TreeCmp 2.0: comparison of trees in polynomial time – manual

# Introduction

A phylogenetic tree represents historical evolutionary relationship between different species or organisms. There are various methods for reconstructing phylogenetic trees. Applying those techniques usually results in different trees for the same input data. An important problem is to determine how distant two trees reconstructed in such a way are from each other. Comparing phylogenetic trees is also useful in mining phylogenetic information databases. The TreeCmp application was designed to compute distances between unweighted arbitrary (not necessary binary) phylogenetic trees. All metrics are implemented using polynomial time algorithms. From the simplest but most efficient Robinson-Foulds metrics. Working in linear time complexity O(|L|), where L - set of taxa. To slower but more sophisticated matching metrics (Matching Split/Cluster/Pair) of polynomial time complexity O(|L|2.5\*log|L|).

# Metrics

The TreeCmp package uses 7 polynomial implementations of metrics for rooted trees:

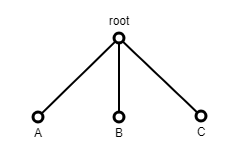
* **Triples metric** (|L|2) is based on comparing sets of triples induced by two trees. A set of triples induced by a rooted tree is a set of the topologies of all 3-species rooted subtrees consistent with this tree. Triplets distance between two trees is the number of different triples in the respective sets.
* **Robinson-Foulds metric** based on clusters (|L|) is one of the most commonly used methods to measure the similarity of phylogenetic trees. This metric is equal to the number of different clusters in compared trees (divided by 2).
* **Matching Pair metric** (|L|2.5\*log|L) uses the concept of the minimum-weight perfect matching in a complete bipartite graph constructed from partitions of all pairs of leaves of the compared phylogenetic trees. This metric becomes the transfer distance between partitions of the set of unordered leaf pairs determined by compared trees.
* **Nodal Splitted metric with L2 norm** (|L|2) is L2 norm of two splitted path lengths matrices with rows and columns indexed by taxa and where every entry is the distance from the least common ancestor of the pair of nodes.
* **Matching Cluster metric** (|L|2.5\*log|L) is based on comparing clusters in two trees. A cluster associated with a vertex v in a rooted tree T with leaves L is a subset of leaves that are descendants of v. To measure the dissimilarity between clusters, MC uses function hC(A, B) = |A⊕B|. For a dummy element, O = ∅, hC(A, O) = |A|.
* **Rooted maximum agreement subtree distance** (|L|2) – Given a set A of n species and two rooted trees leaves uniquely labeled by the items of A, Maximum Agreement Subtree distance is the difference between cardinality of A and cardinality of the largest subset B of A such that the subtrees of compared trees induced by B are isomorphic.
* **Cophenetic Metric with L2 norm** (|L|2) – For every rooted phylogenetic tree T, let its cophenetic vector φ(T) consist of all pairs of cophenetic values (depth of lowest common ancestor of a pair of nodes) between pairs of taxa in T and all depths of taxa in T. Cophenetic Metric comparing cophenetic vectors φ(T) of two trees by means of L2 norm.

and 5 polynomial implementations of metrics for unrooted trees:

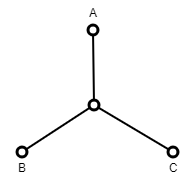
* **Quartet distance** (|L|3) is based on comparing sets of quartets induced by two trees. A set of quartets induced by an unrooted tree is the set of the topologies of all 4-species subsets of its leaves consistent with its topology. Quartet distance between two trees is the number of different quartets in two respective sets.
* **Path difference distance** (|L|2) is the square root of the sum of squares of the differences in the distance between individual pairs of leaves in the compared trees.
* **Robinson-Foulds distance** (|L|) is one of the most commonly used methods to measure the similarity of phylogenetic trees. This metric is equal to the number of different splits in compared trees (divided by 2).
* **Matching Split distance** (|L|2.5\*log|L) is based on comparing splits in two trees. A split A|B of a set L is an unordered pair (ie, A|B = B|A) of its subsets, such that L = A∪B and A∩B = ∅. To compare splits in two trees, MS finds a minimum-weight perfect matching in bipartite graphs whose vertices correspond to splits in these two trees and edges connect each split from one tree to a split in another tree. If the number of splits in the trees differs, the smaller set is extended by the missing number of "dummy" elements.
* **Unrooted maximum agreement subtree distance** (|L|(2+o(1))) – Given a set of species L and two unrooted trees leaves uniquely labeled by the items of L, Unrooted Maximum Agreement Subtree distance is the difference between cardinality of L and cardinality of the largest subset A of L such that the subtrees of compared trees induced by A are isomorphic.

# Input data format

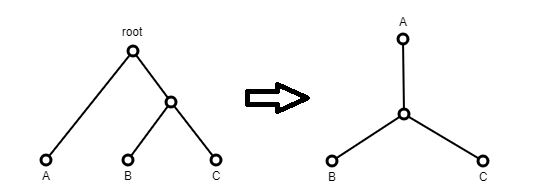
The TreeCmp software was designed to support BEAST (<http://beast.bio.ed.ac.uk/>) and MrBayes (<http://mrbayes.csit.fsu.edu/>) date files, where phylogenetic trees are stored in the NEWICK format. Note that plain text files containing only trees in this format are supported as well. The NEWICK format clearly specifies the vertex that is a candidate for the root. After choosing a metric for rooted trees, this vertex will always be treated as the root. For example, if a unrooted binary tree in the NEWICK format: (A, B, C) is entered in the metric dedicated for a rooted tree, it will be interpreted as a rooted, non-binary tree consisting of a 3-degree root and 3 descendant vertices A, B and C.



In the case when a rooted binary tree is entered to the metric dedicated for a unrooted tree, the root will be treated as an internal vertex or will be automatically shrinked if it’s degree equals 2. For example, if a rooted binary tree in the NEWICK format (A, B, C) is entered in the metric for a unrooted tree, then root will be treated as an internal vertex and tree will be interpreted as a unrooted, not-binary tree (A, B, C).



However, after entering the rooted tree (A, (B, C)) to the metric dedicated for unrooted trees, the root will be shrinked as in the figure below.



In summary, any interference in the given rooted tree will take place only if calculated metric is dedicated to unrooted trees and the root degree is 2. In any other case it will be treated as an internal vertex.

# Compilation

## Compilation from the command line

To compile via the command-line, creating files in the directory out/class, you should be able to do the following from the top level folder

javac -d out/class -cp lib/commons-cli-1.4.jar src/treecmp/\*.java src/treecmp/\*/\*.java src/pal/\*/\*.java

The resulting compiled files can be run directly, for example by issuing the command:

java -cp out/class:lib/commons-cli-1.4.jar treecmp.Main

You will need to replace the colon with a semicolon on windows systems. However, it is usually easier to create a stand-alone .jar executable as below.

## Creating a jar executable from the command line

The jar executable (e.g. TreeCmp.jar) should be created in the bin directory, as it expects to find the config/config.xml file one level above the directory in which the executable resides. Once the class files have been compiled in the out/class directory as above, this jar file can be created in the correct place using the MANIFEST.MF file in src/META-INF:

jar cvfm bin/TreeCmp.jar src/META-INF/MANIFEST.MF -C out/class/ .

As defined in the manifest, extra libraries are expected to be placed in lib directory in the same place as the jar file. An easy way to do this is to move or copy the lib folder (containing jar libraries) into the bin directory. For example, on unix-like systems you could do cp -a lib bin/, on windows xcopy lib bin\.

# Running TreeCmp

The TreeCmp application is distributed as a zip archive. In order to unpack the file any software supporting zip compression, for example free software 7-zip (<http://www.7-zip.org/>), can be used. In order to run the TreeCmp application Java VM in version at least 1.6 is required.

## Directory structure

|  |  |  |
| --- | --- | --- |
| **Description** | | |
| bin |  | contains main jar file: **TreeCmp.jar** and lib folder with necessary open source libraries: pal-1.5.1 (<http://www.cebl.auckland.ac.nz/pal-project/>) and commons-cli-1.2 (<http://commons.apache.org/cli/>) |
| config |  | contains xml configuration file |
| data |  | contains text files with pre-computed data (average value and other statistics) for all the 12 metrics under the two models of generation of random binary trees: the Yule model and the uniform model. |
| examples |  | contains subdirectories with examples |
|  | align | contains an example of creating alignments |
|  | beast | contains an example input file created using BEAST |
|  | mr\_bayes | contains an example input file created using MrBayes |
|  | plain | contains an example input file with plain trees |
|  | plain2 | contains an example input file with plain trees |
|  | prune | contains an example of comparing trees having different sets of taxa |
|  | ref\_tree | contains an example of comparing reference trees to a set of trees |
|  | scaled | contains an example with reporting scaled values of chosen metrics |
| src |  | contains source code of this application |

## Command line syntax

Usage:

java -jar TreeCmp.jar -w <size>|-s|-m|-r <refTreeFile> –d <metrics> -i <inputfile> -o <outputfile> [-N] [-P] [-I] [-A|-O]

Note that options order is important. See section 5 for details regarding output file format for a particular combination of the options.

**Mandatory switches:**

* The comparison mode options (only one option should be specified):
  + –s – overlapping pair comparison mode; every two neighboring trees in the input file are compared,
  + -w <size> – window comparison mode; every two trees within a window with a specified size are compared – the average distance and the standard deviation go to the output file,
  + –m – matrix comparison mode; every two trees in the input file are compared.
  + -r <refTreeFile> – reference trees to all trees mode. Each tree in the input file is compared to all reference trees.

Details of the computation flow in each of these case are explained in the pictures below.





* The metric option (-d). At least one and at most 12 metrics can be specified (numbers in square brackets correspond to the reference list. Metrics should be separated by space character.

Metrics for rooted trees:

* + tt – the Triples metric (Crichlow et al. 1996),
  + rc – the Robinson-Foulds metric based on clusters (Robinson and Foulds 1981),
  + mp – the Matching Pair metric (Bogdanowicz and Giaro 2014),
  + ns – the Nodal Splitted metric with L2 norm (Cardona et al. 2010),
  + mc – the Matching Cluster metric (Bogdanowicz et al. 2012),
  + mt – the Rooted maximum agreement subtree distance (Farach and Thorup 1994),
  + co – the Cophenetic Metric with L2 norm (Cardona, Mir, Rosselló, Rotger and Sánchez 2013).

Metrics for unrooted trees:

* + qt – the Quartet distance (Estabrook 1985),
  + pd – the Path difference distance (Steel and Penny 1993),
  + rf – the Robinson-Foulds distance (Robinson and Foulds 1981),
  + ms – the Matching Split distance (Bogdanowicz and Giaro 2012),
  + um – the Unrooted maximum agreement subtree distance (Farach and Thorup 1994).

Example: -d ms rf

* IO options (both options should be specified):
* -i <inputfile> – input data file with trees in the NEWICK format,
* -o <outputfile> – output data file with the results of computations.

**Optional switches:**

* General options:
  + –N – report normalized distances δ*m* for a particular metric *m* (Bogdanowicz et al. 2012; based on an average value from pre-computed data). This functionality is available for trees with number of leaves between 4 and 1000. Note that normalized tree similarity for a particular metric *m* (*NTSm*) can be expressed by normalized distance as follows: *NTSm*. = 1 - δ*m* (Bogdanowicz et al. 2012).
  + –P – prune compared trees if needed. This option is design to allow comparing trees having different (partially overlapping) sets of taxa. After using this option three additional columns appear in the output file (see section 4 for details).
  + –I – -include summary section in the output file.
* Matching metric specific options (only one option should be specified).
  + –A – Generate alignment files – this option should be used together with selection the MS or MC metrics. As a result additional files containing aligned splits or clusters are generated:
    - [output\_file\_name].out.aln\_MS.txt,
    - [output\_file\_name].out.aln\_MC.txt,

where [output\_file\_name] is the file name specified after -o option.

* + -O – use special implementations of MS/MC metrics optimized for similar trees.

Note that if a rooted tree (with bifurcation in the root) is compared using metrics for unrooted trees the tree will be automatically transform into unrooted one, i.e., the bifurcation will be replaced with an arbitrary trifurcation.

# Output data format

All output files created by the application regardless of chosen mode have similar structure. Output files are tab separated text files (TSV), which means that they can be easily read by various data analysis software (e.g. MS Excel, R, OpenOffice.org). An output file consists of two sections. The first section contains formatted in rows values of distances in selected metrics. The second (optional) section contains summary data computed based on all rows that appears in the first section.

## Basic output file structure

Base output file format for options -s, -m, and -w

|  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- |
| **No** | **Tree1** | **Tree2** | **MetricName\_1** | **MetricName\_2** | **…** | **MetricName\_n** |
| Comparison number | Tree1 number | Tree2 number | Distance value | Distance value | … | Distance value |

Base output file format for option -r,

|  |  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- | --- |
| **No** | **RefTree** | **Tree** | **MetricName\_1** | **MetricName\_2** | **…** | **MetricName\_n** |
| Comparison number | Reference tree number | Tree number | Distance value | Distance value | … | Distance value |

Tree, tree1, tree2 numbers in the output file correspond to the number of the tree in the input file.

The following table contains a mapping between available metrics and column names in the output file that are related to them.

|  |  |  |
| --- | --- | --- |
| **Metric name in the output file** | **Full metric name** | **TreeCmp command line parameter** |
| Triples | the Triples metric | tt |
| R-F\_Cluster (0.5) | the Robinson-Foulds metric based on clusters | rc |
| MatchingPair | the Matching Pair metric | mp |
| NodalSplitted | the Nodal Splitted metric with L2 norm | ns |
| MatchingCluster | the Matching Cluster metric | mc |
| MAST | the Rooted maximum agreement subtree distance | mt |
| CopheneticL2Metric | the Cophenetic Metric with L2 norm | co |
| Quartet | the Quartet distance | qt |
| PathDiffernce | the Path difference distance | pd |
| R-F(0.5) | the Robinson-Foulds distance | rf |
| MatchingSplit | the Matching Split distance | ms |
| UMAST | the Unrooted maximum agreement subtree distance | um |

## Additional columns (-P and -N options)

After using switch -P the following three columns appear additionally in the output file.

|  |  |  |
| --- | --- | --- |
| **Tree1\_taxa** | **Tree2\_taxa (or RefTree\_taxa)** | **Common\_taxa** |
| Number of taxa in the first tree | Number of taxa in the second (or reference) tree | Number of taxa in common |

After using switch -N the following two columns per each chosen metric appear additionally in the output file. These columns contain the value of the distance in a particular metric divided by its empirical average value. If the number of common leaves in compared trees is out of supported range (which is form 4 to 1000), then “N/A” value is inserted.

|  |  |
| --- | --- |
| **MetricName\_toYuleAvg** | **MetricName\_toUnifAvg** |
| (Distance value)/(Empirical average value in the Yule model) | (Distance value)/(Empirical average value in the uniform model) |

For details regarding generating phylogenetic trees under the Yule and uniform models see (McKenzie and Steel 2000; Semple and Steel 2003).

## Summary section format (-I option)

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
| **Name** | **Avg** | **Std** | **Min** | **Max** | **Count** |
| Metric name 1 | Average value | Standard deviation value | Minimal value | Maximal value | Number of analyzed values |
| Metric name 2 | … | … | … | … | … |
| … | … | … | … | … | … |
| Metric name n | … | … | … | … | … |

# Useful Java VM parameters

In the case of an analysis of large trees the following exceptions might occur:

1. Exception in thread "main" java.lang.OutOfMemoryError: Java heap space

To solve the problem increase Java heap space memory limit using JVM option –Xmx

Example:

java –Xmx700m –jar TreeCmp.jar <further options>

1. Exception in thread "main" java.lang.StackOverflowError at pal.io.FormattedInput.skipWhiteSpace(FormattedInput.java:111)

at pal.io.FormattedInput.readNextChar(FormattedInput.java:131)

at pal.tree.ReadTree.readNH(ReadTree.java:81)

…..

at pal.tree.ReadTree.readNH(ReadTree.java:89)

To solve the problem increase Java thread stack size limit using JVM option –Xss

Example:

java –Xss1m –jar TreeCmp.jar <further options>

These options can be used in conjunction.

# Examples

## Running application to compare trees using MS

Input file: \examples\beast\testBSP.newick

Invocation:

java -jar TreeCmp.jar -w 2 -d ms -i testBSP.newick -o testBSP.newick\_w\_2.out -I

Console output:

TreeCmp version 1.0-b291

Active options:

Type of the analysis: window comparison mode (-w) with window size: 2

Metrics:

1. MatchingSplit (ms)

Input file: testBSP.newick

Output file: testBSP.newick\_w\_2.out

Additional options:

I - Include summary section in the output file.

-----

2011-08-27 16:03:17: Start of scanning input file: testBSP.newick

2011-08-27 16:03:17: End of scanning input file: testBSP.newick

2011-08-27 16:03:17: 11 valid trees found in file: testBSP.newick

2011-08-27 16:03:17: Start of calculation...please wait...

2011-08-27 16:03:17: 0.00% completed...

2011-08-27 16:03:17: 20.00% completed...

2011-08-27 16:03:17: 40.00% completed...

2011-08-27 16:03:17: 60.00% completed...

2011-08-27 16:03:17: 80.00% completed...

2011-08-27 16:03:17: 100.00% completed.

2011-08-27 16:03:17: End of calculation.

2011-08-27 16:03:17: Total calculation time: 62 ms.

Output file testBSP.newick\_w\_2.out:

No Tree1 Tree2 MatchingSplit

1 1 2 58.0000

2 3 4 24.0000

3 5 6 10.0000

4 7 8 13.0000

5 9 10 14.0000

---------

Summary:

Name Avg Std Min Max Count

MatchingSplit 23.8 17.73583942191629 10.0 58.0 5

## Computing normalized distances

Reporting distances divided by pre-computed empirical average values for random trees (generated according to Yule and uniform models, -N option) can help in an interpretation of the similarity level of analyzed trees in chosen metric. This functionality is available for trees with number of leaves between 4 and 1000 by using –N option. In the following example, the distance in the MS metric of each tree from a given set to the reference tree is computed. Analyzed trees have 15 leaves.

Input files: \examples\sclaed\ref\_tree.trees

\examples\sclaed\test\_set.trees

Invocation:

java -jar TreeCmp.jar -r ref\_tree.trees -d ms -i test\_set.trees -o test\_set.trees.r.out –N

Output file test\_set.trees.r.out:

No RefTree Tree MatchingSplit MatchingSplit\_toYuleAvg MatchingSplit\_toUnifAvg

1 1 1 43.0000 1.0742 0.9663

2 1 2 43.0000 1.0742 0.9663

3 1 3 41.0000 1.0242 0.9214

4 1 4 40.0000 0.9992 0.8989

5 1 5 43.0000 1.0742 0.9663

6 1 6 41.0000 1.0242 0.9214

7 1 7 43.0000 1.0742 0.9663

8 1 8 41.0000 1.0242 0.9214

9 1 9 39.0000 0.9742 0.8764

10 1 10 40.0000 0.9992 0.8989

11 1 11 0.0000 0.0000 0.0000

12 1 12 6.0000 0.1499 0.1348

Basic interpretation:

* Tree number 11 has the same topology as the reference tree.
* Tree number 12 is very similar to the reference tree in comparison to similarly of random on 15 leaves (the normalized distance is about 0.15 and 0.13 depending on the random model).
* Trees with numbers 1 to 10 are approximately as similar to the reference tree as random trees to each other (the normalized distance is close to 1).

In ordered to perform more advance similarity analysis, e.g. involving different model of generation of random trees, user my need to use TreeCmp twice:

* to compute distances between custom set of random trees generated by other software, e.g. Evolver application form PAML package (<http://abacus.gene.ucl.ac.uk/software/paml.html>) to obtain the empirical average distance in a particular metric or its distribution,
* to compute the distance between analyzed trees.

If the number of compared trees leaves is greater than 1000, it is possible to manually generate a set of random trees and calculate statistics for them. To generate a set of trees we can use PRTGen program – phylogenetic random trees generator. Let's assume that we want to generate 2,000 rooted trees on 1001 leaves using uniform model and save them to the file: trees.newick. Let's use command:

PRTGen –n 1001 –e 2000 –r –f trees.newick

Then, using TreeCmp, we calculate the value of the selected metric (for instance MC) between each subsequent pair of trees. We will get 1000 values:

java –jar TreeCmp.jar -w 2 -d MC -i trees.newick -o results.out

Based on these results, we can calculate desired values: (average, standard deviation, minimum, maximum, and subsequent quantiles: 0.02, 0.05, 0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9, 0.95, 0.97), e.g. in RStudio:

filename<-"<path\_to\_file>\results.out"

m<-read.table(filename,header = TRUE,sep = "\t")

v<-m[,4]

q\_seq<-c(0.02,0.05,0.1,0.2,0.3,0.4,0.5,0.6,0.7,0.8,0.9,0.95,0.97)

q<-quantile(v,q\_seq,type=1,name=FALSE)

row<-c(1001,mean(v),sd(v),min(v),max(v),q)

outfile<-"<path\_to\_file>\row.out"

write(row,file=outfile,append=TRUE,ncolumns=length(row),sep="\t")

Line such obtained in row.out file should be pasted into the appropriate file in the data folder. In that case it will be: unif\_MC.txt. Now Treecmp is ready for computing normalized MC distances for rooted trees on 1001 leaves based on uniform model.

## Finding the most similar trees in the input file

The most convenient comparison mode for such purpose is a matrix mode (-m). In the following example, the Matching Split distance is used.

Input file: \examples\plain2\plain2.trees

(a,(b,c),(d,e));

(a,b,(c,(d,e)));

(((a,b),c),d,e);

(a,(b,(c,d)),e);

Invocation:

java -jar TreeCmp.jar -m -d ms -i plain2.trees -o plain2.trees.m.out

Output file plain2.trees.m.out:

The most similar trees

No Tree1 Tree2 MatchingSplit

1 1 2 2.0000

2 1 3 2.0000

3 1 4 3.0000

4 2 3 0.0000

5 2 4 3.0000

6 3 4 3.0000

Trees number 2, i.e.: (a,b,(c,(d,e))) and 3, i.e.:(((a,b),c),d,e) in the input file are the most similar. In fact, they have the same topology (trees are assumed to be unrooted as metric for unrooted trees is used) because their distance is 0.

## Exporting data to other applications: MS Excel, R

To save a file in MS Excel format, just use the .xlsx extension in output data file name (option: -o <outputfile>.xlsx). Similarly, to save a file in CSV format, use the .csv extension in output data file name (option: -o <outputfile>.csv).

In order to pass data to R (<http://www.r-project.org/>) it is convenient to have the TreeCmp output file in a simple tabular form (therefore, it is recommended to avoid -I option, because it results in generation the summary section, which disturb the tabular order). Such files can be easily read by R environment by using for example the read.table function as follows:

treeCmpData<-read.table("C:\\Program Files\\TreeCmp\\examples\\plain\\plain.trees.m.out", header = TRUE, sep = "\t")

In the example, the file to read “plain.trees.m.out” is placed in “C:\Program Files\TreeCmp\examples\plain” folder.

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