TOPIC 4 ENVIRONMENTAL DISTANCE

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Similarity/dissimilarity and distance

- Species:
 - sites sharing a similar species composition are ecologically similar
 - i.e. high <u>similarity</u> / low <u>dissimilarity;</u>
 - we will cover dissimilarities in Topic 5.
- Environmental variables:
 - sites sharing similar environmental conditions have a low environmental <u>distance</u> between them;
 - how similar sites to each other depends on...
 - measurable environmental differences that influence species composition,
 - it can be due to unmeasured influences, or
 - it can also simply be noise.
- It is the ecologist's role to figure out what influences the similarity/dissimilarity/distances between sites.

Distance matrices

- A **distance matrix** is produced from a data table (**species table** or **environment table**) by calculating one of several dissimilarity indices.
- Also called **association** or **resemblance** matrices.
- See vegdist () vegan for a list of dissimilarity indices.
- The result is a matrix of **pairwise differences** in community composition (as synthesised by the chosen index) or ecological distance between all sites.

Distance matrix for environmental data

- **Euclidian distance** is "the 'ordinary' straight-line distance between two points in Euclidean space" (i.e. in its simplest form a planar area, which you know of as a graph with x- and y-axes)
- So, in 2D and 3D, gives distance in **cartesian units** between points on a plane (x, y) or in a volume (x, y, z).
- Conforms to our physical concept of distance
 - e.g. short geographic distances between points on a map, and
 - (loses accuracy over large distances, as Earth's surface is not on a plane but on a sphere... correct by using great circle distances, e.g. use the Haversine formula).
- Calculated using the Pythagorean theorem
 - differences are squared, so single large differences become very important, and
 - this is not useful for species data.

Distance matrices

- The matrices are **square** and **symmetrical**, and they will have as many rows and columns as the number of sites present in the original species or environment table.
- The **diagonals are zero** (a site is the same as itself, so it has zero dissimilarity).
- The table can be read directly, and each cell represents the species or ecological difference between a pair of sites.
- All information of the species ID (and maybe also abundance) of a site is lost, as this info is condensed into one metric.

Two dimensions [edit]

In the Euclidean plane, if $\mathbf{p} = (p_1, p_2)$ and $\mathbf{q} = (q_1, q_2)$ then the distance is given by

$$d(\mathbf{p},\mathbf{q})=\sqrt{(q_1-p_1)^2+(q_2-p_2)^2}.$$

This is equivalent to the Pythagorean theorem.

Alternatively, it follows from (*2*) that if the polar coordinates of the point **p** are (r_1, θ_1) and those of **q** are (r_2, θ_2) , then the distance between the points is

$$\sqrt{r_1^2+r_2^2-2r_1r_2\cos(heta_1- heta_2)}.$$

Three dimensions [edit]

In three-dimensional Euclidean space, the distance is

$$d(\mathbf{p},\mathbf{q})=\sqrt{(p_1-q_1)^2+(p_2-q_2)^2+(p_3-q_3)^2}$$

n dimensions [edit]

In general, for an *n*-dimensional space, the distance is

$$d(\mathbf{p},\mathbf{q}) = \sqrt{(p_1-q_1)^2 + (p_2-q_2)^2 + \dots + (p_i-q_i)^2 + \dots + (p_n-q_n)^2}.$$



e.g. example with position (such as geographic) coordinates... use **vegan**'s **vegdist**() function





$$d(a, b) = \sqrt{(a_x - b_x)^2 + (a_y - b_y)^2}$$

e.g. example with 3D position coordinates (a.k.a. dimensions)...



7 5.385165 5.656854 5.916080 9.433981 7.874008 4.582576

$$d(a, b) = \sqrt{(a_x - b_x)^2 + (a_y - b_y)^2 + (a_z - b_z)^2}$$

e.g. example with environmental 'dimensions'...

Raw data

| site 🗦 | temperature 🗦 | depth 🗘 | light 🗦 |
|--------|---------------|---------|---------|
| а | 4 | 1 | 3 |
| b | 5 | 5 | 5 |
| с | 6 | 6 | 4 |
| d | 1 | 4 | 9 |
| е | 2 | 3 | 8 |
| f | 8 | 3 | 1 |
| g | 9 | 1 | 5 |

Euclidian distances

| R> | <pre>ex.xyz.euc</pre> | | | | | | |
|--------------|-----------------------|----------|----------|-----------|----------|----------|--|
| R> | ex.xyz.euc | | | | | | |
| | 1 | 2 | 3 | 4 | 5 | 6 | |
| 2 | 4.582576 | | | | | | |
| 3 | 5.477226 | 1.732051 | | | | | |
| 4 | 7.348469 | 5.744563 | 7.348469 | | | | |
| 5 | 5.744563 | 4.690416 | 6.403124 | 1.732051 | | | |
| 6 | 4.898979 | 5.385165 | 4.690416 | 10.677078 | 9.219544 | | |
| 7 | 5.385165 | 5.656854 | 5.916080 | 9.433981 | 7.874008 | 4.582576 | |

 $d(a, b) = \sqrt{(a_{\text{temp}} - b_{\text{temp}})^2 + (a_{\text{depth}} - b_{\text{depth}})^2 + (a_{\text{light}} - b_{\text{light}})^2}$

e.g. example with higher dimension environmental data...

Raw data

| ^ | pH 🗦 | O2 [‡] | temp 🍦 | depth 🗦 |
|----------|------|------------------------|--------|---------|
| a | 7.1 | 6.5 | 12.1 | 1.1 |
| b | 7.5 | 5.5 | 12.3 | 1.3 |
| c | 7.6 | 5.7 | 11.9 | 1.5 |
| d | 7.0 | 5.4 | 11.8 | 1.6 |
| e | 7.1 | 6.3 | 12.0 | 1.8 |
| f | 7.2 | 6.3 | 12.1 | 1.9 |
| g | 6.9 | 6.1 | 12.2 | 2.2 |

Standarised data

(transformation)

| > | ex.env.std | ← decosta | nd(xy.env, r | method = "sta | ndardize") |
|---|------------|------------|--------------|---------------|------------|
| > | ex.env.std | | | | |
| | рН | 02 | temp | depth | |
| а | -0.3872983 | 1.2156767 | 0.2494233 | -1.41749621 | |
| b | 1.1618950 | -1.0842522 | 1.4133987 | -0.88114629 | |
| С | 1.5491933 | -0.6242664 | -0.9145521 | -0.34479637 | |
| d | -0.7745967 | -1.3142450 | -1.4965398 | -0.07662142 | |
| е | -0.3872983 | 0.7556909 | -0.3325644 | 0.45972850 | |
| f | 0.0000000 | 0.7556909 | 0.2494233 | 0.72790346 | |
| g | -1.1618950 | 0.2957051 | 0.8314110 | 1.53242833 | |

Euclidian distances

| > | ex | .env.eu | uc ← veg | ex.env dist(xy.d a | .sta at, method | d = "eucl: | idian") |
|---|----|---------|----------|-----------------------------------|--------------------|------------|----------|
| > | ex | .env.eu | uc | | | | |
| | | a | b | С | d | е | f |
| b | 4. | 123106 | | | | | |
| С | 6. | 324555 | 2.236068 | | | | |
| d | 4. | 242641 | 4.123106 | 5.830952 | | | |
| е | 2. | 828427 | 3.605551 | 5.656854 | 1.414214 | | |
| f | 4. | 472136 | 3.605551 | 4.472136 | 7.071068 | 6.000000 | |
| g | 5. | 000000 | 5.656854 | 6.708204 | 8.544004 | 7.280110 | 2.236068 |