.12377
    42.48847
    54.19997
    56.63629
    52.99964

    1989
    67.79382
    55.85585
    39.12256
    102.36360
    29.00079
    44.56950
    65.43840

    1990
    63.41066
    62.33348
    43.45058
    46.30503
    28.29797
    22.41453
    46.94964
```

### List

The names of the dimensions of IBTS\_Tensor are stored in a list. List can contain all kind of elements, without restriction on length or type (can be elements of different lengths made of characters or numbers). The number of elements is given by function is length(), and the different elements are accessed by [[]].

names\_tensor <- dimnames(IBTS\_tensor) # the list of names is stored in a new variable length(names\_tensor) #there are three elements in the list, one element for each dimension

[1] 3

### names\_tensor[[2]] #show the second element of the list

```
[1] "1985" "1986" "1987" "1988" "1989" "1990" "1991" "1992" "1993" "1994" [11] "1995" "1996" "1997" "1998" "1999" "2000" "2001" "2002" "2003" "2004" [21] "2005" "2006" "2007" "2008" "2009" "2010" "2011" "2012" "2013" "2014" [31] "2015"
```

names\_tensor[[1]][18] #show the 18th element of the first element of the list

### [1] "Gadus morhua"

### Your turn:

- 1. What is the index of Hake (Merluccius merluccius) in the dataset?
- 2. What is the abundance of Hake (Merluccius merluccius) in 1988 in RA 1?
- 3. What is the abundance of Hake between 2010 and 2015 in RA 1?
- 4. Can you show the evolution of Hake abundance between 1985 and 2015 in RA 1?

You should get something like:

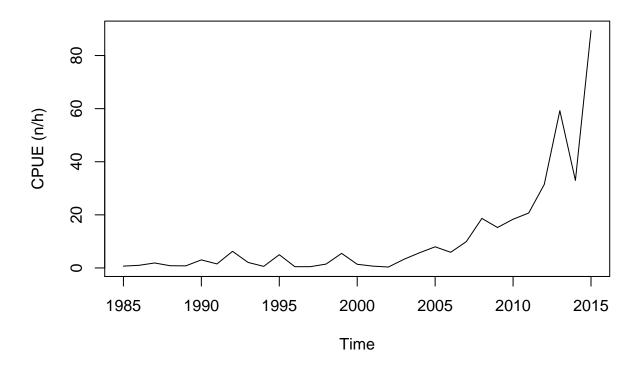
```
[1] 33

[1] 0.83743

2010 2011 2012 2013 2014 2015

18.34881 20.67438 31.44183 59.22724 32.96264 89.40563
```

# Abundance of Hake in RA1



# Solution

# B. Two dimensions: Principal Component Analysis

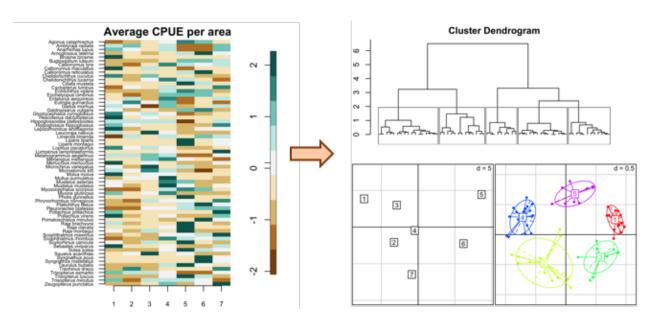


Figure 2:

### Preparing the dataset

### From 3D to 2D

The tensor is flattened into a 2D matrix. This section will study the spatial distribution of fishes in the North Sea. The abundance of fishes will be average over the period 1985-2015, losing the temporal information from the original dataset.

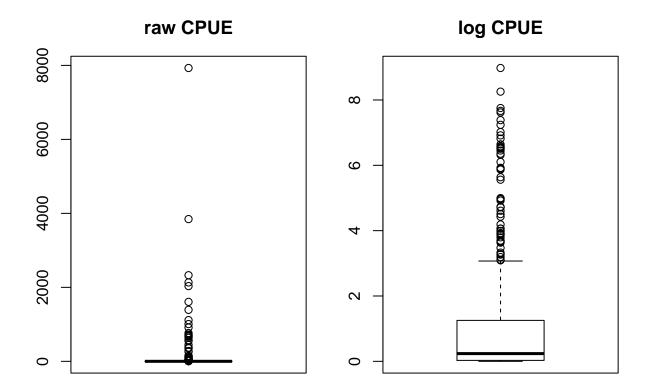
We create a new matrix IBTS\_space which contain the average abundance for the 7 RA (in rows) and 65 species (in columns).

```
IBTS_space <- apply(IBTS_tensor,c(3,1),mean)
dim(IBTS_space) #the new matrix has 7 rows (areas) and 65 columns (species)
[1] 7 65</pre>
```

### Checking the distribution of the data

Principal Component Analysis, and other multivariate analysis in general, are sensible to outliers. So beforehand, one has to be check if the data is not too skewed. If it is the case, it is recommended to log (or square root) transform the data.

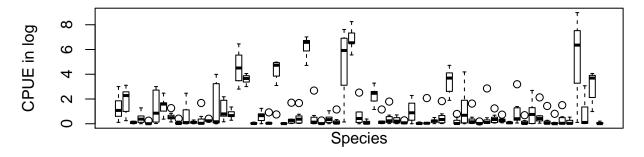
```
#boxplot() is used to look at the distribution
boxplot(as.vector(IBTS_space), main="raw CPUE")
#The CPUE is very skewed, one can not see the difference between the 1st quarter, the median and the 3r
#So data should be log transformed
IBTS_logspace <- log(IBTS_space+1)
#The new distribution of the log transformed CPUE
boxplot(as.vector(IBTS_logspace), main="log CPUE")</pre>
```



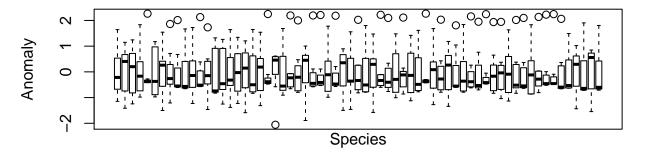
# Scaling the data

The function used to run Principal Components Analysis normalized (i.e. center and scale) the variables by default. It is important to keep that step in mind. An illustration of the normalisation is given in the figures below.

# **Abundance**



# **Anomaly (= abundance normalized per species)**

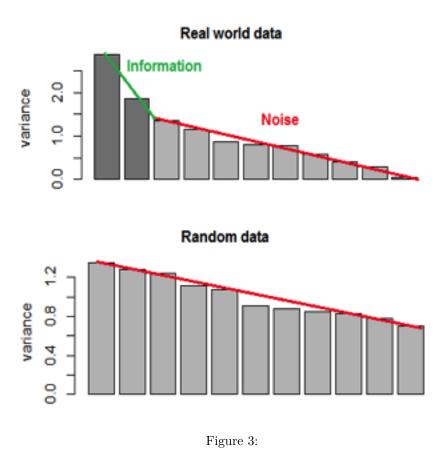


### Principal Component Analysis

### Run a PCA and choose the correct number of PC.

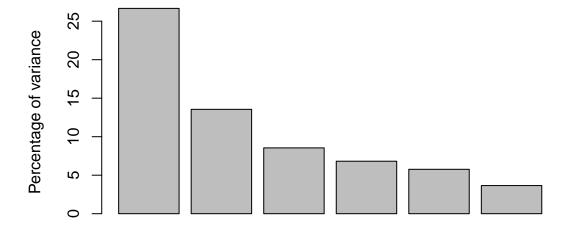
The PCA is run with the function dudi.pca. The data is normalized with the options scale and center set to TRUE. The function is interactive, showing you a scree plot: the variance explained by the successive Principal Components (PC)

The PCA algorithm builds many Principal Component (PC), but not all of the PC perform well in simplifying the information. To choose the correct number of PC to be kept, there are many conflicting methods. The one I recommend is the scree test (Cattell, 1966): graphical detection of a bend in the distribution of the successive variance explained. As a comparison, a PCA run on random data will not find *strong* PC, i.e. PC can not reduce the dimensionality of the data. On the contrary, in real world data, usually we have a bend between successive variance explained by PC. Before the bend, the PCs reduce well the dimensionality of the data and should be kept (i.e. there is a significant pattern); after the bend, PC should be discarded (i.e. it is only noise).



pca\_space=dudi.pca(IBTS\_logspace, scale = TRUE, center = TRUE)

Select the number of axes:



PC

In our case, a bend can be seen on the scree plot after the second PC. So, **please type 2 and press Enter**. The function inertia.dudi show how much variance is explained by the successive PC.

```
#To see how much variance the axes explain:
inertia.dudi(pca_space)$TOT
```

```
inertia cum ratio
1 26.662246 26.66225 0.4101884
2 13.552817 40.21506 0.6186933
3 8.547308 48.76237 0.7501903
4 6.815873 55.57824 0.8550499
5 5.766595 61.34484 0.9437667
6 3.655162 65.00000 1.0000000
```

In our case, the first PC explains 41% of the variance, the two first PC together explain 62% of the variance.

### Interpretation of the PC.

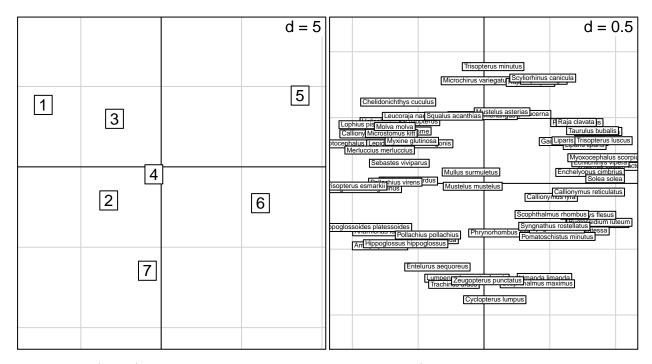
The variable pca\_space, result of the dudi.pca, contains two important matrices: cowith the projetion of columns in the two PC; and li with the projection of the rows in the two PCs. All PC kept in the analysis should be interpreted.

We can see the projection in a table directly by typing the name of these object, or by using graphical function s.label().

### pca\_space\$li # or pca\_space\$co

```
Axis1 Axis2
1 -7.3353689 3.8519039
2 -3.2640990 -2.0883788
3 -2.9030102 2.9885736
4 -0.4560783 -0.4644314
5 8.6404760 4.4337734
6 6.1761023 -2.2590725
7 -0.8580218 -6.4623682
```

```
par(mfrow=c(1,2), mar=c(0,0,0,0))
#Show the weight of the variables:
s.label(pca_space$li, xax=1, yax=2)
s.label(pca_space$co, xax=1, yax=2, clabel = 0.4)
```



The first PC (x-axis) make the difference between the northern NS (RA 1, 2 and 3 have negative weights, projected on the left side) and the southern NS (RA 5 and 6 have positive weights, projected on the right side).

The second PC (y-axis) make the difference between the South-Eastern NS (RA 7 have negative weight, projected on the lower side) and the rest of northern or western NS (RA 1 and 5 have positive weights, projected on the upper side).

The species projection (right side) allow us to see differences between species, but the high number of species makes it difficult to characterize the species. One solution is to group these species, i.e. simplify the number of species into a smaller number of groups, and then characterize these new groups.

### Clustering the species

Clustering is a subject by itself, here we will only see quickly one of its most famous method: Hierarchical clustering. It works in 4 steps:

- 1. Compute the distances between each objects.
- 2. Build a tree according to a given joining criteria
- 3. Choose the number of cluster depending on the topology of the dendogram,
- 4. Create the clusters and interpret them

```
#1. Compute the distance between species
dist_species=dist(pca_space$co, method = "euclidean")

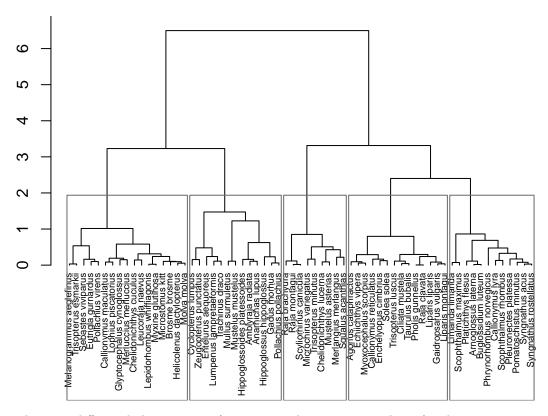
#2. Build a tree with Ward method
den=hclust(dist_species,method = "ward.D2")
```

```
#3. Plot the dendogram
par(mar=c(2,3,3,1))
plot(den, hang=-1, ax = T, ann=T, xlab="", cex=0.6)

#Choosing the number of cluster
nclust<-5

#Visualize the cutting
rect.hclust(den, k=nclust, border="dimgrey")</pre>
```

# **Cluster Dendrogram**

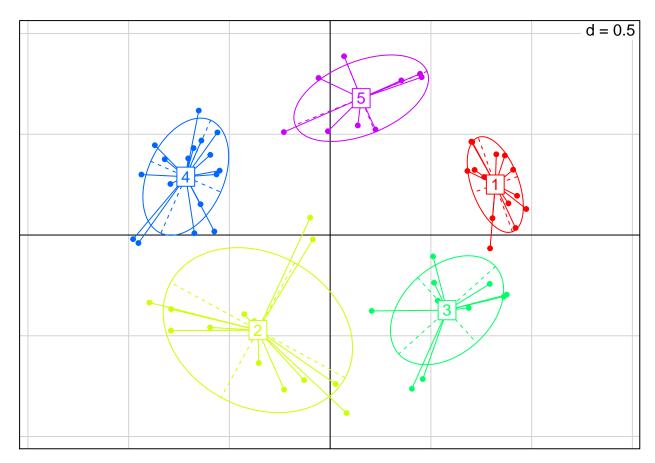


There are different linkage criteria (criterion used to group two objects). The most common ones are: Single (minimum distance between elements of each cluster), Complete (maximum distance between elements of each cluster), Average (mean distance between elements of each cluster, also called UPGMA) and Ward (decrease in variance for the cluster being merged). Ward linkage (Ward, 1963) is known to be more suitable for spherical data, and in most of the case, gives the best results. I invite you to try other linkage criteria by changing the parameter method = in the function hclust. Do you find the same clusters?

In the graph above, the question is where to put a horizontal line on the dendogram to create the clusters. The number of clusters should not be too sensitive of the height of the line. In our case, 5 clusters seem appropriate. We can now visualize how the hierarchical clustering grouped the species in 5 clusters on the two first PC.

```
#4. Create the clusters
clust_space <- as.factor(cutree(den, k=nclust))

#Visualize the cluster in the PC axis
s.class(pca_space$co,fac=clust_space, col=rainbow(nclust),xax=1,yax=2)</pre>
```



These clusters should be interpreted with the previous interpretation of the PC. For example,

- Cluster 1 (in red) has high value in PC1, and above average value in PC2. So it groups species that lives in the southern North Sea (high PC1), with a preference in the south-western side (high PC2).
- Cluster 2 (in yellow in the graph above) groups species mainly located in entrance to the Skagerrak, RA7 (low PC2) but that can spread in the north (low PC1). This cluster is heterogeneous, with the largest ellipse in the figure above.
- Cluster 3 (in green) groups species mainly located in southern NS (high PC1), with a preference on its eastern side (low PC2).

### Your turn:

- 1. How would you interpret the clusters 4 and 5?
- 2. How many species are located mainly in the south-west of the North Sea (RA 5, i.e. grouped in cluster 1)?
- 3. In which cluster is grouped Saithe (Pollachius virens)?
- 4. Depending on your time and your interest, you can either:
  - change the initial matrix and run the same analysis, but categorizing species on their temporal variation over the North Sea (averaged over RA) with the matrix defined as: IBTS\_time <- apply(IBTS tensor,c(2,1),mean)
  - change the linkage method in the clustering algorithm and compare the clusters
  - or go directly to the next step: multivariate analysis in 3 dimension.

### Solution

1. Cluster 4 (in blue) groups species spread exclusively in the northern NS (high PC1 and high PC2). Cluster 5 (in purple) groups species in majority either in the northern extremity (RA1) or the southwestern community (RA5) but not in RA 7 (high PC2)

# C. Three dimension: Tensor Decomposition

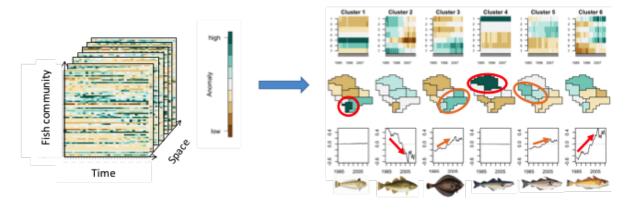


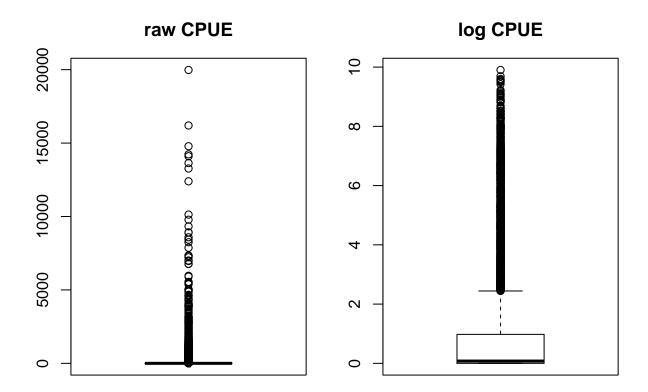
Figure 4:

### Preparing the dataset

### Checking the distribution of the data

Like with PCA, we have to check the skewness of the data and log transform it if it is highly skewed.

```
#boxplot() is used to look at the distribution
boxplot(IBTS_tensor, main="raw CPUE")
#Data should be log transformed
IBTS_logtensor <- log(IBTS_tensor+1)
#The new distribution of the log tranformed CPUE
boxplot(IBTS_logtensor, main="log CPUE")</pre>
```



### Scaling the data

Contrary to PCA, the normalization of a tensor is not straightforward, and have to be done manually before running a PTA. Here, we decided to normalize the values per species and saved them in a new tensor IBTS logscale.

```
#Scaling per species
#Create a new empty array
IBTS_logscale<-array(0,dim=dim(IBTS_tensor))
#Loop scanning each species
for (i in 1:dim(IBTS_tensor)[1]){
    #Calculating the mean and sd of the log CPUE for species i
    ma<-mean(IBTS_logtensor[i,,])
    sa<-sd(IBTS_logtensor[i,,])
    #Saving the anomaly in the array
    IBTS_logscale[i,,]<-(IBTS_logtensor[i,,]-ma)/sa
}
#Copy the labels to the new array
dimnames(IBTS_logscale)<-dimnames(IBTS_tensor)</pre>
```

### Run a PTA and choosing the number of PT.

### Run the PTA

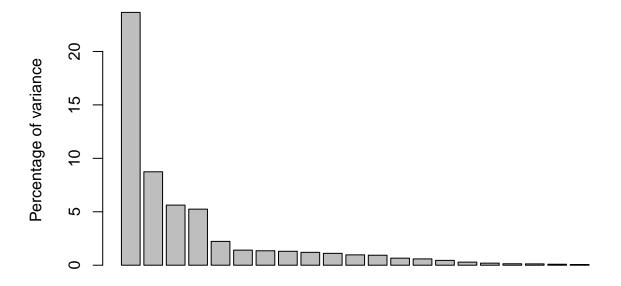
The PTA is run with the function PTA3. The number of principal tensor is indicated by nbPT and nbPT2. One chooses the number of principal tensors at each *level* of analysis by nbPT, the last level (2-modes analysis) is fixed by nbPT2.

The Principal Tensor Analysis computed three main principal tensor and their two mode associated principal tensors.

```
pta<-PTA3(IBTS_logscale, nbPT = 3, nbPT2 = 3, minpct = 0.1)</pre>
 ---Final iteration--- 28
 --Singular Value-- 57.63059 -- Local Percent -- 23.65587 %
 ---Final iteration--- 30
 --Singular Value-- 27.13744 -- Local Percent -- 12.05884 %
 ---Final iteration--- 92
 --Singular Value-- 13.52781 -- Local Percent -- 5.369746 %
 ----Execution Time---- 0.45
summary.PTAk(pta,testvar = 0)
++++ PTA- 3 modes ++++
               data= IBTS_logscale
                                      65 31 7
                -----Percent Rebuilt---- 56.37922 %
                -----Percent Rebuilt from Selected ---- 56.37922 %
              -no- --Sing Val-- --ssX-- --local Pct-- --Global Pct--
                        57.6306 14040.00
vs111
                                               23.6559
                                                            23.655875
                 1
65 vs111 31 7
                 3
                         9.0978
                                 3464.06
                                                2.3894
                                                              0.589533
65 vs111 31 7
                 4
                         5.2054
                                 3464.06
                                                0.7822
                                                              0.192991
31 vs111 65 7
                 6
                        35.0333
                                 6511.03
                                               18.8500
                                                              8.741668
31 vs111 65 7
                7
                                 6511.03
                        28.0998
                                               12.1271
                                                             5.623930
7 vs111 65 31
                9
                        17.6930
                                 4600.42
                                                6.8047
                                                              2.229654
7 vs111 65 31
                10
                        11.4570
                                 4600.42
                                                2.8533
                                                             0.934920
vs222
                11
                        27.1374
                                 6107.06
                                               12.0588
                                                              5.245304
65 vs222 31 7
                                                              0.124215
                13
                         4.1761
                                  789.39
                                                2.2093
65 vs222 31 7
                14
                         3.5684
                                  789.39
                                                1.6131
                                                              0.090693
31 vs222 65 7
                16
                        13.7724
                                 1313.69
                                               14.4387
                                                              1.350994
31 vs222 65 7
                17
                        12.4827
                                 1313.69
                                               11.8612
                                                              1.109821
7 vs222 65 31
                19
                        14.0761
                                 2068.85
                                                9.5771
                                                              1.411229
7 vs222 65 31
                20
                        13.0077
                                 2068.85
                                                8.1785
                                                              1.205129
vs333
                        13.5278
                                 3408.02
                                                5.3697
                                                              1.303431
65 vs333 31 7
                23
                         4.3630
                                  220.69
                                                              0.135581
                                                8.6256
65 vs333 31 7
                24
                         3.0676
                                  220.69
                                                4.2639
                                                              0.067022
31 vs333 65 7
                26
                         7.9591
                                  336.79
                                               18.8094
                                                              0.451196
31 vs333 65 7
                27
                         6.3762
                                  336.79
                                               12.0715
                                                              0.289568
7 vs333 65 31
                29
                        11.6585
                                  996.61
                                               13.6383
                                                              0.968089
7 vs333 65 31
                30
                         9.6144
                                  996.61
                                                9.2751
                                                              0.658381
 ++++
 over 21 PT
```

# Choosing the number of PT.

To select the significant PT, we build the scree plot from the global variance explained.



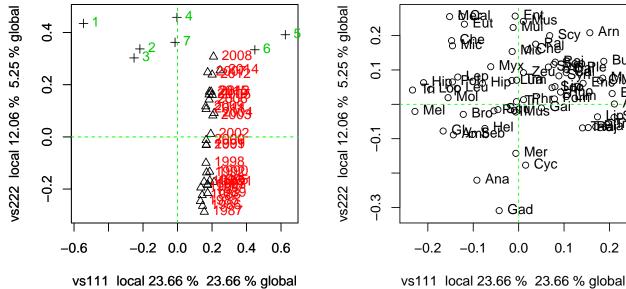
PT

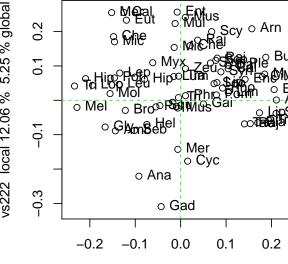
A bend can be seen after 4 PT, so we will select the 4 PT with the best explaining power. Numerically, it corresponds to the PT 1, 6, 7 and 11 of our object pta

## Interpretation of PT.

The plotting function per default allow to use the argument mod to select which dimension to plot, nb1 and nb2 to select which PT will be shown on x-axis and y-axis. For example, we plot the time and space components (the second and third dimension of the array, so mod=c(2,3)) projected on vs111 and vs222 (respectively the elements 1 and 11 of pta, so nb1 = 1, nb2 = 11):

```
par(mfrow=c(1,2))
plot(pta, mod=c(2,3), nb1 = 1, nb2 = 11, xpd=NA, lengthlabels = 4)
plot(pta, mod=1, nb1 = 1, nb2 = 11, lengthlabels = 3)
```





We can see from the plot above that vs111 characterize the space (green crosses representing RA are spread over vs111) and vs222 characterize the time (red triangles representing years are spread over vs222).

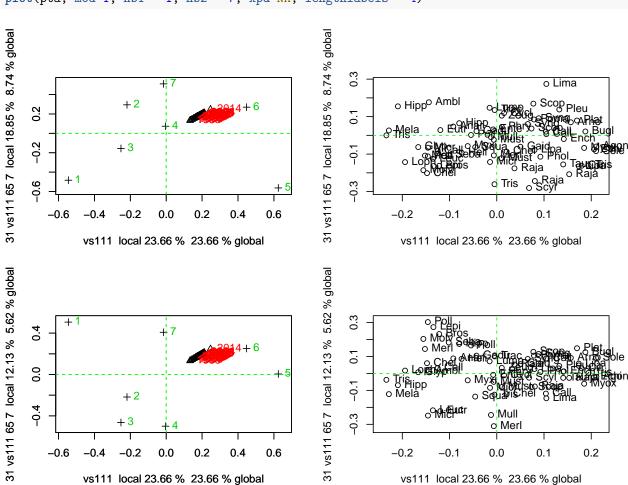
vs111 makes the difference between the northern NS (RA 1, 2 and 3 have negative weights, projected on the left side) and the southern NS (RA 5 and 6 have positive weights, projected on the right side).

vs222 shows a temporal trend and makes the difference between the period 1985-1998 (years before 1998 have negative weights, projected on the lower side) and the recent period 2003-2015 (years after 2003 have positive weights, projected on the upper side).

The species projection (right side, created with mod=1) allow us to see the species plotted on these two PT. For example, we can see Cod (abbreviated *Gad*) in the upper part: it is the species with the strongest decrease in abundance (the highest point in vs222), and doesn't have clear spatial north-south pattern. However, the high number of species makes it difficult to characterize each species (suggesting the need for clustering).

Similar plot can be done for the two other significant PT

```
par(mfrow=c(2,2))
plot(pta, mod=c(2,3), nb1 = 1, nb2 = 6, xpd=NA, lengthlabels = 4)
plot(pta, mod=1, nb1 = 1, nb2 = 6, xpd=NA, lengthlabels = 4)
plot(pta, mod=c(2,3), nb1 = 1, nb2 = 7, xpd=NA, lengthlabels = 4)
plot(pta, mod=1, nb1 = 1, nb2 = 7, xpd=NA, lengthlabels = 4)
```



31vs111 are temporal mode PT associated with vs111, meaning that vs111 and 31vs111 share the same temporal component. This feature can be seen with the straight line representing the projection of the years on these PT.

Another way to represent the spatiotemporal variations of the PT is to use a 2D representation, with x-axis

being the time and y-axis being the space. The graphical functions are out of scope of this tutorial but presented here to help the interpretation.

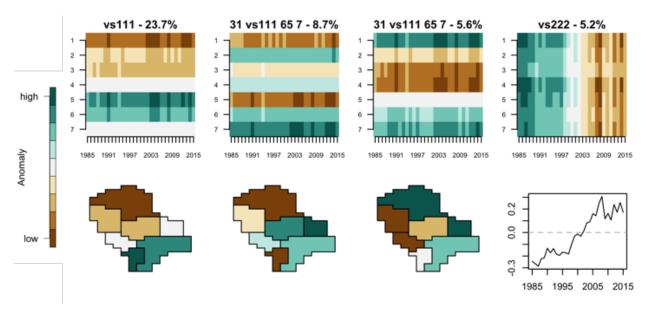


Figure 5:

### Clustering

We use the same approach as before, but with the species projected on the 4 PT.

```
#Create the matrix with the projection of species on the 4 PT
keep <- c(1, 6, 7, 11) # PT that are kept in the analysis
coo<-t(pta[[1]]$v[c(keep),])
labkeep <- paste0(pta[[3]]$vsnam[keep], " - ", round((100 * (pta[[3]]$d[keep])^2)/pta[[3]]$ssX[1],1), "

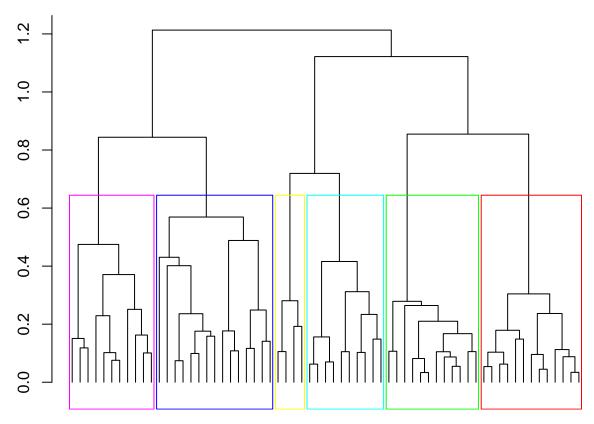
#1. Compute the distance between species
dist1=dist(coo, method = "euclidean")

#2. Build a tree with Ward linkage
den=hclust(dist1,method = "ward.D2")

#3. Plot the dendogram
par(mar=c(1,3,1,1))
plot(den, hang=-1, ax = T, ann=F, xlab="", sub="",labels = FALSE)

#Choose the number of clusters
nclust<-6

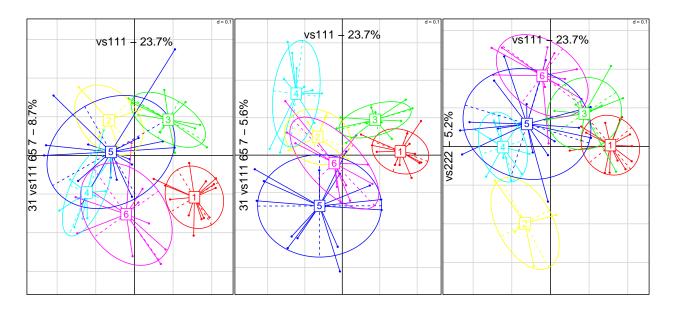
#Visualize the cutting
rect.hclust(den, k=nclust, border=rainbow(nclust)[c(6,5,2,4,3,1)])</pre>
```



The dendogram suggest to create 6 clusters. It is now important to interpret the clusters by projecting them on the PT.

```
#4. Create the clusters
clust_3D <- as.factor(cutree(den, k=nclust))

#Visualize them
par(mfrow=c(1,3))
s.class(coo, fac = clust_3D, xax=1, yax=2, col=rainbow(nclust), clabel = 2)
text(min(coo[,1])-0.03,0, labkeep[2], srt=90, xpd=NA, cex=1.5)
text(0,max(coo[,2])+0.02, labkeep[1], xpd=NA, cex=1.5)
s.class(coo, fac = clust_3D, xax=1, yax=3, col=rainbow(nclust), clabel = 2)
text(min(coo[,1])-0.03,0, labkeep[3], srt=90, xpd=NA, cex=1.5)
text(0,max(coo[,3])+0.02, labkeep[1], xpd=NA, cex=1.5)
s.class(coo, fac = clust_3D, xax=1, yax=4, col=rainbow(nclust), clabel = 2)
text(min(coo[,1])-0.03,0, labkeep[4], srt=90, xpd=NA, cex=1.5)
text(0,max(coo[,4])+0.02, labkeep[1], xpd=NA, cex=1.5)
```



- Cluster 1 is the southern community with high abundance in RA5 and no temporal pattern.
- Cluster 2 is the community with decreasing abundance and no clear spatial pattern.
- Cluster 3 is the South-East community with small increase in abondance
- Cluster 4 is the northern community with high abundance in RA1 and no temporal trend.
- Cluster 5 is the North-West community with a small increase in abondance
- Cluster 6 is the increasing community, mainly pronounced in RA 1, 3 and 5

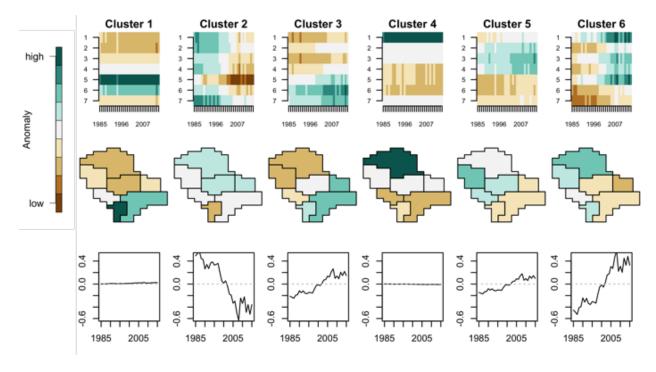


Figure 6:

# Summary

Multivariate analysis are methods to simplify complex and multidimensional dataset. The dimensions are simplified with the objective of keeping most of the variance. In this process, the information and the noise are separated; the main patterns being revealed by the principal components. Multivariate analysis are data-mining tools, in other words they are not predictive or mechanistic tools but data-driven. They can help to visualize and characterize what information is hidden in large dataset.

Steps	2D matrix	3D tensor
1. Check distribution	If too skewed, log or square root transform log(), sqrt()	
2. Scale	Automatically done in dudi.pca()	Depending on each dataset
3. Run the analysis	dudi.pca()	PTA3() or PTAk()
and select # of components with scree plot	Automatically	summary.PTAk()
4. Interpret	s.corcircle() s.label() scatter()	plot(, mod=, nb1= , nb2=)
5. Cluster (if needed)	dist(): compute the distance between individuals hclust(): create the dendogram cutree(): create the clusters s.class(): visualize the clusters on the components	
Package:	ade4	PTA-k

Figure 7:

# References

Cattell, R. B. (1966). The scree test for the number of factors. Multivariate behavioral research, 1(2), 245-276.

Cichocki, A., Mandic, D., De Lathauwer, L., Zhou, G., Zhao, Q., Caiafa, C., & Phan, H. A. (2015). Tensor decompositions for signal processing applications: From two-way to multiway component analysis. IEEE Signal Processing Magazine, 32(2), 145-163.

Leibovici, D. G. (2010). Spatio-temporal multiway decompositions using principal tensor analysis on k-modes: The R package PTAk. Journal of Statistical Software, 34(10), 1-34.

Ward Jr, J. H. (1963). Hierarchical grouping to optimize an objective function. Journal of the American statistical association, 58(301): 236-244.