SuperTRI3

A) Short description of the method

The SuperTRI method is based on the branch support analyses of the independent data sets, in which the reliability of the nodes is assessed using three measures: the supertree Bootstrap percentage (SBP) and two other values calculated from the separate analyses: the mean branch support (mean Bootstrap percentage [MBP] or mean posterior probability [MPP]) and the reproducibility index (Ropiquet et al., 2009). The SuperTRI approach shows less sensitivity to the phylogenetic methods and models, and it is more accurate to interpret the relationships among taxa.

supertri3.py

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supertri3.py is a python script that will help you producing a matrix representation weighted by bootstrap proportions or bayesian posterior probabilities. Such a matrix representation can be used to produce supertrees.

supertri3.py also computes the support indices described in Ropiquet et al. (2009): reproducibility and mean robustness (or mean bayesian posterior probability) and may report these indices on nexus trees.

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supertri3.py is a python script. To be able to run it, you must first check wether python is installed.

In MacOSX there is a command-line terminal called "Terminal"; it should be in the "Utilities" directory of the "Applications" directory.

B) Choose a place where to install supertri.py

It is advisable to put files in a place where you you can find it later. There are usually some conventions on where to put executable files. For example, on an UNIX-like operating system (UNIX, Linux, MacOSX), executables are often placed in directories ending in "bin" (bin for "binary": that is, programs compiled in machine language, note that supertri3.py is not in machine language; you may open this python script in a text editor and see human-readable commands, python is the program that will translate these commands into machine language).

There are system-wide bin directories, like /usr/local/bin, that are usually in the PATH of every user (this means that programs in such a place will be found and executed when the user types the name of the program in a command-line terminal). To place a program in /usr/local/bin you need to have administrative rights. Under MacOSX, it is possible that directories like /usr/local/bin are not visible from the finder, but only from the command-line terminal. In such a case, you will have to use the command "cp" to copy supertri3.py to the place you want.

If you are not allowed to place a program in /usr/local/bin, you may place the program in a personal bin directory. A good place would be a bin directory in your home directory. For example, if your username is dupont and you are in a MacOSX system, this would be /Users/dupont/bin. On a Linux system, this would be /home/dupont/bin.

C) There are several ways to execute a python script

Once you have placed supertri3.py in a suitable place, if you want to be able to call the command directly, you'll have to check wether that place is in your PATH or not (see A) 2)). You may also call python and tell it to run the supertri3.py script.

1) The directory containing supertri3.py is in your PATH

This is probably the most convenient situation.

Open a command-line terminal.

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Go to the directory where you placed supertri.py, using the command "cd" ("change directory"):
---- MacOSX-like example -----
Welcome to Darwin !
Computer:~ dupont$ cd /Users/dupont/bin
Computer:~/bin dupont$
-----
Make the script executable, using the command "chmod":
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Computer:~/bin dupont\$ chmod +x supertri3.py

Computer:~/bin dupont\$

```
Now you can call supertri3.py in command-line from any directory (provided your
PATH contains the directory where supertri3.py has been placed):
Computer: ~/bin dupont$ cd
Computer: ~ dupont$ supertri3.py
usage: supertri3.py [options] <arguments>
available options:
[list of options here]
In interactive mode, the user will be prompted for the arguments,
otherwise, the arguments should be given in the command line.
Traceback (most recent call last):
  File "/Users/dupont/bin/supertri.py", line 995, in ?
    main()
  File "/Users/dupont/bin/supertri.py", line 788, in main
    raise Exception("No dataset names provided. Either use option -i or provide
a file with option --marklist.")
Exception: No dataset names provided. Either use option -i or provide a file
with option --marklist.
Well, it seems to work. Go to D).
2) You want to call python explicitly
Open a command-line terminal.
Go to the directory where you placed supertri.py, using the command "cd":
---- MacOSX-like example ----
Welcome to Darwin !
Computer:~ dupont$ cd /Users/dupont/bin
Computer:~/bin dupont$
Tell python to execute the script:
Computer:~/bin dupont$ python supertri3.py
usage: supertri3.py [options] <arguments>
available options:
[list of options here]
In interactive mode, the user will be prompted for the arguments,
otherwise, the arguments should be given in the command line.
Traceback (most recent call last):
  File "supertri3.py", line 995, in ?
    main()
  File "supertri3.py", line 788, in main
    raise Exception("No dataset names provided. Either use option -i or provide
a file with option --marklist.")
Exception: No dataset names provided. Either use option -i or provide a file
with option --marklist.
____
```

3) You want to execute the script from anywhere, but it is not in your PATH Open a command-line terminal. Go to the directory where you placed supertri3.py, using the command "cd": ---- MacOSX-like example ----Welcome to Darwin ! Computer:~ dupont\$ cd /Users/dupont/bin Computer:~/bin dupont\$ Make the script executable, using the command "chmod": Computer:~/bin dupont\$ chmod +x supertri3.py Computer: ~/bin dupont\$ Go to the directory where you placed your analyses results, using the command "cd": Computer: ~/bin dupont\$ cd /Users/dupont/Documents/Analyses/Smurfs/bootstrap Computer:~/Documents/Analyses/Smurfs/bootstrap dupont\$ Execute the script, calling it with its absolute path: Computer:~/Documents/Analyses/Smurfs/bootstrap dupont\$ /Users/dupont/bin/supertri3.py usage: supertri3.py [options] <arguments> available options: [list of options here] In interactive mode, the user will be prompted for the arguments, otherwise, the arguments should be given in the command line. Traceback (most recent call last): File "/Users/dupont/bin/supertri.py", line 995, in ? main() File "/Users/dupont/bin/supertri.py", line 788, in main raise Exception("No dataset names provided. Either use option -i or provide a file with option --marklist.") Exception: No dataset names provided. Either use option -i or provide a file with option --marklist.

D) Input files

A preliminary piece of advice:

When working with command-line tools, it is usually better not to have files with names containing spaces, accentuated letters or special characters other than underscore ("_"). Spaces are used to separate commands and arguments. Special characters may have special meanings. Accentuated letters are not always handled correctly.

To produce a valid nexus matrix representation of your results, weighted by support values, supertri3.py needs at least the following:

- files describing the bipartitions with "." and "*" and the support values, produced by the analyses of your different datasets.
- a file containing the taxon names, in the same order as in the matrix used for the analyses, without blanks in the names, and with a name on each line of the file.
- if there are taxa missing from one or more datasets, prepare for each dataset a file containing the names of the taxa missing for this dataset.

The results files should be named the following way: Dataset.parts for the results of a MrBayes analysis, where Dataset is the name of the dataset that was analysed.

Dataset.log for the log of a PAUP bootstrap analysis, where Dataset is the name of the dataset that was analysed.

Since the description of bipartitions with "." and "*" is made without taxon names, the analyses should all have been done whith the taxa in the same order in the matrix. Otherwise, the representations of the bipartitions from two different analyses will not be comparable. This order is given to supertri.py by the file with the names of the taxa.

The files for the missing taxa should contain the taxa names as they appear in the list of the taxa. These files should be named the following way:
Dataset.abs, where Dataset is the name of the dataset lacking the taxa in the file.

It is recommended, but optional to prepare a file with the names of the datasets, without blanks in the names of the datasets, and with one name per line. These names should correspond to the prefixes of the files containing the results.

It is advisable not to have other files than those concerned with the analyses in the directory where you put the data for supertri3.py.

E) Executing supertri3.py

As many command-line tools, supertri3.py works with a system of options.

Suppose the names of the taxa are written in a file named "taxons". To indicate supertri3.py to use this file, the command will be called with the --taxlist option:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons

This is not enough to make it work. If you want to provide a list of datasets on the command line, with a file named "genes", you will have to use the --datasets option:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons --datasets genes

You may also run the program in interactive mode, with the -i option, so that you can provide the names of the datasets interactively:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons -i

You may also specify the type of results files with the --suffix option:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons --datasets genes -- suffix .log

You can chose the name of the matrix to write with the option -o:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons --datasets genes -- suffix .log -o MRP.nex

The default name will be composed by the names of the datasets followed by ".mrpp.nex" for MrBayes data or ".mrbp.nex" for bootstrap log data. You can chose the root of the output trees with the option --root:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons --datasets genes --suffix .log -o MRP.nex --root The_outgroup_taxon

You must ensure that the name you give is present in the file listing the taxa, but not in any .abs file.

You can give trees on which to map the indices with the option --intree:

Computer:~ dupont\$ python3 supertri3.py --taxlist taxons --datasets genes --suffix .log -o MRP.nex --root The_outgroup_taxon --intree synthesistree.tre

The trees you give with this option must be in nexus format and contain the names of the taxa in the parenthesized representation of the tree; not in a translation block. You can provide several trees in the same file, or call the option several times with different file names.

If you want to map the indices on a tree obtained from the matrix representation produced by a first run of supertri.py, you'll have to run supertri3.py again, with the --intree option, with the previously used parameters.

The indices will be written as node labels in two separate nexus trees. They will also be written in a .tgf file for graphical export of .svg or .eps trees. In the .tgf file, the first index above the branch is the mean support and the second is the reproducibility index.

F) Miscellaneous remarks

This program was not written by a professional programmer. We hope it works and we hope the code is clear enough to allow you to modify it if you know python.

If python works and you are able to launch supertri3.py but you still have problems running this program, check carefully that your files conform to the requirements of part D).

There may also be problems due to file end-of-line coding.

Before reporting bugs, note the error messages, if any, and try to also determine the version of python and the version of supertri3.py that you are using. Please also indicate the type of operating system you are using (Windows, Mac, Linux, with details about the particular version of the operating system).

G) References

Citation

Ropiquet A., Li B. & Hassanin A. (2009) SuperTRI: a new approach based on branch support analyses of multiple independent data sets for assessing reliability of phylogenetic inferences. Comptes Rendus Biologies 332: 832—847.

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