

User Manual for SplitsTree4 V4.8

Daniel H. Huson and David Bryant

April 17, 2007



Contents

Contents	1
1 Introduction	4
2 Getting Started	5
3 Obtaining and Installing the Program	5
4 Program Overview	6
5 Splits, Trees and Networks	7
6 Opening, Reading and Writing Files	10
7 Estimating Distances	10
8 Building and Processing Trees	11
9 Building and Drawing Networks	11
10 Main Window	12

10.1 Network Tab	13
10.2 Data Tab	13
10.3 Source Tab	13
10.4 Tool bar	13
11 Graphical Interaction with the Network	13
12 Main Menus	14
12.1 File Menu	14
12.2 Edit Menu	16
12.3 View Menu	16
12.4 Data Menu	17
12.5 Distances Menu	19
12.6 Trees Menu	19
12.7 Network Menu	20
12.8 Analysis Menu	21
12.9 Draw Menu	21
12.10 Window Menu	22
12.11 Configuring Methods	23
13 Popup Menus	23
14 Tool Bar	24
15 Pipeline Window	24
15.1 Taxa Tab	25
15.2 Unaligned Tab	25
15.3 Characters Tab	25
15.4 Distances Tab	26
15.5 Quartets Tab	26
15.6 Trees Tab	26
15.7 Splits Tab	26
16 Export and Export Images Dialogs	27
16.1 Export	27
16.2 Export Image	28

17 Preferences Window	28
18 Additional Windows and Dialog	30
18.1 Open File	30
18.2 Choose Datatype	30
18.3 Save As	30
18.4 Find/Replace Window	31
18.5 Node and Edge Window	31
18.6 Highlight Confidence Window	31
18.7 Bootstrap Window	31
18.8 Message Window	32
18.9 About Window	32
19 Nexus Blocks	32
19.1 Taxa Block	32
19.2 Unaligned Block	33
19.3 Characters Block	33
19.4 Distances Block	34
19.5 Quartets Block	34
19.6 Trees Block	35
19.7 Splits Block	35
19.8 Network Block	36
19.9 Bootstrap Block	37
19.10Sets Block	37
19.11ST Assumptions Block	38
20 File Formats	38
21 All Methods	39
22 Command-Line Options and Mode	47
23 Examples	49
24 Acknowledgments	49
References	49

1 Introduction

Disclaimer: This software is provided "AS IS" without warranty of any kind. This is developmental code, and we make no pretension as to it being bug-free and totally reliable. Use at your own risk. We will accept no liability for any damages incurred through the use of this software. Use of the SplitsTree4 is free, however the program is not open source.

Type-setting conventions: In this manual we use e.g. **Network**→NeighborNet to indicate the NeighborNet menu item in the Network menu. We use e.g. **Main:Source** to indicate the Source tab in the Main window.

How to cite: If you publish results obtained in part by using SplitsTree4 , then we require that you acknowledge this by citing the program as follows:

- D.H. Huson and D. Bryant, *Application of Phylogenetic Networks in Evolutionary Studies*, Molecular Biology and Evolution, 23(2):254-267, 2006. software available from www.splitstree.org, or
- D.H. Huson, *SplitsTree: A program for analyzing and visualizing evolutionary data*. Bioinformatics, 14(10):68–73, 1998.

Evolutionary relationships are usually represented using phylogenetic trees, based on a model of evolution dominated by mutations and speciation events. More realistic models must also account for gene genesis, loss and duplication events, hybridization, horizontal gene transfer or recombination. Here, phylogenetic networks have a role to play.

Moreover, network methods also provide a value tool for phylogenetic inference even when reticulation events do not play an important role. The combined effect of *sampling error* and *systematic error* makes phylogenetics an uncertain science, and network methods provide tools for representing and quantifying this uncertainty.

The aim of SplitsTree4 is to provide a framework for evolutionary analysis using both trees and networks. The program takes as input a set of taxa represented by characters (that is, aligned sequences), distances, quartets, trees or splits and produces as output trees or networks using a number of different methods.

This document provides both an introduction and a reference manual for SplitsTree4 .

2 Getting Started

This section describes how to get started.

First, download an installer for the program from www.splitstree.org, see Section 3 for details.

Use the **File→Open** menu item to open a file containing data such as character sequences, distances or trees. The file must be in one of the following formats: **Nexus**, **ClustalW**, **Phylip**, **FastA** or **Newick**. Alternatively, select the **Main:Source** tab of the **Main** window and enter data in one of these formats by hand.

Example files are provided with the program. They are contained in the **examples** sub directory of the installation directory. The precise location of the installation directory depends upon your operating system.

Use the different menu items to determine which methods are applied to the input data. The **Distances**, **Trees** and **Network** menus contain items that determine how to compute distances, trees or a network from the given data. Some of the methods provided have parameters that can be set using the **Analysis→Configure Pipeline** or **Analysis→Configure Recent Methods** items.

The computed data can be viewed in text form in the **Main:Data** tab and can be saved using the **File→Save As** item.

The computed network or tree is displayed in the **Main:Network** tab. Attributes of the network can be changed by selecting nodes or branches (also called edges) and using the **Nodes and Edges** dialog which is reachable using the **View→Nodes and Edges** item.

Individual blocks of data can be saved in a number of different formats using the **File→Export** item. The network displayed in the **Main:Network** tab can be saved in a graphics format using the **File→Export Image** item.

3 Obtaining and Installing the Program

SplitsTree4 is written in Java and requires a Java runtime environment version 1.4.2 or newer, freely available from www.java.org.

SplitsTree4 is installed using an installer program that is freely available from www.splitstree.org. There are four different installers, targeting different operating systems:

- **splitstree_windows_4.8.exe** provides an installer for Windows.
- **splitstree_macos_4.8.sit** provides an installer for Mac OS.
- **splitstree_linux_4.8.rpm** provides a RPM package for Linux.
- **splitstree_unix_4.8.sh** provides a shell installer for Linux and Unix.

4 Program Overview

In this section, we give an overview over the main design goals and features of this program. Basic knowledge of the underlying design of the program should make it easier to use the program.

SplitsTree4 is written in the programming language Java. The advantages of this is that we can provide versions that run under the Linux, MacOS, Windows and Unix operating systems. Additionally, this makes it possible for the program to support plug-ins that can add new functionality to the program, such as new methods and import or export modules. A potential draw-back is that an algorithm implemented in Java will generally run slower than the same algorithm implemented in *C* or *C++*.

Earlier versions 1 – 3 of the program [16] were written in *C++* and only contain a small part of what is now available with SplitsTree4 .

SplitsTree4 uses multi-threading and supports multiple documents. This means that that you can work on more than one data set simultaneously, in different windows, and run multiple calculations simultaneously, making use of multiple processors, when available.

A *document* consists of an individual data set and possesses its own **Main** window. The document is discarded, when its window is closed. The program is based on the **Nexus** format, as introduced in [25] and the data associated with a document is organized into *blocks*, and each such block of data is represented by a corresponding “block” in the Nexus format. The blocks are:

- **Taxa**: the names of all taxa.
- **Unaligned**: unaligned sequences.
- **Characters**: aligned character sequences.
- **Distances**: pairwise distances between taxa.
- **Quartets**: (possibly weighted) quartet topologies.
- **Trees**: list of (possibly partial) trees.
- **Splits**: (possibly weighted) splits.
- **Network**: phylogenetic tree or network.
- **ST_Assumptions**: contains all methods and options used to compute data.
- **ST_Bootstrap**: bootstrap support of splits.

The first eight blocks **Taxa**, **Unaligned**, **Characters**, **Distances**, **Quartets**, **Trees**, **Splits** and **Network** are organized as a *pipeline* and data is processed from left-to-right along this pipeline. Any non-empty document must contain a **Taxa** block and will usually contain an **ST_Assumptions** block. All other blocks are optional and the presence or absence of some block depends on the set of computations that the user has selected.

We will use the term *source block* to denote the left-most block in the pipeline (excluding the **Taxa** block). The source block contains the original data that is provided to the program. Any computations performed by the program will update blocks from left to right along the pipeline, starting after the source block. Note that some types of computations do not fit into this pipeline design, for example **SplitsTree4** cannot provide any method that takes a **Splits** block and produces a **Trees** block, because the latter occurs before the former in the pipeline.

Typically, the user will provide a source-block and will then use different menu items to determine how the program will compute data along the pipeline until a **Network** block has been computed and an image of the **network** has been drawn in the **Main** window.

The program is designed to keep all blocks in the pipeline *synchronized* by enforcing that the different blocks only contain data that has been computed via the pipeline. For example, if you attempt to load data from a file in **Nexus** format that contains a **Taxa** and two or more other blocks, e.g. a **Characters** block or **Distances** block, then the program will request you to choose which of the two latter blocks you want to keep. (This does not apply if the file was created by **SplitsTree4** using the **File→Save** or **File→Save As** command, in which case the blocks in the file are synchronized, and so none must be discarded and no computations are necessary).

Once a source block has been provided, the program will proceed to perform a chain of *default calculations*, which differ, depending on the type of the source block. In the case of a **Characters** block, the program will use **UncorrectedP**, **NeighborNet** and then **EqualAngle** to compute a network for the data. In the case of a **Distances** block, **NeighborNet** and **EqualAngle** are used. In the case of a **Trees** block, the first tree in the block is displayed using the **EqualAngle** algorithm.

The data contained in the different blocks of the pipeline is displayed in the **Main:Data** tab. The data contained in the source block is edited using the **Main:Source** tab.

The program is designed to operate in two different modes: in a GUI mode, the program provides a GUI for the user to interact with the program. In **command-line mode**, the program reads commands from a file or from standard input and writes output to files or to standard output.

5 Splits, Trees and Networks

Here we give a brief introduction to some of the concepts from phylogenetics that the program is based on.

Evolutionary relationships between taxa are usually represented using a phylogenetic *tree*. Such trees are often computed from molecular sequence data, either directly, using methods such as parsimony, or indirectly, by first computing a distance matrix and then applying a method such as Neighbor-Joining.

Such approaches implicitly or explicitly model the evolution of a single gene under the assumption that the process is dominated by two types of events, mutations and speciation events. Under more complex models of evolution, i.e. involving gene loss and duplication, or hybridization, horizontal gene transfer or recombination, a single phylogenetic tree will often not be an appropriate repre-

sensation of the phylogentic history or of the different incompatible phylogenetic signals. Also, the presence of noise in a data set, or uncertainty due to inadequacies of the underlying model up on which a tree inference is based, may also make it necessary to use a more general graph, that is, a *network* , to represent the data.

The aim of **SplitsTree4** is to provide a frame-work for computing phylogenetic networks. As the name of the program suggests, it is based on the fundamental mathematical concept of a *split* . For example, if we are given an alignment of binary sequences

```
a 010011010110
b 100001011110
c 011001101110
d 010001101111
```

then each non-constant column in the alignment defines a split of the taxon set consisting of those taxa with the value 0 and those with the value 1. For example, the first column partitions the taxa into two sets $\{a, c, d\}$ and $\{b\}$, and thus gives rise to the split $\frac{\{a, c, d\}}{\{b\}} (= \frac{\{b\}}{\{a, c, d\}})$, while the fourth column does not define a split because the characters are constant.

Mathematically, for a set of X taxa any phylogenetic tree T defines a set of such splits called the *split encoding* $\Sigma(T)$ of T , as follows: deletion of any *branch* (also called *edge*) e in the tree produces two subtrees, T_A and T_B , say, and they give rise to a split $S = \frac{A}{B} = \frac{B}{A}$, consisting of the set A of all taxa contained in T_A and the set B of all taxa contained in T_B . In the literature, a split $\frac{A}{B}$ is sometimes denoted by $A|B$ or $\{A, B\}$.

For example, consider the tree T displayed in Figure 1. Its split encoding $\Sigma(T)$ contains 5 *trivial splits* that separate a single taxon from all other taxa, and 2 *non-trivial splits* that contain at least two taxa in both parts. The trivial splits are:

$$\frac{\{a\}}{\{b, c, d, e\}}, \frac{\{b\}}{\{a, c, d, e\}}, \frac{\{c\}}{\{a, b, d, e\}}, \frac{\{d\}}{\{a, b, c, e\}} \text{ and } \frac{\{e\}}{\{a, b, c, d\}},$$

and the non-trivial ones are:

$$\frac{\{a, b\}}{\{c, d, e\}} \text{ and } \frac{\{a, b, e\}}{\{c, d\}}.$$

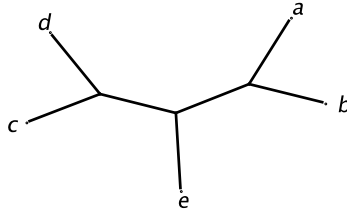


Figure 1: An unrooted phylogenetic tree.

A fundamental result in mathematical phylogeny states that any given set of splits Σ corresponds to some phylogenetic tree T , with $\Sigma(T) = \Sigma$, if and only if Σ is *compatible* [6]. Thus, any

phylogenetic tree may be viewed as a graph whose task it is to give a visual representation of a given set of compatible splits.

A phylogenetic tree can be thought of as an idealized representation of the historical relationships among a set of taxa; and tree building methods are attempts to find a the set of compatible splits that are most consistent with the data according to some algorithm. Often there are multiple tree that are equally consistent with the data, i.e. multiple sets of compatible splits. In general that collection of splits will be incompatible. Moreover, there exist inference methods such as [split decomposition](#) and [Neighbor-Net](#) that compute a set of incompatible splits from the data in the form of a given distance matrix. Note that split decomposition produces a set of splits that are called *weakly compatible* , while Neighbor-Net produces a set of splits that is *circular* , as originally defined in [1].

A *split network* is a more general type of phylogenetic graph that can represent any collection of splits, whether incompatible or not [1]. For a compatible set of splits, it is always possible to represent each split by a single [branch](#), and thus the resulting graph is a tree. In general, however, this will not be possible and in a split network usually a whole band of *parallel branches* (also called *parallel edges*) is required to represent a single split. A phylogenetic tree is therefore a special case of a split network. In [SplitsTree4](#) , if you click on any branch in the representation of a split network, then all branches corresponding to the same split will be highlighted. If one were to delete all branches corresponding to a given split S , then the remaining graph will consist of precisely two components, G_A and G_B , and, as above, we have $S = \frac{A}{B}$, where A is the set of all taxa contained in G_A and B is the set of all taxa contained in G_B .

There is an important difference between phylogenetic trees and more general split networks: Any rooted tree has a direct interpretation in evolutionary terms: the leaves represent taxa and the internal nodes represent speciation events. In a (possibly rooted) split network, the internal nodes do not have such a direct interpretation. Instead, split networks must be viewed on a more abstract level as networks giving a visual representation of incompatible signals, that is, showing how “tree-like” or “certain” parts of a phylogeny are.

A *reticulate network* is used to describe *reticulate evolution* and there are two types of such networks, one is the *hybridization network* and the other is the *recombination network* . A hybridization network describes evolution in the presence of *hybridization* events such as *allopolyploidization* or *diploid* hybridization, and in both cases the net effect is that a new species arises that has obtained some of its genes via one ancestral species and the other genes via another ancestral species. A recombination network is used to describe evolution in the presence of *recombination* events. As presently implemented in [SplitsTree4](#) , such recombination networks can only be computed for *binary* sequences. Such sequences may, for example, indicate the presence or absence of a certain restriction site at a certain position. There are two main differences between hybridization networks and recombination networks. First, the former are based on gene trees, whereas the latter are based on sequences. Second, a recombination network is additionally annotated by inferred sequences at each of the nodes and the mutations that happen along each of the branches. Moreover, the recombination nodes are usually annotated with a description of which parts of the two ancestor sequences are used in the recombination event.

In contrast to split networks, these types of networks can be directly interpreted in evolutionary terms. Recent research has revealed that there are strong connections between split networks and

reticulate networks [21, 22] and these are used in SplitsTree4 to provide methods for computing reticulate networks.

A more detailed discussion of the fundamental concepts can be found in [17] and [20].

6 Opening, Reading and Writing Files

To open a file, select the **File→Open** menu item and then browse to the desired file. Alternatively, if the file was recently opened by the program, then it may be contained in the **File→Open Recent** submenu.

The native file format of SplitsTree4 is based on the **Nexus** format, see [25]. However, the program can also parse a number of other formats, including **ClustalW** format for aligned sequences, **Phylip** format for sequences and distances, **FastA** format for aligned distances and **Newick** format for trees. Earlier versions had separate Open and Import menu items; these have now been combined.

If a file is opened while the **Main:Network** tab or **Main:Data** tab is open, then the program will attempt to parse and execute the file. If the file is a complete Nexus file previously generated by SplitsTree4, then the network described in the file will be displayed. If the file is in some other format, then, depending on the type of content, the program will perform a chain of **default calculations**.

If a file is opened while the **Main:Source** tab is opened, then the contents of the file is read into the program and displayed as text in the Source-tab, but no computation is performed.

To save the complete data associated with a given window in SplitsTree4's native Nexus format, use the **File→Save** or **File→Save As** menu items. To save selected blocks of data, or to export data in a different file format, use the **File→Export**. A picture of the computed **network** can be saved using the **File→Export Image** menu item.

7 Estimating Distances

Many methods in phylogenetics begin by estimating distances between the taxa. This is done by taking sequences two at a time and estimating the average number of mutations that occurred on the paths from them to their most recent common ancestor. If the rate of mutation was constant, then this will be proportional to their divergence time. SplitsTree4 provides a large number of methods for estimating distances, for sequences and other types of data.

The **UncorrectedP** method computes the proportion of positions at which two sequences differ. For DNA or RNA sequences, there are choices over how ambiguous state codes (such as W, M, V) are handled. Ignore, means that these states are treated as missing states. 'Average' means that the contribution at a site is averaged over all possible resolutions of the ambiguous codes, with the exception that sites having the same ambiguous code contribute zero. 'Match' looks at each possible state in each sequence, counts one if the state is not a possible resolution of the ambiguous

code in the other sequence, and normalises the count by the number of states for each ambiguous code.

The **Distances** menu lists several standard distance estimation methods. Most of these have parameters that can be changed. When you select the menu item the **Pipeline** window will open and will display a panel for setting the parameters of the method. If you always use the same parameters for a given method, or if the method has no parameters, then selecting the “Don’t show this dialog to configure this method again” will prevent the dialog from appearing again. All methods can be configured using the **Pipeline** window, accessed by **Analysis→Configure Pipeline** or **Analysis→Configure Recent Methods**.

8 Building and Processing Trees

The **Tree** menu provides a number of methods for constructing trees, both from distance and character data. The following distance-based methods are available: **NJ**, **BioNJ**, **UPGMA**, **Buneman** and **Refined Buneman**.

The program provides ‘front-ends’ for several tree based programs, currently **PhyML** and **PhyIip**. **SplitsTree4** will pass these programs the current sequences, and place any trees produced in the **SplitsTree4 Trees** block. Our distribution of **SplitsTree4** does not provide any external programs, so you must install them separately. In the dialog that is associated with such a method you must specify the location of the external application, e.g. using the **PhyML Path** text field. Once entered, the program will remember this entry as the default value.

SplitsTree4 provides a number of methods for processing a collection of trees. The **pmenu-itemTreesTreeSelector** menu contains items to select individual trees or to compute a consensus of trees. If the **Trees→TreesSelector** is selected, then the **Trees→Previous Tree** and **Trees→Next Tree** items can be used to move from one tree to the next in the **Trees** block. Pressing the shift-key when selecting either of these methods will move to the first tree, or the last tree, respectively. The **Trees→ConsensusTree** can be used to compute the majority or strict consensus tree.

As with distances, when you select such a menu item the **Pipeline** window will open and will display a panel for setting the parameters of the method. If you always use the same parameters for a given method, or if the method has no parameters, then selecting the “Don’t show this dialog to configure this method again” will prevent the dialog from appearing again. All methods can be configured using the **Pipeline** window, accessed by **Analysis→Configure Pipeline** or **Analysis→Configure Recent Methods**.

9 Building and Drawing Networks

The **Network** menu provides methods for computing phylogenetic networks from character sequences, distances and trees.

Methods that compute a [split network](#) directly from character sequences provided in the [Characters](#) block are [ParsimonySplits](#), [MedianNetwork](#) [SpectralSplits](#). Note that the Median network method requires binary sequences. However, given DNA or RNA, this program will detect all sites that contain precisely two character-states and will build a Median network from these.

Two methods for computing split networks from distances provided in the [Distances](#) block are [SplitDecomposition](#) and [NeighborNet](#).

If a set of phylogenetic trees in the [Trees](#) block all contain the full set of taxa listed in the [Taxa](#) block, then the [ConsensusNetwork](#) method can be applied to compute a consensus network. If, however, the [Trees](#) block contains partial trees, that is, trees that do not necessarily all involve identical sets of taxa, then the [SuperNetwork](#) or [FilteredSuperNetwork](#) method can be used to compute a super network.

The [Networks](#) menu also provides two methods for computing *reticulate networks*. If the [Trees](#) block contains more than one tree, then the [HybridizationNetwork](#) method can be used to compute a *hybridization network* for the given data set. More precisely, the [HybridizationNetwork](#) requires that the [Trees](#) and [Splits](#) blocks are present and non-empty. If the [Characters](#) block contains binary sequence data, then the [RecombinationNetwork](#) method can be used to compute a *recombination network*.

The [Draw](#) menu determines which algorithm is used to construct the final visualization of the [tree](#) or [network](#). Existing methods are [EqualAngle](#), [RootedEqualAngle](#), [Phylogram](#) and [ReticulateNetwork](#). Additionally, the [Draw→HideSelectedSplits](#) can be used to remove selected splits from the network and the [Draw→Select Trees](#) can be used to highlight different trees contained in a split network.

10 Main Window

With [SplitsTree4](#), multiple documents can be opened and worked on simultaneously. Each document is represented by a [Main](#) window. Additional windows or dialogs are sometimes opened to perform certain tasks.

The [Main](#) window has three tabbed panes:

- the [Main:Network](#) tab is used to display the computed tree or network,
- the [Main:Data](#) tab can be used to inspect all data associated with a given document in text format, and
- the [Main:Source](#) tab can be used to inspect or edit the “source” data from which all other data is derived.

We now discuss each separately:

10.1 Network Tab

The `Main:Network` tab is used to display the computed tree or network. A [status line](#) of text along the bottom of the network pane gives a summary of the data, such as number of taxa, length of sequences, etc, and a summary of the methods used to compute the network. Optionally, a [scale bar](#) can be displayed in the upper left corner of the Network-tab to indicate the scale of branches. The scale bar is can be turned on or of using the [Preferences:General](#) tab. The scale bar can be configured using the [Preferences:Status Line](#) tab.

10.2 Data Tab

The `Main:Data` tab provides a textual display of the data associated with the given document in the program's native [Nexus](#) format, organized in a linear list of items that can be either collapsed or expanded. This view of the data is read-only.

10.3 Source Tab

The `Main:Source` tab provides an editable view of the source data associated with the given view. This view provides three different functionalities, namely:

- data can be entered by hand or by copy-and-paste, in any of the file formats that the program supports, see [Section 20](#),
- once the data has been executed, the source data is displayed in [Nexus](#) format, and
- if an error is encountered while parsing an input file, the file is opened and the line in which the error was detected is selected.

10.4 Tool bar

The *tool bar* associated with the window provides buttons for quick access to many of the menu items. It can be configured using the `Preferences:Toolbar` tab.

11 Graphical Interaction with the Network

We now describe how the user can interactively modify the layout and attributes of the displayed [network](#). The [View](#) menu is used to rotate, move, and zoom in and out. Alternatively, the view can be changed using a wheel mouse, together with the following modifier keys:

none	zoom in and out.
Shift	rotate
ALT/option	move vertically
Shift & ALT/option	move horizontally

Here are additional modifications that can be performed:

- The graph can be dragged around by clicking and dragging a node.
- If all selected nodes lie on one side of a band of [parallel branches](#) representing a single split, then clicking and dragging on one of the nodes will change the angle of the branches.
- By default, clicking on an edge will select *all* edges representing the same split, and all nodes on the smaller side of the represented split. To select all nodes on the other side of the split, use the `View→Invert Selection` menu item. To allow selection of individual edges, deselect the `Split-Selection Mode` check box in the [Preferences:General](#) tab.
- Select a node by clicking on it. Shift-click outside of the network and drag a rectangle to select several nodes at once. Hold the shift key and click to add or remove further nodes from the set selected.
- Click on a text label to edit it.
- Double-click on an edge to edit its label.
- By default, node labels are automatically positioned to avoid overlap. Any label that has been interactively repositioned by the user is no longer automatically positioned. To apply automatic positioning all node labels, including those that have been moved by the user, deselect and then select the `View→Auto Layout Node Labels` check box.
- Many aspects of the visual representation of nodes and edges can be modified using the [Node and Edge Window](#), which is opened using the `View→Nodes and Edges` menu item.

12 Main Menus

We now discuss all menus of the [Main](#) window. Some of the menus differ, depending upon which tab is selected in the [Main](#) window.

12.1 File Menu

The `File` menu contains the usual file-related items:

- The `File→New` item opens a new [Main](#) window and selects its [Main:Source](#) tab.

- The **File→Open** item provides an **Open File** dialog to open a file containing input data in one of the supported formats.

If the file contains character sequences, then the program must know whether the sequences are DNA, RNA, protein or “standard” (0,1) data. In **Nexus** files, the datatype is explicitly given. In other file formats this not always the case. If the program cannot guess the data type (e.g., if all character states are 'A'), then the program will display a **Choose Datatype** dialog and prompt the user to specify the data type.

If the current document is non-empty, then the selected file is opened in a new **Main** window.

- The **File→Open Recent** submenu provides access to recently opened or saved files.
- The **File→Replace** item is used to replaced the current data by a new data set. This is not available under MacOS.
- The **File→Clones** item clones the current window.
- The **File→Close** item closes the current document.
- The **File→Save** item saves the current document to its file, if known.
- The **File→Save As** item provides a **Save As** dialog and saves the current document to selected file.
- The **File→Export** item opens the **Export** dialog which is used to save individual data in a number of different formats.
- The **File→Export Image** item opens the **Export Image** dialog which is used to save the current network in a number of different formats.
- The **File→Tools** item provides a submenu of tools. The **File→Tools→Load Trees** item can be used to merge a set of trees into one document. This item opens a dialog for opening multiple files. The program parses all given files and extracts any trees found in them and opens a new **Document** containing all trees found. The **File→Tools→Load Multi-Labeled Tree** item can be used to read a single tree containing multiple labels and to convert it into a single-labeled split network (and will be made available upon publication of the paper describing this new method). The **File→Tools→Concatenate Sequences** item can be used to concatenate sequences from different files into one document. It requires that each of the input files uses exactly the same set of taxon labels. The **File→Tools→Group Identical Haplotypes** tools is used to detect multiple identical sequences or taxa. These are recognised from the **Characters** or, if there is no **Characters** block, from the distances block. Only selected sites are used to test if two taxa are distinguishable. A new document is created; identical taxa are collapsed into a single taxa with label 'TYPE_x' (x ranging from 1,2,3,..., and the original taxa stored in the info field of the **Taxa** block. Note that files with hundreds or thousands of identical sequences can cause SplitsTree to stall. One solution is to cancel the computations after data is read in, and to then use this tool to create a new smaller document containing only distinct sequences.
- The **File→Print** item is used to print the current network.
- The **File→Quit** item quits the program, after asking whether to save unsaved changes.

12.2 Edit Menu

The **Edit** menu contains the usual edit-related items:

- The **Edit→Undo** item is used to undo text-editing, interactive network manipulation and any item chosen from the **View** menu.
- The **Edit→Redo** item is used to redo text-editing, interactive network manipulation and any item chosen from the **View** menu.
- The **Edit→Cut** item is used to cut text.
- The **Edit→Copy** item is used to copy text or to copy the current network as an image.
- The **Edit→Paste** item is used to paste text.
- The **Edit→Select All** item is used to select all nodes and edges of a network.
- The **Edit→Invert Selection** item is used to invert the selection of nodes.
- The **Edit→Taxon Sets** submenu is used to select all nodes labeled by a given taxon set or to add new taxa sets to the **SETS** block. The **SETS** block can be copied to other files to provide a convenient method to quickly label or select taxa in different taxa groups. The **Edit→TaxonSets→All** item selects all taxa. The **Edit→TaxonSets→New taxa set...** item creates a new taxa set with the currently selected taxa and a name provided by the user. Note that the name should be a valid NEXUS name. If a set exists with this name, it is overwritten. The **Edit→TaxonSets→Clear All Taxa Sets...** removes all taxa sets from the **SETS** block.
- The **Edit→Find/Replace** item opens the **Find/Replace** dialog.
- The **Edit→Go to Line** item is used to scroll to a particular line in the **Main:Source** tab.
- The **Edit→Preferences** opens the **Preferences** window.

12.3 View Menu

The **View** menu contains items that control aspects of the visualization of a network.

This menu changes depending on which tab on is in. In the **Main:Network** tab, the **View** menu contains the following items, which are all undo-able and redo-able:

- The **View→Reset** item resets the layout of the network, if used while in the **Main:Network** tab, and it clears the text area, while in the **Main:Source** tab.
- The **View→Zoom In** item is used to zoom into the network.

- The `View→Zoom Out` item is used to zoom out from the network.
- The `View→Rotate Left` and `View→Rotate Right` items are used to rotate the network.
- The `View→Flip` item changes the layout of the network to its mirror image.
- The `View→Nodes and Edges` item opens the `Nodes and Edges` window that is used to modify the graphical attributes of the nodes and edges of the network.
- The `View→Highlight Confidence` item opens the `Confidence` window that can be used to control how confidence values for splits or edges are highlighted in the network.
- The `View→Auto Layout Node Labels` check box item controls whether node labels are automatically positioned or whether the user can interactively reposition them.

In the `Main:Data` tab, the `View` menu contains the following items:

- The `View→Expand All` and `View→Collapse All` items expand and collapse all nodes in the displayed data tree.
- The `View→Characters`, `View→Distances` and `View→Splits` submenus control the format used to write the `Characters`, `Distances` and `Splits` blocks, respectively.

In the `Main:Source` tab, the `View` menu only contains a `View→Reset` item, which is used to clear the source text area.

12.4 Data Menu

The `Data` menu contains the following items:

- The `Data→Keep Only Selected Taxa` item removes all taxa from the analysis, whose nodes in the network are unselected. All data associated with the removed taxa is removed from the original `source block` and all subsequent data is recomputed.
- The `Data→Exclude Selected Taxa` item removes all taxa from the analysis, whose nodes in the network are selected. All data associated with the removed taxa is removed from the original `source block` and all subsequent data is recomputed.
- The `Data→Restore All Taxa` item restores all taxa that were previously removed using either of the previous two menu items, or by the next menu item.
- The `Data→Filter Taxa` item opens the `Pipeline:Taxa:Filter` tab that can be used to interactively include or exclude taxa from the analysis.
- The `Data→Exclude Gap Sites` item excludes from computation all sites in a `Characters` block that contain a gap or missing character in any of the sequences.

- The **Data→Exclude Parsimony-Uninformative Sites** item excludes from computation all sites in a **Characters** block that are parsimony-uninformative, that is, which are constant across all but one sequence.
- The **Data→Exclude Constant Sites** item excludes from computation all sites in a **Characters** block that are constant across all sequences.
- The **Data→Restore All Sites** item restores all sites that were excluded using the above menu items.
- The **Data→Filter Characters** item opens the **Pipeline:Characters:Filter** tab that can be used to interactively include or exclude sites from the analysis.
- The **Data→Greedy Make Compatible** item uses a greedy approach to make the splits in the **Splits** block **compatible**: in decreasing order of weight, the algorithm adds the next split to the set of kept splits, if it is compatible with all splits that have already been kept.
- The **Data→Greedy Make Weakly Compatible** item uses a greedy approach to make the splits in the **Splits** block **weakly compatible**: in decreasing order of weight, the algorithm adds the next split to the set of kept splits, if it is weakly-compatible with all splits that have already been kept.
- The **Data→Exclude Selected Splits** item removes all splits from the analysis, whose edges in the network are selected. The **Splits** and **Network** block are modified accordingly. See also: **Draw→Hide Selected Splits**.
- The **Data→Restore All Splits** item restores all splits that were excluded using the previous menu item.
- The **Data→Filter Splits** item opens the **Pipeline:Splits:Filter** tab that can be used to interactively include or exclude splits from the analysis.
- The **Data→Filter Trees** item opens the **Pipeline:Tree:Filter** tab that can be used to interactively include or exclude trees from the analysis.
- The **Data→Set Tree Names** item opens a dialog that allows one to set the tree names.

In the **Main:Source** tab, the **Data** contains three commands for processing the text contained in the *source text area* displayed in the tab:

- The **Data→Execute** item parses the text in **Nexus** format and executes it.
- The **Data→Convert to Nexus** item parses the text in any of the supported file formats (see Section 20), including **old Nexus**, and converts it into **Nexus** format and replaces the text in the source text area by the converted text.

If either of the first two items is used and the text contained in the text area is not in **Newick** format, then a dialog is displayed that allows the user to request to import the text before parsing it.

The **Data→Load** item is particularly useful if you would like to load a large data set and would like to prevent the program from performing any *initial calculations* . First open the file using the **File→Open** item while in the **Main:Source** tab, then use the **Data→Load** item to “load” the data into the program and then switch to the **Main:Network**. No calculations will be performed until you choose some menu item.

12.5 Distances Menu

The **Distances** menu contains the following items:

- The **Distances→UncorrectedP** item requests the program to compute distances using the **UncorrectedP** method.
- The **Distances→LogDet** item requests the program to compute distances using the **LogDet** method.
- The **Distances→HKY85** item requests the program to compute distances using the **HKY85** method.
- The **Distances→JukesCantor** item requests the program to compute distances using the **JukesCantor** method.
- The **Distances→K2P** item requests the program to compute distances using the **K2P** method.
- The **Distances→K3ST** item requests the program to compute distances using the **K3ST** method.
- The **Distances→F81** item requests the program to compute distances using the **F81** method.
- The **Distances→F84** item requests the program to compute distances using the **F84** method.
- The **Distances→ProteinMLdist** item requests the program to compute distances using the **ProteinMLdist** method.
- The **Distances→NeiMiller** item requests the program to compute distances using the **NeiMiller** method.
- The **Distances→GeneContentDistance** item requests the program to compute distances using the **GeneContentDistance** method.

12.6 Trees Menu

The **Trees** menu contains the following items:

- The `Trees→NJ` item requests the program to compute a tree using the [NJ](#) method.
- The `Trees→BIONJ` item requests the program to compute a tree using the [BIONJ](#) method.
- The `Trees→UPGMA` item requests the program to compute a tree using the [UPGMA](#) method.
- The `Trees→BunemanTree` item requests the program to compute a tree using the [BunemanTree](#) method.
- The `Trees→RefinedBunemanTree` item requests the program to compute a tree using the [RefinedBunemanTree](#) method.
- The `Trees→TreeSelector` item requests the program to compute a tree using the [TreeSelector](#) method.
- If the [Trees→TreeSelector](#) is selected, then the `Trees→Previous Tree` and `Trees→Next Tree` items are used to move from one tree to the next in the [Trees](#) block. Pressing the shift-key when selecting either of these methods will move to the first tree, or the last tree, respectively.
- The `Trees→ConsensusTree` requests the program to compute compute the majority or strict consensus tree.
- The `Trees→PhylipParsimony` item requests the program to compute a tree using the [PhylipParsimony](#) method.
- The `Trees→PhyML` item requests the program to compute a tree using the [PhyML](#) method.

12.7 Network Menu

The `Networks` menu contains the following items:

- The `Networks→NeighborNet` item requests to compute a network using the [NeighborNet](#) method.
- The `Networks→SplitDecomposition` item requests to compute a network using the [SplitDecomposition](#) method.
- The `Networks→ParsimonySplits` item requests to compute a network using the [ParsimonySplits](#) method.
- The `Networks→ConsensusNetwork` item requests to compute a network using the [ConsensusNetwork](#) method.
- The `Networks→SuperNetwork` item requests to compute a network using the [SuperNetwork](#) method.
- The `Networks→FilteredSuperNetwork` item requests to compute a network using the [FilteredSuperNetwork](#) method.

- The `Networks→MedianNetwork` item requests to compute a network using the `MedianNetwork` method.
- The `Networks→SpectralSplits` item requests to compute a network using the `SpectralSplits` method.
- The `Networks→HybridizationNetwork` item requests to compute a network using the `HybridizationNetwork` method.
- The `Networks→RecombinationNetwork` item requests to compute a network using the `RecombinationNetwork` method.

12.8 Analysis Menu

The `Analysis` menu contains the following items:

- The `Analysis→Bootstrap` item opens the `Bootstrap` dialog.
- The `Analysis→Show Bootstrap Network` item opens a new `Main Viewer` depicting a `network` that is based on all splits that occurred in any of the bootstrap replicates. Note that this item is enabled only after bootstrapping has been completed.
- The `Analysis→Show Confidence Network` item opens a new `Main Viewer` containing a `network` that represents a 95% confidence set for the trees or networks estimated. Note that this item is enabled only after bootstrapping has been completed.
- The `Analysis→Estimate Invariable Sites` uses the capture-recapture method of [24].
- The `Analysis→Compute Phylogenetic Diversity` is enabled when a set of taxa are computed. It computes the sum of the weights for all splits that separate these taxa into two non-empty groups. On a tree, this is equivalent to the phylogenetic diversity measure of [9], as these splits will be exactly those lying on a path between two taxa.
- The `Analysis→Conduct Phi test for Recombination` tests for recombination using the Phi test of [4].
- The `Analysis→Configure Pipeline` item opens the `Pipeline` window which can be used to configure the parameters of a given method.
- The `Analysis→Configure Recent Methods` submenus lists all recently used methods and can be used to open the `Pipeline` window in the appropriate tab to configure the parameters of a given method.

12.9 Draw Menu

The `Draw` menu contains the following items:

- The `Draw→EqualAngle` item requests to draw a network using the [EqualAngle](#) method.
- The `Draw→RootedEqualAngle` item requests to draw a rooted network using the [RootedEqualAngle](#) method.
- The `Draw→ConvexHull` item requests to draw a network using the [ConvexHull](#) method.
- The `Draw→Phylogram` item requests to draw a tree using the [Phylogram](#) method.
- The `Draw→NoGraph` item requests that no network be computed.
- The `Draw→Select Characters` item opens the [Pipeline:Characters:Select](#) tab that can be used to select individual characters (that is, sites in the given sequence alignment) to be displayed on the nodes of the network. set of selected input trees.
- The `Draw→Hide Selected Splits` item can be used to remove all selected splits directly from the depicted network. This does not change the [Splits](#) block. See also: [Data→Exclude Selected Splits](#).
- The `Draw→Hide Non-Selected Splits` item can be used to remove all non-selected splits directly from the depicted network. This does not change the [Splits](#) block. See also: [Data→Exclude Selected Splits](#).
- The `Draw→Hide Incompatible Splits` item can be used to remove all splits that are incompatible with the set of currently selected splits directly from the depicted network. This does not change the [Splits](#) block.
- The `Draw→Redraw All Splits` item redraws the network using all splits, including those excluded using the previous menu item.
- The `Draw→Select Trees` item opens the [Pipeline:Trees:Select](#) tab that can be used to select all splits in a given network that are contained in a given set of selected input trees.

12.10 Window Menu

The `Window` menu contains the following items:

- The `Window→About` item shows information about the version of `SplitsTree4`.
- The `Window→How to Cite` item gives instructions on how to cite the program.
- The `Window→Nexus Syntax` submenu provides the syntax of all [Nexus](#) blocks that `SplitsTree4` can process.
- The `Window→Command Syntax` item summarizes the syntax of the commands that can be used with the [command-line mode](#) version of the program, or which can be present in an input file.
- The `Window→Enter a Command` item allows entry of a single command.

- The `Window→Message Window` item opens the message window.

The bottom of the `Window` menu contains a list of all open windows.

12.11 Configuring Methods

By default, selecting a menu item that applies some method to the given data, such as e.g. the `Trees→NeighborJoining`, will open up a dialog in the `Pipeline` window which one can use to configure the method by selecting appropriate options. For methods that have no options, or that are always used with the same options, selecting the *Don't show this dialog for this method again* checkbox will tell the program not to open the dialog in the future, but rather to immediately apply the named method. Note that configuration dialogs hidden in this way can still be accessed using the `Analysis→Configure Recent Methods` submenu.

13 Popup Menus

Right-clicking on the `Main` window will open a popup menu. There are three different menus that will appear, depending upon what is hit by the mouse click.

If the mouse is clicked on a node of a network, then this opens the `Node` popup menu, which has the following items:

- The `Copy Label` copies the node label to the system clipboard.
- The `Edit Label` opens a dialog to edit the current node label.
- The `Exclude Selected Taxa` excludes the selected taxa from all computations.
- The `Show Name` labels the selected node by the names of any correspond taxa.
- The `Show Id` labels the selected node by the ids of any correspond taxa.
- The `Hide Label` hides the label of the selected node.
- The `Configure` opens the `Nodes and Edges:Nodes` tab of the `Nodes and Edges` window.

If the mouse is clicked on a edge of a network, then this opens the `Edge` popup menu, which has the following items:

- The `Copy Label` copies the edge label to the system clipboard.
- The `Edit Label` opens a dialog to edit the current edge label.
- The `Show Id` labels the selected edge by the id of the corresponding split.

- The `Show Weight` labels the selected edge by the weight of the corresponding split.
- The `Show Confidence` labels the selected edge by the confidence of the corresponding split.
- The `Show Interval` labels the selected edge by the confidence interval of the corresponding split.
- The `Hide Label` hides the label of the selected edge.
- The `Configure` opens the `Nodes and Edges:Edge` tab of the `Nodes and Edges` window.

If the mouse is clicked on the drawable region of the window, but not on any node or edge, then the `Window` popup menu will open, which has the following items:

- The `Zoom In` item is a short-cut to the `View→Zoom In` menu item.
- The `Zoom Out` item is a short-cut to the `View→Zoom Out` menu item.
- The `Reset` item is a short-cut to the `View→Reset` menu item.
- The `Reset Label Positions` item resets all node labels to their default positions.
- The `Restore All Taxa` item restores all excluded taxa.
- The `Restore All Splits` item restores all excluded splits.

14 Tool Bar

For easier access to frequently used options and methods, `SplitsTree4` provides a *tool bar* at the top of the `Main` window. The tool bar can be configured using the `Preferences:Toolbar` tab. By default, the tool bar contains the following items: `File→Open`, `File→Clone`, `File→Save As`, `File→Print`, `File→Export Image`, `Edit→Find/Replace`, `View→Reset`, `View→Zoom In`, `View→Zoom Out`, `View→Rotate Right`, `View→Rotate Left`, `Edit→Preferences`, `Window→Message Window`, `Draw→EqualAngle`, `Draw→RootedEqualAngle`, `Draw→Phylogram`, `Trees→Previous Tree` and `Trees→Next Tree`.

15 Pipeline Window

The `Pipeline` window is accessed using the `Analysis→Configure Pipeline` item. It controls all aspects of the computational `pipeline` that `SplitsTree4` uses. It is organized into nine tabs, reflecting the order of the blocks in the pipeline: `Pipeline:Taxa`, `Pipeline:Unaligned`, `Pipeline:Characters`, `Pipeline:Distances`, `Pipeline:Quartets`, `Pipeline:Trees` and `Pipeline:Splits`.

This is a potential source of confusion, because in the main menus of the [Main](#) window we organize the methods by the type of their *output*, rather than by the type of their *input*, as is done here.

Each tab in controls how a [block](#) of the given type is processed to produce a block of a subsequent type. Each of the tabs contains upto three sub-tabs, as mentioned below. Here we now describe each of the main tabs:

15.1 Taxa Tab

The `Pipeline:Taxa` tab consists of precisely one `Pipeline:Taxa:Filter` sub-tab, which is used to include (show) or exclude (hide) taxa from all calculations. All taxa listed in the *show list* are included in all calculations, where as all taxa listed in the *hide list* are removed from the data set. If a set of taxa are selected in the network displayed in the [Main](#) window, then these can be added to the show or hide list by pressing the appropriate `From Graph` button. Press `Show All` or `Hide All` to show all hide all known taxa.

Please note that the set of “shown” or “hidden” taxa will change automatically when viewing different [partial trees](#) to reflect the set of taxa contained in the current tree.

15.2 Unaligned Tab

The `Pipeline:Unaligned` tab consists of precisely one sub-tab. The `Pipeline:Unaligned:Method` sub-tab is used to choose and configure a method to apply to the current [Unaligned](#) block. Currently, the program provides three methods:

- The [NoAlign](#) item simply adds gaps to the ends of sequences to make the all have the same length.
- The [ClustalW](#) item can be used to call the program ClustalW externally to compute an alignment of the sequences.
- The [Muscle](#) item can be used to call the program Muscle externally to compute an alignment of the sequences.

15.3 Characters Tab

The `Pipeline:Characters` tab has three sub-tabs. The `Pipeline:Characters:Method` sub-tab is used to choose and configure a method to apply to the current [Characters](#) block. See [Section 21](#) for a description of all available methods.

The `Pipeline:Characters:Filter` sub-tab is used to exclude certain sites from the analysis. By default, the sites remain present in the [Characters](#) block and are simply masked during calculations. To remove the sites completely, press the `Delete` button.

The `Pipeline:Characters:Select` sub-tab is used to select certain sites in the [Characters](#) block for display of sites in the network.

15.4 Distances Tab

The `Pipeline:Distances` tab has only one sub-tab. The `Pipeline:Distances:Method` sub-tab is used to choose and configure a method to apply to the current [Distances](#) block. See [Section 21](#) for a description of all available methods.

15.5 Quartets Tab

The `Pipeline:Quartets` tab has only one sub-tab. The `Pipeline:Quartets:Method` sub-tab is used to choose and configure a method to apply to the current [Quartets](#) block. See [Section 21](#) for a description of all available methods.

15.6 Trees Tab

The `Pipeline:Trees` tab has three sub-tabs. The `Pipeline:Trees:Method` sub-tab is used to choose and configure a method to apply to the current [Trees](#) block. See [Section 21](#) for a description of all available methods.

The `Pipeline:Trees:Filter` sub-tab is used to exclude trees sites from the analysis.

The `Pipeline:Trees:Select` sub-tab is used to highlight a set of trees in the displayed network by selecting all splits in the network that are contained in the set of chosen trees.

15.7 Splits Tab

The `Pipeline:Splits` tab has two sub-tabs. The `Pipeline:Splits:Method` sub-tab is used to choose and configure a method to apply to the current [Splits](#) block. See [Section 21](#) for a description of all available methods.

The `Pipeline:Splits:Filter` sub-tab is used to modify the weights of splits using a *least squares* optimization, or to exclude certain splits from the analysis. Possible filters are:

Greedy Compatible	uses a greedy approach to makes the splits compatible : in decreasing order of weight, the algorithm adds the next split to the set of kept splits, if it is compatible with all splits that have already been kept.
Closest Tree	makes the splits compatible by computing the “closest tree”
Greedy Weakly Compatible	uses a greedy approach to makes the splits weakly compatible : in decreasing order of weight, the algorithm adds the next split to the set of kept splits, if it is weakly-compatible with all splits that have already been kept.
Weight Threshold	removes any split whose weight does not exceed the given threshold. Pressing the [Set button will open a histogram and slider to set the threshold.
Confidence Threshold	removes all splits whose confidence does not exceed the given threshold. A confidence value usually lies in the range $[0, 1]$ and can be obtained by bootstrapping, for example. Pressing the [Set button will open a histogram and slider to set the threshold.
Set Maximum Dimension	greedily removes a subset of splits that cause boxes in the network of dimension higher than the given threshold, as described in [18] .
None	applies no filter.
Use the Exclude Selected Splits to remove all splits that are currently selected in the displayed network.	

16 Export and Export Images Dialogs

16.1 Export

The **Export** dialog is opened using the [File→Export](#) item. It is used to save individual blocks of data in any of the formats described in [Section 20](#). The dialog show two lists. On the left, the set of available [Nexus](#) blocks. Is listed. Depending on the selection of blocks made by the user, on the right, the set of available export formats is listed.

Most of the dependences between blocks and formats should be apparent. One interesting feature is the following. If one has enabled graph editing using the [Allow Graph Editing](#) option in the [Preferences:General](#) tab and has interactively constructed a phylogenetic tree, then this tree can

be saved in [Newick](#) format. Hence, SplitsTree4 can be used for editing trees.

16.2 Export Image

The `Export Image` dialog is opened using the [File→Export Image](#) item. This dialog is used to save a picture of the current network in a number of different formats. The following graphics formats are supported:

- JPEG, “Joint Photographic Experts Group”.
- GIF, “Graphics Interchange Format”.
- EPS, “Encapsulated PostScript”.
- SVG, “Scalable Vector Graphics”.
- PNG, “Portable Network Graphics”.
- BMP, “Bitmap”.

There are two radio buttons, the `Save whole image` to save the whole image, and the `Save visible image` to save only the part of the image that is currently visible in the main viewer. If the chosen format is EPS, then selecting the `Convert text to graphics` check box will request the program to render all text as graphics, rather than fonts.

Pressing the apply button will open a standard file save dialog to determine where to save the graphics file.

17 Preferences Window

The `Preferences` window is opened using the [Edit→Preferences](#) item.

This window controls program *preferences*. Many of the preferences are persistent, that is, they remain effective even after closing and re-opening the program. To “remember” these choices, SplitsTree4 creates and maintains a *properties file*, which is placed in the users home directory and is called something like *.SplitsTree.def*.

It has five tabs.

The `Preferences:General` tab controls general aspects of the program:

Allow Graph Editing	if set, the user can add or delete nodes and edges from displayed network. To create a node, control-double click on the canvas. To create an edge between two nodes, control-click on a node and then drag create an edge. If dragging ends on a second node, then this is connected to the first, otherwise, a new node is created. To delete a node or delete an edge select it and press the delete key.
Maintain Edge Lengths	Usually, when interacting with a displayed network, we want the network to maintain edge lengths.
Show Scale Bar	The program uses a small <i>scale bar</i> in the top left hand corner of the Main:Network tab to indicate the scale of the network.
use Split-Selection Mode	When this feature is on, clicking on an edge will select all edges that correspond to the same split.

Any choices made here can either be applied to the current document or can be made the default for all subsequently opened documents.

The `Preferences:Defaults` tab controls the default sizes, shapes, colors and fonts for nodes and edges, see also `Nodes and Edges` .

The `Preferences:Layout` tab controls aspects of the layout of trees and split networks:

Recompute	The taxon layout is recomputed each time the Splits block is modified.
Stabilize	Use the union of the set of current splits and of the previously computed splits to compute a taxon layout. This is useful when one wants to compare two different networks computed on the same taxon set: by switching back and forth between the two computations, the program will find a joint taxon layout for both graphs, if one exists. In this case, both graphs will be laid out in such a way that in both graphs the taxa appear in approximately the same region of the canvas.
Snowball	Similar to the previous method, this method uses all splits ever computed in a given dataset to produce a taxon layout. This doesn't work very well.
Use this layout	Use the provided taxon layout to draw the network.

The `Preferences:Toolbar` tab allows the user to interactively configure the [tool bar](#).

The `Preferences:Status Line` tab allows the user to configure the *status line* displayed along the bottom of the [Main:Network](#) tab.

Here, two items are of particular interest. If the displayed graph was computed either by the [BunemanTree](#) or [SplitDecomposition](#) method, then the pair-wise distances in the graph may under estimate the true distances in the given [Distances](#) block. The difference between the two can be expressed in terms of the *fit* value, which is activated using the `Fit` check box, which is defined as the sum of all pairwise distances in the graph divided by the sum of all pairwise distances in the given matrix, times 100. This is **not** applicable to other tree or network building methods.

For other methods, please use the `LSFit` item, which computes the *least squares fit* between the pairwise distances in the graph and the pairwise distances in the matrix.

18 Additional Windows and Dialog

Here we list all other windows.

18.1 Open File

The `Open File` dialog is opened using the `File→Open` item. Use it to open any file containing phylogenetic data in one of the formats described in Section 20.

If an error is encountered, then the file is opened in the `MainSource` tab. If possible, the offending line is highlighted.

18.2 Choose Datatype

When opening a file containing character sequences, or importing sequences from the `Main:Source` tab, the program must know whether the sequences are to be interpreted as DNA, RNA, protein or “standard” (0,1) data. In `Nexus` files, the datatype is explicitly given. In other file formats this is usually not the case. The program employs a simple heuristic to guess the datatype. If this fails (e.g., if all character states are 'A'), then the program will display a `Choose Datatype` dialog and prompt the user to specify the datatype.

The choices are:

- `dna`,
- `rna`,
- `protein`,
- `standard`, which means 0,1 data, and
- `unknown`, which the program cannot deal with.

18.3 Save As

The `Save As` dialog is opened using the `File→Save As` item. Its purpose is to save the complete state of a document in `Nexus` format. To save parts of the document in `Nexus` format, or in some other supported format listed in Section 20, use the `Export` dialog.

18.4 Find/Replace Window

The **Find/Replace** window is opened using the **Edit→Find/Replace** item. Its purpose is to find and/or replace text. Such searches can be performed either in the **Main:Source** tab or in the **Main:Network** tab. In the latter case, find and replace is performed in all visible node and edge labels. Note that changing a node label only changes the label displayed in the network. The original taxon label provided in the **Taxa** block remains unchanged.

18.5 Node and Edge Window

The **Nodes and Edges** window is opened using the **View→Nodes and Edges** item. Its purpose is to modify the appearance of nodes and their labels, and edges and their labels, in the displayed network. This window has two tabs:

The **Nodes and Edges:Nodes** tab is used to control the appearance of nodes. It can be used to set node size, node shape, node color and node font. Moreover, it can be used to determine whether nodes are labeled by taxon names, taxon ids, or both.

The **Nodes and Edges:Edges** tab is used to control the appearance of edges. It can be used to set edge width, edge color, and edge font. Moreover, it can be used to determine whether edges are labeled by split weights, split ids, split confidences, or some combination. If the splits block contains confidence intervals for the edges (for example, in a confidence network) then these intervals can be displayed by selecting the intervals check box.

Please note that any changes made *only apply* to the currently selected nodes or edges. If there are no edges or nodes selected, then changes will apply to *all* nodes and edges. Any change made in this dialog box can be reversed using **Edit→Undo**.

18.6 Highlight Confidence Window

The **Highlight Confidence** Window is opened using the **View→Highlight Confidence** window item.

Use this window to request that edges and/or edge shading is done to reflect the confidence values associated with each split.

18.7 Bootstrap Window

The **Bootstrap** window is opened using the **Analysis→Bootstrap** item. Enter the Number of Replicates and then press Run to execute.

18.8 Message Window

The `Message` window is opened using the `Window→Message Window` item. The program writes all messages to this window.

18.9 About Window

The `About` Window is opened using the `Window→About` item. It reports the version of the program.

19 Nexus Blocks

In this section we describe the *Nexus* format, as implemented in `SplitsTree4`, based on the definition provided in [25]. Unfortunately, there exist two variants of the Nexus format, which we will call *old Nexus* and *new Nexus*. `SplitsTree4` is based on the latter, as this is what is defined in [25]. Given a file formatted in old Nexus, the program will often be able to parse it as it contains code to automatically convert from old to new Nexus format. However, the algorithm that does this does not provide a full implementation of the old Nexus format and thus sometimes it will be necessary to reformat a file by hand.

It is easy to tell the difference between old Nexus and new Nexus: if a file in Nexus format contains a `Taxa` block, then it is new Nexus. Most blocks in both formats have the same name and similar syntax. One main difference is that the `Characters` block in new Nexus is called a `Data` block in old Nexus.

In the following syntax descriptions, we used upper case letters for keywords, square brackets for optional statements and curly brackets to indicate a list of choices.

19.1 Taxa Block

The `Taxa` block is the only mandatory block in a Nexus file. Its purpose is to list the names of all taxa. It has the following syntax:

```
BEGIN TAXA;  
DIMENSIONS NTAX=number-of-taxa;  
[TAXLABELS taxon_1 taxon_2 ... taxon_ntax;]  
[TAXINFO info_1 info_2 ... info_ntax;]  
END;
```


The TAXLABELS statement is optional, if it is followed by a [source](#) block that contains all taxa labels.

19.2 Unaligned Block

The `Unaligned` block contains unaligned sequences. It has the following syntax:

```
BEGIN UNALIGNED;
[DIMENSIONS NTAX=number-of-taxa;]
[FORMAT
  [DATATYPE={STANDARD|DNA|RNA|NUCLEOTIDE|PROTEIN}]
  [RESPECTCASE]
  [MISSING=symbol]
  [SYMBOLS="symbol symbol ..."]
  [LABELS={LEFT|NO}]
;]
MATRIX
  [taxonlabel1] sequence ,
[taxonlabel2] sequence ,
  ...
[taxonlabelN] sequence
;
END;
```

19.3 Characters Block

The `Characters` block contains aligned character sequences. It has the following syntax:

```
BEGIN CHARACTERS;
DIMENSIONS [NTAX=number-of-taxa] NCHAR=number-of-characters;
[FORMAT
  [DATATYPE={STANDARD|DNA|RNA|PROTEIN}]
  [RESPECTCASE]
  [MISSING=symbol]
  [GAP=symbol]
  [SYMBOLS="symbol symbol ..."]
  [LABELS={NO|LEFT}]
  [TRANSDPOSE={NO|YES}]
  [INTERLEAVE={NO|YES}]
  [TOKENS=NO]
;]
```

```

[CHARWEIGHTS wgt_1 wgt_2 ... wgt_nchar;]
[CHARSTATELABELS character-number [character-name]
[ /state-name [ state-name... ] ], ...;]
MATRIX
    sequence data in specified format
;
END;

```

19.4 Distances Block

The `Distances` block contains a matrix of pairwise distances. It has the following syntax:

```

BEGIN DISTANCES;
[DIMENSIONS [NTAX=number-of-taxa];]
[FORMAT
    [TRIANGLE={LOWER|UPPER|BOTH}]
    [[NO] DIAGONAL]
    [LABELS={LEFT|NO}]
;]
MATRIX
    distance data in specified format
;
END;

```

19.5 Quartets Block

The `Quartets` block contains a list of quartets. It has the following syntax:

```

BEGIN Quartets;
DIMENSIONS [NTAX=number-of-taxa] NQUARTETS=number-of-quartets;
[FORMAT
    [LABELS={LEFT|NO}]
    [WEIGHTS={YES|NO}]
;]
MATRIX
[label1] [weight1] a1 b1 : c1 d1,
...
[labeln] [weightn] an bn : cn dn,
;
END;

```

19.6 Trees Block

The `Trees` block contains a list of phylogenetic trees. It has the following syntax:

```
BEGIN Trees
[PROPERTIES PARTIALTREES={YES|NO};]
[TRANSLATE
    nodeLabel1 taxon1,
    nodeLabel2 taxon2,
    ...
    nodeLabelN taxonN
;]
[TREE name1 = tree1-in-Newick-format;]
[TREE name2 = tree2-in-Newick-format;]
...
[TREE nameM = treeM-in-Newick-format;]
END;
```

19.7 Splits Block

The `splits` block contains a list of splits. It has the following syntax:

```
BEGIN Splits;
[DIMENSIONS [NTAX=number-of-taxa] [NSPLITS=number-of-splits];]
[FORMAT
    [LABELS={LEFT|NO}]
    [WEIGHTS={YES|NO}]
    [CONFIDENCES={YES|NO}]
    [INTERVALS={YES|NO}]
;]
[THRESHOLD=non-negative-number;]
[PROPERTIES
    [FIT=non-negative-number]
    [leastsquares]
    [{COMPATIBLE|CYCLIC|WEAKLY COMPATIBLE|INCOMPATIBLE}]
;]
[CYCLE [taxon_i_1 taxon_i_2 ... taxon_i_ntax];]
[SPLITSLABELS label_1 label_2 ... label_nsplits;]
MATRIX
    [label_1] [weight_1] [confidence_1] split_1,
    [label_2] [weight_2] [confidence_2] split_2,
    ....
```

```

        [label_nsplits] [weight_nsplits] [confidence_nsplits] split_nsplits,
;
END;

```

19.8 Network Block

The `Network` block contains the definition of a phylogenetic network. It has the following syntax:

```

BEGIN NETWORK;
DIMENSIONS NTAX=number-taxa NVERTICES=number-vertices NEDGES=number-edges;
[DRAW
    [ROTATE=rotation]
;]
[TRANSLATE
    [vertex_1 taxon_1,
    vertex_2 taxon_2,
    ...
    vertex_ntax taxon_ntax,]
;]
VERTICES
    1 x_1 y_1 [WIDTH=n] [HEIGHT=n] SHAPE=[RECT|OVAL] [FGC=color]
[BGC=color] [LINE=n],
    2 x_2 y_2 [WIDTH=n] [HEIGHT=n] SHAPE=[rect|OVAL] [FGC=color]
[BGC=color] [LINE=n],
    ...
    nvertices x_nvertices y_nvertices [WIDTH=n] [HEIGHT=n]
SHAPE=[RECT|OVAL] [FGC=color] [BGC=color] [LINE=n],
;
VLABELS
    vertex_id label [X= xoffset Y=yoffset] [FGC=color] [FONT=font],
    ...
    vertex_id label [X= xoffset Y=yoffset] [FGC=color] [FONT=font],
EDGES
    1 vertex_id vertex_id [ECLASS=n] [FGC=color] [LINE=n],
    2 vertex_id vertex_id [ECLASS=n] [FGC=color] [LINE=n],
    ...
    nedges vertex_id vertex_id [FGC=color] [LINE=n],
[ELABELS
    edge_id label [X= xoffset Y=yoffset] [FGC=color] [FONT=font],
    ...
    edge_id label [X= xoffset Y=yoffset] [FGC=color] [FONT=font],
;]
[INTERNAL
    edge_id [x y] [x y] ...,

```

```

    ...
    edge_id [x y] [x y] ...,
;]
END;

```

19.9 Bootstrap Block

The `Bootstrap` block contains the results of a bootstrap analysis. It has the following syntax:

```

BEGIN ST_BOOTSTRAP;
[DIMENSIONS [NTAX=number-of-taxa] [NCHAR=number-of-characters]
[NSPLITS=number-of-splits];]
[FORMAT
    [LABELS={LEFT|NO}]
    [SPLITS={NO|YES}]
    [ALL={YES|NO}]
;]
[RUNS=the-number-of-runs;]
[LENGTH={sample-length | SAME}];]
[SEED=random-number-seed;]
[SAVEWEIGHTS={yes|no};]
[FIXSPLITS={yes|no};]
[COMPUTEDIST={yes|no};]
[OUTPUTFILE=file-name;]
[MATRIX
    [label_1] value_1 [split_1,]
    [label_2] value_2 [split_2,]
    ....
    [label_nsplits] value_nsplits [split_nsplits,]
    [label_nsplits+1] value_(nsplits+1) [splits_(nsplits+1),]
    ....
    [label_n] value_n [splits_n,]
;]
END;

```

19.10 Sets Block

The `Sets` block can be used to define sets of taxa or characters. It has the following syntax:

```

BEGIN Sets;
[TAXSET taxset-name = taxon-list;]
[CHARSET charset-name = character-list;]

```

```
[CHARPARTITION charpart-name = 1:charset-name|character-list [...];
...
END;
```

19.11 ST Assumptions Block

The `ST_Assumptions` block controls the processing of the data along the [pipeline](#). It contains all choices made by the user that affect computations. It has the following syntax:

```
BEGIN ST_ASSUMPTIONS;
[UNALIGNTRANSFORM=name [parameters];]
[CHARTRANSFORM=name [parameters];]
[DISTTRANSFORM=name [parameters];]
[SPLITSTRANSFORM=name [parameters];]
[SPLITSPOSTPROCESS
[[NO] LEASTSQUARES]
[FILTER={GREEDYCOMPATIBLE|WEAKLYCOMPATIBLE|WEIGHT VALUE=value
|CONFIDENCE VALUE=value|DIMENSION VALUE=value|NONE};]
[EXTAXA={NONE|list-of-original-taxa-labels};]
[EXCHAR={NONE|list-of-original-char-positions};]
[EXCLUDE [[NO] GAPS] [[NO] NONPARSIMONY
[{NO CONSTANT|CONSTANT [number]}]
[[NO] CODON1] [[NO] CODON2] [[NO] CODON3];]
[EXTREES={NONE|list-of-original-tree-labels};]
[LAYOUTSTRATEGY={STABILIZE|SNOWBALL|KEEP};]
[[NO] AUTOLAYOUTNODELABELS];]
[UPTODATE;]
END;
```

20 File Formats

By default, `SplitsTree4` reads and writes data in [Nexus](#) format. The program can read and export the following additional formats: *FastA* , *Phylip* and *ClustalW* :

Name	file suffix	data type
new Nexus	.nex, .nxs	taxa, unaligned sequences, aligned sequences (characters), distances, quartets, trees, splits, networks
old Nexus	.nex, .nxs	aligned characters, distances, trees [25]
<i>FastA</i>	.fa, .fasta	unaligned sequences, or aligned characters
<i>Phylip</i>	.phy, .dst, .dist	aligned characters or distances [10]
<i>ClustalW</i>	.aln	aligned characters [32]

21 All Methods

Here we list all methods supported by `SplitsTree4`. We describe the usage and list the input and output [Nexus](#) blocks.

`Binary2Splits` : This method converts binary characters to splits.

Usage: `Binary2Splits AddAllTrivial=<boolean> MinSplitWeight=<int>`
Input: `Characters`
Output: `Splits`

`BioNJ` : This method computes the Bio-NJ tree [\[12\]](#).

Usage: `BioNJ`
Input: `Distances`
Output: `Trees`

`BunemanQuartets` : This method computes all quartets with positive Buneman index [\[6\]](#).

Usage: `BunemanQuartets threshold=<double>`
Input: `Distances`
Output: `Quartets`

`BunemanTree` : This method computes the Buneman tree [\[6\]](#).

Usage: `BunemanTree`
Input: `Distances`
Output: `Splits`

`ClustalW` : This method externally runs the *ClustalW sequence alignment* program [\[32\]](#).

Usage: ClustalW GapOpen=<int> GapExtension=<double>
WeightMatrix=<java.lang.String> OptionalParameter=<java.lang.String>
PathToCommand=<java.lang.String>

Input: Unaligned
Output: Characters

Coalescent : This method transforms a set of quartets to a set of splits representing a binary phylogenetic tree by applying the *coalescent method* described in [26].

Usage: Coalescent
Input: Quartets
Output: Splits

ConsensusNetwork : This method computes the consensus splits of trees [2, 15] to produce a *consensus network*.

Usage: ConsensusNetwork Threshold=<double> EdgeWeights=<String>
Input: Trees
Output: Splits

ConsensusTree : This method computes different types of *consensus trees*.

Usage: ConsensusTree EdgeWeights=<String> Method=<String>
Input: Trees
Output: Splits

ConvexHull : This method computes a splits graph using the convex hull extension algorithm [7].

Usage: ConvexHull Weights=<boolean> ScaleNodesMaxSize=<int>
Input: Splits
Output: Network

DNA2Splits : This method converts DNA characters to splits by setting the majority state against all other states.

Usage: DNA2Splits AddAllTrivial=<boolean> MinSplitWeight=<int>
Input: Characters
Output: Splits

DQuartets : This method compute all quartets with positive isolation index [1].

Usage: DQuartets threshold=<double>
Input: Distances
Output: Quartets

EqualAngle : This method computes a planar split network for a circular (sub-)set of splits [7]. If the `RunConvexHull` option is chosen, then the convex hull algorithm is subsequently applied to obtain a graph for the complete set of splits. This method provides a number of heuristics for obtaining a better layout [11]: set `DaylightIterations` to ≈ 5 to apply the *equal daylight* heuristic, set `OptimizeBoxesIterations` to ≈ 10 to apply the *box opening* heuristic and set `SpringEmbedderIterations` to ≈ 500 to apply a modified spring embedder.

Usage: `EqualAngle UseWeights=<boolean> RunConvexHull=<boolean>`
`DaylightIterations=<int> OptimizeBoxesIterations=<int>`
`SpringEmbedderIterations=<int>`
Input: `Splits`
Output: `Network`

FilteredSuperNetwork : This method computes a *super-network* from *partial trees* using the *Z-closure* algorithm [18] and a *distortion filter* [23].

Usage: `FilteredSuperNetwork MinNumberTrees=<int> MaxDistortionScore=<int> EdgeWeights=<String>`
`AllTrivial=<boolean> UseTotalScore=<boolean>`
Input: `Trees`
Output: `Splits`

F81 : This method calculates distances using the *Felsenstein-81* model [31].

Usage: `F81 Maximum_Likelihood=<boolean> Estimate_Base_Frequencies=<boolean>`
`Base_Freqs=<N double1 double2 ... doubleN> Normalize=<boolean>`
Input: `Characters`
Output: `Distances`

F84 : This method Calculates distances using the *Felsenstein-84* model [31].

Usage: `F84 Maximum_Likelihood=<boolean> Estimate_Base_Frequencies=<boolean>`
`Normalize=<boolean> TRatio=<double> A=<double> C=<double> G=<double>`
`T_U=<double>`
Input: `Characters`
Output: `Distances`

GapDist : This method calculates the “gap distance” from a set of sequences.

Usage: `GapDist`
Input: `Characters`
Output: `Distances`

GeneContentDistance : This method compute distances based on *gene content* [19].

Usage: GeneContentDistance UseMLDistance=<boolean>
Input: Characters
Output: Distances

Hamming : This method calculates distances using the Hamming distance. This is identical to the [UncorrectedP](#) method.

Usage: Hamming ignoregaps=<boolean>
Input: Characters
Output: Distances

HKY85 : This method calculates distances using the *Hasegawa, Kishino and Yano model*.

Usage: HKY85 Estimate_Base_Frequencies=<boolean> Normalize=<boolean>
TRatio=<double> A=<double> C=<double> G=<double> T_U=<double>
P_Invar=<double> Gamma=<double>
Input: Characters
Output: Distances

HybridizationNetwork : This method computes a *hybridization network* [21]. The present implementation operates by calling the [ReticulateNetwork](#) method.

Usage: HybridizationNetwork Which=<int> Method=<String> ShowSplits=<boolean>
OutGroup=<String> MaxAngle=<int> MaxReticulationsPerTangle=<int>
ShowSequences=<boolean> ShowMutations=<boolean>
Input: Splits
Output: Network

JukesCantor : This method computes distances using the *Jukes Cantor model* [31].

Usage: JukesCantor Maximum_Likelihood=<boolean>
Input: Characters
Output: Distances

K2P : This method calculates distances using the *Kimura-2P* model [31].

Usage: K2P Maximum_Likelihood=<boolean> TRatio=<double>
Input: Characters
Output: Distances

K3ST : This method calculates distances using the *Kimura-3ST* model [31].

Usage: K3ST Maximum_Likelihood=<boolean> TRatio=<double> AC_vs_ATRatio=<double>
Input: Characters
Output: Distances

LogDet : This method Calculates the LogDet-distance [30].

Usage: **LogDet** **Impute_Gaps**=<boolean>

Input: Characters

Output: Distances

LogHamming : This method calculates distances using the log-Hamming distance.

Usage: **LogHamming** **ignoregaps**=<boolean>

Input: Characters

Output: Distances

MedianNetwork : This method computes an (unreduced) *median network* [2]. It uses all sites in the character alignment that contain exactly two different states, and no gaps or missing states. If **UseRYAlphabet** is selected, then DNA and RNA sequences are translated using $R = \{A, G\}$ and $Y = \{C, T, U\}$. If **UseRelaxedSupport** is selected, a character need only be constant on one side of a split to count toward the support of the split.

Usage: **MedianNetwork** **AddAllTrivial**=<boolean> **MinimalSupport**=<int>

UseRYAlphabet=<boolean> **LabelEdges**=<boolean> **UseRelaxedSupport**=<boolean>

Input: Characters

Output: Splits

Muscle : This method externally runs the Muscle sequence alignment program [8].

Usage: **Muscle** **Maxiters**=<int> **ClusterMethod_1**=<String>

ClusterMethod_2=<String> **DistanceMeasure_1**=<String>

DistanceMeasure_2=<String> **ObjectiveScore**=<String>

LogFile=<boolean> **LogFileName**=<String>

OptionalParameter=<String> **PathToCommand**=<String>

Input: Unaligned

Output: Characters

NeiMiller : This method computes distances from restriction-sites using the Nei and Miller method [27].

Usage: **NeiMiller**

Input: Characters

Output: Distances

NJ : This method computes the Neighbour-Joining tree [28].

Usage: **NJ**

Input: Distances

Output: Trees

NeighborNet : This method computes the *Neighbor-Net* splits [5] to produce a Neighbor-Net network.

Usage: NeighborNet Variance=<String> Minimize_AIC=<boolean>
Input: Distances
Output: Splits

Noalign : This method obtains a trivial *sequence alignment* by adding gaps to the end of each sequence to make all sequences have the same length.

Usage: Noalign
Input: Unaligned
Output: Characters

NoGraph : This method prevents the program from constructing a final network.

Usage: NoGraph
Input: Splits
Output: Network

NoSplits : This method prevents the program from constructing splits from a set of trees.

Usage: NoSplits
Input: Trees
Output: Splits

ParsimonySplits : This method computes the set parsimony splits [1].

Usage: ParsimonySplits
Input: Characters
Output: Splits

PhylipParsimony : This method computes the *maximum parsimony* tree from DNA sequences using an external call to the Phylip package [10].

Usage: PhylipParsimony PhylipPath=<String> SearchMode=<String>
NumberOfTreesToSave=<int> InputOrderSeed=<int> InputOrderJumbles=<int>
OutgroupRoot=<String> ThresholdParsimony=<double>
TranversionParsimony=<boolean> WeightsFile=<String>
Input: Characters
Output: Trees

Phylogram : This method computes a traditional phylogenetic tree.

Usage: Phylogram Weights=<boolean> Cladogram=<boolean> Outgroup=<String>
UseOutgroup=<boolean>
Input: Splits
Output: Network

PhyML : This method calculates *maximum likelihood* trees from DNA sequences using PHYML [13].

Usage: PhyML PHYMLPath=<String> TreePath=<String> Bootstrap=<boolean>
NumberOfBootstrapReplicates=<int> PrintBootstrap=<boolean>
SubstitutionModel=<java.lang.String>
OptimiseEquilibriumFrequencies=<boolean>
EquilibriumFrequenciesEmpirical=<boolean> FrequencyA=<double>
FrequencyC=<double> FrequencyG=<double> FrequencyT=<double>
SubstitutionParameterAC=<int> SubstitutionParameterAG=<int>
SubstitutionParameterAT=<int> SubstitutionParameterCG=<int>
SubstitutionParameterCT=<int> SubstitutionParameterGT=<int>
EmpiricalBaseFrequencyEstimates=<boolean>
GammaDistributionParameter=<double>
GammaDistributionParameterFixed=<boolean>
InvariableSitesProportion=<double>
InvariableSitesProportionFixed=<boolean>
NumberOfSubstitutionCategories=<int> OneSubstitutionCategory=<boolean>
OptimiseStart=<boolean> OptimiseStartKeepingTopology=<boolean>
TstvRatio=<double> TstvRatioFixed=<boolean> UseBioNJstart=<boolean>
OptimiseRelativeRateParameters=<boolean>

Input: Characters

Output: Trees

ProteinMLdist : This method calculates maximum likelihood protein distance estimates using the following models: *cpREV45* , *Dayhoff* , *JTT* , *mtMAM* , *mtREV24* , *pmb* , *Rhodopsin* and *WAG* [31].

Usage: ProteinMLdist Gamma=<double> Model=<String> PInvar=<double>
Estimate_variances=<boolean>

Input: Characters

Output: Distances

PTreeSplits : This method computes the parsimony splits tree [1].

Usage: PTreeSplits

Input: Characters

Output: Splits

RecombinationNetwork : This method compute a *recombination network* from binary sequences [22]. If the given data is not 0,1 data, but DNA, then only those sites are considered that show precisely two different states. The present implementation operates by calling the [ReticulateNetwork](#) method.

Usage: RecombinationNetwork MinSplitWeight=<int> Which=<int> Method=<String>

ShowSplits=<boolean> OutGroup=<String> MaxAngle=<int>
 MaxReticulationsPerTangle=<int> ShowSequences=<boolean>
 ShowMutations=<boolean>
Input: Characters
Output: Network

ReticulateNetwork : This method compute a *reticulate network* from splits [22].

Usage: ReticulateNetwork MinSplitWeight=<int> Which=<int> Method=<String>
 ShowSplits=<boolean> OutGroup=<String> MaxAngle=<int>
 MaxReticulationsPerTangle=<int> ShowSequences=<boolean>
 ShowMutations=<boolean>
Input: Characters
Output: Network

RefinedBunemanTree : This method computes the Refined Buneman Tree [3]. This module was implemented by Lasse Westh-Nielsen and Christian N. S. Pedersen.

Usage: RefinedBunemanTree
Input: Distances
Output: Splits

ReticulateNetwork : This method computes a *reticulate network* [21].

Usage: ReticulateNetwork Which=<int> Method=<String> ShowSplits=<boolean>
 OutGroup=<String> MaxAngle=<int> MaxReticulationsPerTangle=<int>
 ShowSequences=<boolean> ShowMutations=<boolean>
Input: Splits
Output: Network

RootedEqualAngle : This method computes a rooted split network using the rooted equal angle algorithm [11].

Usage: RootedEqualAngle OptimizeDaylight=<boolean> UseWeights=<boolean>
 RunConvexHull=<boolean> DaylightIterations=<int> OutGroup=<String>
 MaxAngle=<int> SpecialSwitch=<boolean>
Input: Splits
Output: Network

RYSplits : This method computes all RY splits.

Usage: RYSplits
Input: Characters
Output: Splits

SpectralSplits : This method computes all splits arising using spectral analysis [14].

Usage: SpectralSplits Threshold=<double> Method=<String> Weight_ATvsGC=<double>
Weight_AGvsCT=<double> Weight_ACvsGT=<double>

Input: Characters

Output: Splits

SplitDecomposition : This method computes the *split decomposition* [1].

Usage: SplitDecomposition

Input: Distances

Output: Splits

SuperNetwork : This method computes a *super-network* from *partial trees* using the *Z-closure* algorithm [18].

Usage: SuperNetwork EdgeWeights=<String> ZRule=<boolean> NumberOfRuns=<int>
SuperTree=<boolean> ApplyRefineHeuristic=<boolean>

Input: Trees

Output: Splits

TreeSelector : This method is used to select a single tree from a set of trees.

Usage: TreeSelector Which=<int>

Input: Trees

Output: Splits

UncorrectedP : This method calculates uncorrected (observed, "P") distances. This is identical to the [Hamming](#) method.

Usage: Uncorrected_P ignoregaps=<boolean>

Input: Characters

Output: Distances

UPGMA : This method computes UPGMA tree [29].

Usage: UPGMA

Input: Distances

Output: Trees

22 Command-Line Options and Mode

SplitsTree4 has the following *command-line options* :

```

-g <switch> (default=true): GUI mode
-p <String> (default="$HOME/.SplitsTree.def"): Properties file
-i <String> (default=""): Input file
-x <String> (default=""): Execute this command at startup
-V <switch> (default=false): show version string
-S <switch> (default=false): silent mode
-d <switch> (default=false): debug mode
-s <switch> (default=true): show startup splash screen
-h <switch> (default=false): Show usage

```

Launching the program with option `-g` will make the program run in non-GUI *command-line mode*, first reading commands from a file supplied with the `-i` option, if any, then executing any command given with the `-x` option, and then finally reading additional commands from standard input.

Please be aware that the command-line version of the program uses the same [properties file](#) as the interactive version. So, any a [preferences](#) set using the interactive version of the program will also apply to the command-line version of the program. If this is not desired, then please use the `-p` option to supply a different properties file.

Commands provided to the program from within a file must be grouped together in a `SplitsTree` block so:

```

begin SplitsTree;
commands...
end;

```

Here we list all commands known to `SplitsTree4`:

```

EXECUTE FILE=file - open and execute a file in Nexus-format
OPEN FILE=file - open (but don't execute) a file in Nexus-format
IMPORT FILE=file [DATATYPE={PROTEIN|RNA|DNA|STANDARD|UNKNOWN}] - open (but don't execute) a file in
    old-Nexus format
LOAD FILE=file - open or import a file
LOAD TREEFILES=file1 ... fileN - load trees from one or more files
LOAD CHARFILES=file1 ... fileN - concatenate sequences from one or more files
LOAD FILE=file - open or import a file
SAVE FILE=file [REPLACE={YES|NO}] [APPEND={YES|NO}] [DATA={ALL|list-of-blocks}]
    - save all data or named blocks to a file in Nexus format
EXPORT FILE=file FORMAT=format [REPLACE={YES|NO}] [APPEND={YES|NO}]
    [DATA=list-of-blocks]
    - export data in the named formatAA
EXPORTGRAPHICS [format={EPS|PNG|GIF|JPG|SVG}] file=file [REPLACE={YES|NO}]
    [TEXTASSHAPES={YES|NO}] [TITLE=title]
    - export graphics in specified format (default is EPS)
UPDATE - rerun computations to bring data up-to-date
BOOTSTRAP RUNS=number-of-runs - perform bootstrapping on character data

```



```

DELETEEXCLUDED; - delete all sites from characters block that are
                    currently excluded
ASSUME assumption - set an assumption using any statement defined
                    in the ST_ASSUMPTIONS block
SHOW [DATA=list-of-blocks] - show the named data blocks
CYCLE {KEEP|cycle} - set the graph layout cycle to KEEP or to a given cycle
HELP - show this info
HELP DATA=list-of-blocks - show syntax of named blocks
HELP TRANSFORM=transform - show usage of a specific data transformation
VERSION - report version
ABOUT - show info on version and authors
QUIT - exit program

```

To begin or end a multi-line input, enter a backslash '\'

23 Examples

Example files are provided with the program. They are contained in the `examples` sub directory of the installation directory. The precise location of the installation directory depends upon your operating system.

24 Acknowledgments

We would like to thank Barry G. Hall, Pete Lockhart, David Morrison and Mike Steel for many helpful comments.

This product includes software developed by the Apache Software Foundation (<http://www.apache.org/>), namely the *batik* library for generating image files. It also contains *Jama*, a Java matrix package (<http://math.nist.gov/javanumerics/jama/>) and *MRJAdapter*, a Java package used to help construct user interfaces for the Apple Macintosh.

References

- [1] H.-J. Bandelt and A. W. M. Dress. A canonical decomposition theory for metrics on a finite set. *Advances in Mathematics*, 92:47–105, 1992.
- [2] H.-J. Bandelt, P. Forster, B. C. Sykes, and M. B. Richards. Mitochondrial portraits of human population using median networks. *Genetics*, 141:743–753, 1995.

- [3] G.S. Brodal, R. Fagerberg, A. Östlin, C.N.S. Pedersen, and S.S. Rao. Computing refined buneman trees in cubic time. *Lecture Notes in Computer Science*, 2812:259–270, 2003. Springer Verlag.
- [4] T. Bruen, H. Philippe, and D. Bryant. A quick and robust statistical test to detect the presence of recombination. *Genetics*, (in press), 2005.
- [5] D. Bryant and V. Moulton. NeighborNet: An agglomerative method for the construction of planar phylogenetic networks. In R. Guigó and D. Gusfield, editors, *Algorithms in Bioinformatics, WABI 2002*, volume LNCS 2452, pages 375–391, 2002.
- [6] P. Buneman. The recovery of trees from measures of dissimilarity. In F. R. Hodson, D. G. Kendall, and P. Tautu, editors, *Mathematics in the Archaeological and Historical Sciences*, pages 387–395. Edinburgh University Press, 1971.
- [7] A. W. M. Dress and D. H. Huson. Constructing splits graphs. *IEEE/ACM Transactions in Computational Biology and Bioinformatics*, 1(3):109–115, 2004.
- [8] R.C. Edgar. MUSCLE: multiple sequence alignment with high accuracy and high throughput. *Nucleic Acids Research*, 32(5):1792–97, 2004.
- [9] D.P. Faith. Conservation evaluation and phylogenetic diversity. *Biol. Conserv.*, 61:1–10, 1992.
- [10] J. Felsenstein. PHYLIP (phylogeny inference package) version 3.6. Distributed by the author. Department of Genome Sciences, University of Washington, Seattle, 2004.
- [11] P. Gambette and D.H. Huson. Improved layout of phylogenetic networks. Submitted, 2005.
- [12] O. Gascuel. BIONJ: An improved version of the NJ algorithm based on a simple model of sequence data. *Mol. Biol. Evol.*, 14:685–695, 1997.
- [13] S. Guindon and O. Gascuel. A simple, fast, and accurate algorithm to estimate large phylogenies by maximum likelihood. *Syst Biol.*, 52(5):696–704, 2003.
- [14] M.D. Hendy and D. Penny. Spectral analysis of phylogenetic data. *Journal of Classification*, 10:5–24, 1993.
- [15] B. Holland and V. Moulton. Consensus networks: A method for visualizing incompatibilities in collections of trees. In G. Benson and R. Page, editors, *Proceedings of “Workshop on Algorithms in Bioinformatics”*, volume 2812 of *LNBI*, pages 165–176. Springer, 2003.
- [16] D. H. Huson. SplitsTree: A program for analyzing and visualizing evolutionary data. *Bioinformatics*, 14(10):68–73, 1998.
- [17] D. H. Huson and D. Bryant. Application of phylogenetic networks in evolutionary studies. *Molecular Biology and Evolution*, 23(2):254–267, 2006. Software available from www.splitstree.org.
- [18] D. H. Huson, T. DeZulian, T. Klopper, and M. A. Steel. Phylogenetic super-networks from partial trees. *IEEE/ACM Transactions in Computational Biology and Bioinformatics*, 1(4):151–158, 2004.

- [19] D. H. Huson and M. Steel. Phylogenetic trees based on gene content. 20(13):2044–9, 2004.
- [20] D.H. Huson. Introduction to phylogenetic networks. Tutorial presented at ISMB, 2005.
- [21] D.H. Huson, T. Kloepper, P.J. Lockhart, and M.A. Steel. Reconstruction of reticulate networks from gene trees. In *Proceedings of the Ninth International Conference on Research in Computational Molecular Biology (RECOMB)*, 2005.
- [22] D.H. Huson and T.H. Kloepper. Computing recombination networks from binary sequences. To appear in: ECCB, 2005.
- [23] D.H. Huson, M.A. Steel, and J. Whitfield. Reducing distortion in phylogenetic networks. To appear in: WABI 2006, 2006.
- [24] P. J. Lockhart, M. A. Steel, and D. H. Huson. Invariable site models and their uses in phylogeny reconstruction. *Syst. Biol.*, 49(2):225–232, 2000.
- [25] D.R. Maddison, D.L. Swofford, and W.P. Maddison. NEXUS: an extendible file format for systematic information. *System. Bio.*, 46(4):590–621, 1997.
- [26] E. Mossel and M. Steel. A phase transition for a random cluster model on phylogenetic trees. Submitted, 2003.
- [27] M. Nei and J.C. Miller. A simple method for estimating average number of nucleotide substitutions within and between populations from restriction data. *Genetics*, 125:873–879, 1990.
- [28] N. Saitou and M. Nei. The Neighbor-Joining method: a new method for reconstructing phylogenetic trees. *Molecular Biology and Evolution*, 4:406–425, 1987.
- [29] R. R. Sokal and C. D. Michener. A statistical method for evaluating systematic relationships. *University of Kansas Scientific Bulletin*, 28:1409–1438, 1958.
- [30] M.A. Steel. Recovering a tree from the leaf colorations it generates under a markov model. *Appl. Math. Lett.*, 7(2):19–24, 1994.
- [31] D. L. Swofford, G. J. Olsen, P. J. Waddell, and D. M. Hillis. Chapter 11: Phylogenetic inference. In D. M. Hillis, C. Moritz, and B. K. Mable, editors, *Molecular Systematics*, pages 407–514. Sinauer Associates, Inc., 2nd edition, 1996.
- [32] J.D. Thompson, D.G. Higgins, and T.J. Gibson. CLUSTAL W: improving the sensitivity of progressive multiple sequence alignment through sequence weighting, position-specific gap penalties and weight matrix choice. *Nucl. Acids Res.*, 22:4673–4680, 1994.

Index

- .SplitsTree.def, [28](#)
- .aln, [39](#)
- .dist, [39](#)
- .dst, [39](#)
- .fa, [39](#)
- .fasta, [39](#)
- .nex, [39](#)
- .nxs, [39](#)
- .phy, [39](#)
- [, [27](#)
- About, [22](#), [32](#)
- About command, [48](#)
- All, [16](#)
- allopolyploidization, [9](#)
- Allow Graph Editing, [29](#)
- Analysis, [21](#)
- Analysis→Bootstrap, [21](#), [31](#)
- Analysis→Compute Phylogenetic Diversity, [21](#)
- Analysis→Conduct Phi test for Recombination, [21](#)
- Analysis→Configure Pipeline, [5](#), [11](#), [21](#), [24](#)
- Analysis→Configure Recent Methods, [5](#), [11](#), [21](#), [23](#)
- Analysis→Estimate Invariable Sites, [21](#)
- Analysis→Show Bootstrap Network, [21](#)
- Analysis→Show Confidence Network, [21](#)
- Assume command, [48](#)
- Auto Layout Node Labels, [14](#), [17](#)
- batik, [49](#)
- binary, [9](#)
- Binary2Splits, [39](#)
- BIONJ, [20](#)
- BioNJ, [39](#)
- blocks, [6](#)
- BMP, [28](#)
- Bootstrap, [21](#), [31](#), [37](#)
- Bootstrap command, [48](#)
- box opening, [41](#)
- branch, [8](#)
- BunemanQuartets, [39](#)
- BunemanTree, [20](#), [39](#)
- Characters, [17](#), [33](#)
- Choose Datatype, [15](#), [30](#)
- circular, [9](#)
- Clear All Taxa Sets..., [16](#)
- Clone, [24](#)
- Clones, [15](#)
- Close, [15](#)
- Closest Tree, [27](#)
- ClustalW, [38](#), [39](#)
- Coalescent, [40](#)
- coalescent method, [40](#)
- Collapse All, [17](#)
- Command Syntax, [22](#)
- command-line mode, [48](#)
- command-line options, [47](#)
- compatible, [8](#)
- Compute Phylogenetic Diversity, [21](#)
- Concatenate Sequences, [15](#)
- concatenate sequences, [15](#)
- Conduct Phi test for Recombination, [21](#)
- Confidence Threshold, [27](#)
- Configure, [23](#), [24](#)
- Configure Pipeline, [5](#), [11](#), [21](#), [24](#)
- Configure Recent Methods, [5](#), [11](#), [21](#), [23](#)
- consensus network, [12](#), [40](#)
- consensus trees, [40](#)
- ConsensusNetwork, [20](#), [40](#)
- ConsensusTree, [11](#), [20](#), [40](#)
- constant, [18](#)
- Convert text to graphics, [28](#)
- Convert to Nexus, [18](#)
- ConvexHull, [22](#), [40](#)
- Copy, [16](#)
- Copy Label, [23](#)
- cpREV45, [45](#)
- create a node, [29](#)
- create an edge, [29](#)
- Cut, [16](#)
- Data, [17](#), [32](#)
- data type, [15](#)
- Data→Convert to Nexus, [18](#)
- Data→Exclude Constant Sites, [18](#)
- Data→Exclude Gap Sites, [17](#)

- Data→Exclude Parsimony-Uninformative Sites, [18](#)
- Data→Exclude Selected Splits, [18](#), [22](#)
- Data→Exclude Selected Taxa, [17](#)
- Data→Execute, [18](#)
- Data→Filter Characters, [18](#)
- Data→Filter Splits, [18](#)
- Data→Filter Taxa, [17](#)
- Data→Filter Trees, [18](#)
- Data→Greedy Make Compatible, [18](#)
- Data→Greedy Make Weakly Compatible, [18](#)
- Data→Keep Only Selected Taxa, [17](#)
- Data→Load, [19](#)
- Data→Restore All Sites, [18](#)
- Data→Restore All Splits, [18](#)
- Data→Restore All Taxa, [17](#)
- Data→Set Tree Names, [18](#)
- datatype, [15](#), [30](#)
- Dayhoff, [45](#)
- DaylightIterations, [41](#)
- default calculations, [7](#)
- Delete, [25](#)
- delete a node, [29](#)
- delete an edge, [29](#)
- Delete excluded command, [48](#)
- diploid, [9](#)
- Disclaimer, [4](#)
- display of sites, [26](#)
- Distances, [17](#), [19](#), [34](#)
- Distances→F81, [19](#)
- Distances→F84, [19](#)
- Distances→GeneContentDistance, [19](#)
- Distances→HKY85, [19](#)
- Distances→JukesCantor, [19](#)
- Distances→K2P, [19](#)
- Distances→K3ST, [19](#)
- Distances→LogDet, [19](#)
- Distances→NeiMiller, [19](#)
- Distances→ProteinMLdist, [19](#)
- Distances→UncorrectedP, [19](#)
- distortion filter, [41](#)
- DNA2Splits, [40](#)
- document, [6](#)
- Don't show this dialog for this method again, [23](#)
- DQuartets, [40](#)
- Draw, [21](#)
- Draw→ConvexHull, [22](#)
- Draw→EqualAngle, [22](#), [24](#)
- Draw→Hide Incompatible Splits, [22](#)
- Draw→Hide Non-Selected Splits, [22](#)
- Draw→Hide Selected Splits, [18](#), [22](#)
- Draw→HideSelectedSplits, [12](#)
- Draw→NoGraph, [22](#)
- Draw→Phylogram, [22](#), [24](#)
- Draw→Redraw All Splits, [22](#)
- Draw→RootedEqualAngle, [22](#), [24](#)
- Draw→Select Characters, [22](#)
- Draw→Select Trees, [12](#), [22](#)
- Edge, [23](#)
- edge, [8](#)
- edge color, [31](#)
- edge font, [31](#)
- edge width, [31](#)
- Edge→Configure, [24](#)
- Edge→Copy Label, [23](#)
- Edge→Edit Label, [23](#)
- Edge→Hide Label, [24](#)
- Edge→Show Confidence, [24](#)
- Edge→Show Id, [23](#)
- Edge→Show Interval, [24](#)
- Edge→Show Weight, [24](#)
- Edit, [16](#)
- Edit Label, [23](#)
- Edit→Copy, [16](#)
- Edit→Cut, [16](#)
- Edit→Find/Replace, [16](#), [24](#), [31](#)
- Edit→Go to Line, [16](#)
- Edit→Invert Selection, [16](#)
- Edit→Paste, [16](#)
- Edit→Preferences, [16](#), [24](#), [28](#)
- Edit→Redo, [16](#)
- Edit→Select All, [16](#)
- Edit→Taxon Sets, [16](#)
- Edit→TaxonSets→All, [16](#)
- Edit→TaxonSets→Clear All Taxa Sets..., [16](#)
- Edit→TaxonSets→New taxa set..., [16](#)
- Edit→Undo, [16](#), [31](#)
- editing trees, [28](#)
- Enter a Command, [22](#)
- EPS, [28](#)

EPS command, [48](#)
 equal daylight, [41](#)
 EqualAngle, [22](#), [24](#), [41](#)
 Estimate Invariable Sites, [21](#)
 examples, [5](#), [49](#)
 Exclude Constant Sites, [18](#)
 Exclude Gap Sites, [17](#)
 Exclude Parsimony-Uninformative Sites, [18](#)
 Exclude Selected Splits, [18](#), [22](#), [27](#)
 Exclude Selected Taxa, [17](#), [23](#)
 Execute, [18](#)
 execute, [18](#)
 Expand All, [17](#)
 Export, [5](#), [10](#), [15](#), [27](#)
 Export command, [48](#)
 Export Image, [5](#), [10](#), [15](#), [24](#), [28](#)

 F81, [19](#), [41](#)
 F84, [19](#), [41](#)
 FastA, [38](#), [39](#)
 Felsenstein-81, [41](#)
 Felsenstein-84, [41](#)
 File, [14](#)
 File command, [48](#)
 File→Clone, [24](#)
 File→Clones, [15](#)
 File→Close, [15](#)
 File→Export, [5](#), [10](#), [15](#), [27](#)
 File→Export Image, [5](#), [10](#), [15](#), [24](#), [28](#)
 File→New, [14](#)
 File→Open, [5](#), [10](#), [15](#), [19](#), [24](#), [30](#)
 File→Open Recent, [10](#), [15](#)
 File→Print, [15](#), [24](#)
 File→Quit, [15](#)
 File→Replace, [15](#)
 File→Save, [7](#), [10](#), [15](#)
 File→Save As, [5](#), [7](#), [10](#), [15](#), [24](#), [30](#)
 File→Tools, [15](#)
 File→Tools→Concatenate Sequences, [15](#)
 File→Tools→Group Identical Haplotypes, [15](#)
 File→Tools→Load Multi-Labeled Tree, [15](#)
 File→Tools→Load Trees, [15](#)
 Filter Characters, [18](#)
 Filter Splits, [18](#)
 Filter Taxa, [17](#)
 Filter Trees, [18](#)

 FilteredSuperNetwork, [20](#), [41](#)
 Find/Replace, [16](#), [24](#), [31](#)
 Fit, [29](#)
 fit, [29](#)
 Flip, [17](#)
 From Graph, [25](#)

 gap, [17](#)
 GapDist, [41](#)
 gene content, [41](#)
 GeneContentDistance, [19](#), [41](#)
 GIF, [28](#)
 Go to Line, [16](#)
 graphical attributes, [17](#)
 Greedily Make Compatible, [18](#)
 Greedily Make Weakly Compatible, [18](#)
 Greedy Compatible, [27](#)
 Greedy Weakly Compatible, [27](#)
 Group Identical Haplotypes, [15](#)

 Hamming, [42](#)
 Hasegawa, Kishino and Yano model, [42](#)
 Help command, [48](#)
 Hide All, [25](#)
 Hide Incompatible Splits, [22](#)
 Hide Label, [23](#), [24](#)
 hide list, [25](#)
 Hide Non-Selected Splits, [22](#)
 Hide Selected Splits, [18](#), [22](#)
 HideSelectedSplits, [12](#)
 Highlight Confidence, [17](#), [31](#)
 HKY85, [19](#), [42](#)
 How to Cite, [22](#)
 How to cite, [4](#)
 hybridization, [9](#)
 hybridization network, [9](#), [12](#), [42](#)
 HybridizationNetwork, [21](#), [42](#)

 import, [18](#)
 Import command, [48](#)
 initial calculations, [19](#)
 intervals, [31](#)
 Invert Selection, [14](#), [16](#)

 Jama, [49](#)
 JPEG, [28](#)
 JTT, [45](#)

Jukes Cantor model, [42](#)
 JukesCantor, [19](#), [42](#)

 K2P, [19](#), [42](#)
 K3ST, [19](#), [42](#)
 Keep Only Selected Taxa, [17](#)
 Kimura-2P, [42](#)
 Kimura-3ST, [42](#)

 layout, [29](#)
 least squares, [26](#)
 least squares fit, [30](#)
 Linux, [5](#), [6](#)
 Load, [19](#)
 Load command, [48](#)
 Load Multi-Labeled Tree, [15](#)
 Load Trees, [15](#)
 LogDet, [19](#), [43](#)
 LogHamming, [43](#)
 LSFit, [30](#)

 Mac OS, [5](#)
 MacOS, [6](#)
 Main, [12](#)
 Main:Data, [13](#)
 Main:Network, [13](#)
 Main:Source, [13](#)
 Maintain Edge Lengths, [29](#)
 maintain edge lengths, [29](#)
 maximum likelihood, [45](#)
 maximum parsimony, [44](#)
 median network, [43](#)
 MedianNetwork, [21](#), [43](#)
 merge a set of trees, [15](#)
 Message, [32](#)
 Message Window, [23](#), [24](#), [32](#)
 missing character, [17](#)
 MRJAdapter, [49](#)
 mtMAM, [45](#)
 mtREV24, [45](#)
 multi-threading, [6](#)
 multiple documents, [6](#)
 Muscle, [43](#)

 Neighbor-Net, [44](#)
 NeighborJoining, [23](#)
 NeighborNet, [20](#), [44](#)

NeiMiller, [19](#), [43](#)
 Network, [36](#)
 network, [8](#)
 Networks, [20](#)
 Networks→ConsensusNetwork, [20](#)
 Networks→FilteredSuperNetwork, [20](#)
 Networks→HybridizationNetwork, [21](#)
 Networks→MedianNetwork, [21](#)
 Networks→NeighborNet, [20](#)
 Networks→ParsimonySplits, [20](#)
 Networks→RecombinationNetwork, [21](#)
 Networks→SpectralSplits, [21](#)
 Networks→SplitDecomposition, [20](#)
 Networks→SuperNetwork, [20](#)
 New, [14](#)
 new Nexus, [32](#)
 New taxa set..., [16](#)
 Next Tree, [11](#), [20](#), [24](#)
 Nexus, [32](#)
 Nexus Syntax, [22](#)
 NJ, [20](#), [43](#)
 Noalign, [44](#)
 Node, [23](#)
 node color, [31](#)
 node font, [31](#)
 node shape, [31](#)
 node size, [31](#)
 Node→Configure, [23](#)
 Node→Copy Label, [23](#)
 Node→Edit Label, [23](#)
 Node→Exclude Selected Taxa, [23](#)
 Node→Hide Label, [23](#)
 Node→Show Id, [23](#)
 Node→Show Name, [23](#)
 Nodes and Edges, [5](#), [14](#), [17](#), [29](#), [31](#)
 Nodes and Edges:Edges, [31](#)
 Nodes and Edges:Nodes, [31](#)
 NoGraph, [22](#), [44](#)
 non-trivial splits, [8](#)
 None, [27](#)
 NoSplits, [44](#)
 Number of Replicates, [31](#)

 old Nexus, [32](#)
 Open, [5](#), [10](#), [15](#), [19](#), [24](#), [30](#)
 Open File, [30](#)

- Open Recent, [10](#), [15](#)
- OptimizeBoxesIterations, [41](#)
- p, [23](#)
- parallel branches, [9](#)
- parallel edges, [9](#)
- parsimony-uninformative, [18](#)
- ParsimonySplits, [20](#), [44](#)
- partial trees, [12](#), [41](#), [47](#)
- Paste, [16](#)
- Phylip, [38](#), [39](#), [44](#)
- PhylipParsimony, [20](#), [44](#)
- Phylogram, [22](#), [24](#), [44](#)
- PHYML, [45](#)
- PhyML, [20](#), [45](#)
- PhyML Path, [11](#)
- Pipeline, [24](#)
- pipeline, [6](#)
- Pipeline:Characters, [25](#)
- Pipeline:Characters:Filter, [25](#)
- Pipeline:Characters:Method, [25](#)
- Pipeline:Characters:Select, [26](#)
- Pipeline:Distances, [26](#)
- Pipeline:Distances:Method, [26](#)
- Pipeline:Quartets, [26](#)
- Pipeline:Quartets:Method, [26](#)
- Pipeline:Splits, [26](#)
- Pipeline:Splits:Filter, [26](#)
- Pipeline:Splits:Method, [26](#)
- Pipeline:Taxa, [25](#)
- Pipeline:Taxa:Filter, [25](#)
- Pipeline:Trees, [26](#)
- Pipeline:Trees:Filter, [26](#)
- Pipeline:Trees:Method, [26](#)
- Pipeline:Trees:Select, [26](#)
- Pipeline:Unaligned, [25](#)
- Pipeline:Unaligned:Method, [25](#)
- pmb, [45](#)
- PNG, [28](#)
- Preferences, [16](#), [24](#), [28](#)
- preferences, [28](#)
- Preferences:Defaults, [29](#)
- Preferences:General, [28](#)
- Preferences:Layout, [29](#)
- Preferences:Status Line, [29](#)
- Preferences:Toolbar, [13](#), [29](#)

- Previous Tree, [11](#), [20](#), [24](#)
- Print, [15](#), [24](#)
- properties file, [28](#)
- ProteinMLdist, [19](#), [45](#)
- PTreeSplits, [45](#)
- Quartets, [34](#)
- Quit, [15](#)
- Quit command, [48](#)
- recombination, [9](#)
- recombination network, [9](#), [12](#), [45](#)
- RecombinationNetwork, [21](#), [45](#)
- Redo, [16](#)
- Redraw All Splits, [22](#)
- RefinedBunemanTree, [20](#), [46](#)
- Replace, [15](#)
- Reset, [16](#), [17](#), [24](#)
- Reset Label Positions, [24](#)
- Restore All Sites, [18](#)
- Restore All Splits, [18](#), [24](#)
- Restore All Taxa, [17](#), [24](#)
- restriction-sites, [43](#)
- reticulate evolution, [9](#)
- reticulate network, [9](#), [46](#)
- reticulate networks, [12](#)
- ReticulateNetwork, [46](#)
- Rhodopsin, [45](#)
- RootedEqualAngle, [22](#), [24](#), [46](#)
- Rotate Left, [17](#), [24](#)
- Rotate Right, [17](#), [24](#)
- Run, [31](#)
- RunConvexHull, [41](#)
- RYSplits, [46](#)
- sampling error, [4](#)
- Save, [7](#), [10](#), [15](#)
- Save As, [5](#), [7](#), [10](#), [15](#), [24](#), [30](#)
- Save command, [48](#)
- Save visible image, [28](#)
- Save whole image, [28](#)
- scale bar, [29](#)
- Select All, [16](#)
- Select Characters, [22](#)
- Select Trees, [12](#), [22](#)
- sequence alignment, [39](#), [43](#), [44](#)

- Set Maximum Dimension, [27](#)
- Set Tree Names, [18](#)
- Sets, [37](#)
- Show All, [25](#)
- Show Bootstrap Network, [21](#)
- Show Confidence, [24](#)
- Show Confidence Network, [21](#)
- Show cycle command, [48](#)
- Show Id, [23](#)
- Show Interval, [24](#)
- show list, [25](#)
- Show Name, [23](#)
- Show Scale Bar, [29](#)
- Show Weight, [24](#)
- source block, [7](#)
- source text area, [18](#)
- SpectralSplits, [21](#), [46](#)
- split, [8](#)
- split confidences, [31](#)
- split decomposition, [47](#)
- split encoding, [8](#)
- split ids, [31](#)
- split network, [9](#)
- split weights, [31](#)
- Split-Selection Mode, [14](#)
- SplitDecomposition, [20](#), [47](#)
- Splits, [17](#)
- splits, [35](#)
- SplitsTree, [48](#)
- splitstree_linux_4.8.rpm, [5](#)
- splitstree_macos_4.8.sit, [5](#)
- splitstree_unix_4.8.sh, [5](#)
- splitstree_windows_4.8.exe, [5](#)
- SpringEmbedderIterations, [41](#)
- ST_Assumptions, [38](#)
- status line, [29](#)
- super network, [12](#)
- super-network, [41](#), [47](#)
- SuperNetwork, [20](#), [47](#)
- SVG, [28](#)
- synchronized, [7](#)
- systematic error, [4](#)
- Taxa, [32](#)
- TAXLABELS, [33](#)
- taxon ids, [31](#)
- taxon names, [31](#)
- Taxon Sets, [16](#)
- tool bar, [13](#), [24](#)
- Tools, [15](#)
- tree, [7](#)
- tree names, [18](#)
- Trees, [19](#), [35](#)
- Trees→BIONJ, [20](#)
- Trees→BunemanTree, [20](#)
- Trees→ConsensusTree, [11](#), [20](#)
- Trees→NeighborJoining, [23](#)
- Trees→Next Tree, [11](#), [20](#), [24](#)
- Trees→NJ, [20](#)
- Trees→PhylipParsimony, [20](#)
- Trees→PhyML, [20](#)
- Trees→Previous Tree, [11](#), [20](#), [24](#)
- Trees→RefinedBunemanTree, [20](#)
- Trees→TreeSelector, [20](#)
- Trees→TreesSelector, [11](#), [20](#)
- Trees→UPGMA, [20](#)
- TreeSelector, [20](#), [47](#)
- TreesSelector, [11](#), [20](#)
- trivial splits, [8](#)
- Type-setting conventions, [4](#)
- Unaligned, [33](#)
- UncorrectedP, [19](#), [47](#)
- Undo, [16](#), [31](#)
- Unix, [5](#), [6](#)
- Update command, [48](#)
- UPGMA, [20](#), [47](#)
- use Split-Selection Mode, [29](#)
- Version command, [48](#)
- View, [16](#)
- View→Auto Layout Node Labels, [14](#), [17](#)
- View→Characters, [17](#)
- View→Collapse All, [17](#)
- View→Distances, [17](#)
- View→Expand All, [17](#)
- View→Flip, [17](#)
- View→Highlight Confidence, [17](#), [31](#)
- View→Invert Selection, [14](#)
- View→Nodes and Edges, [5](#), [14](#), [17](#), [31](#)
- View→Reset, [16](#), [17](#), [24](#)
- View→Rotate Left, [17](#), [24](#)

- View→Rotate Right, [17](#), [24](#)
- View→Splits, [17](#)
- View→Zoom In, [16](#), [24](#)
- View→Zoom Out, [17](#), [24](#)

- WAG, [45](#)
- weakly compatible, [9](#)
- Weight Threshold, [27](#)
- Window, [22](#), [24](#)
- Window→About, [22](#), [32](#)
- Window→Command Syntax, [22](#)
- Window→Enter a Command, [22](#)
- Window→How to Cite, [22](#)
- Window→Message Window, [23](#), [24](#), [32](#)
- Window→Nexus Syntax, [22](#)
- Window→Reset, [24](#)
- Window→Reset Label Positions, [24](#)
- Window→Restore All Splits, [24](#)
- Window→Restore All Taxa, [24](#)
- Window→Zoom In, [24](#)
- Window→Zoom Out, [24](#)
- Windows, [5](#), [6](#)

- Z-closure, [41](#), [47](#)
- Zoom In, [16](#), [24](#)
- Zoom Out, [17](#), [24](#)