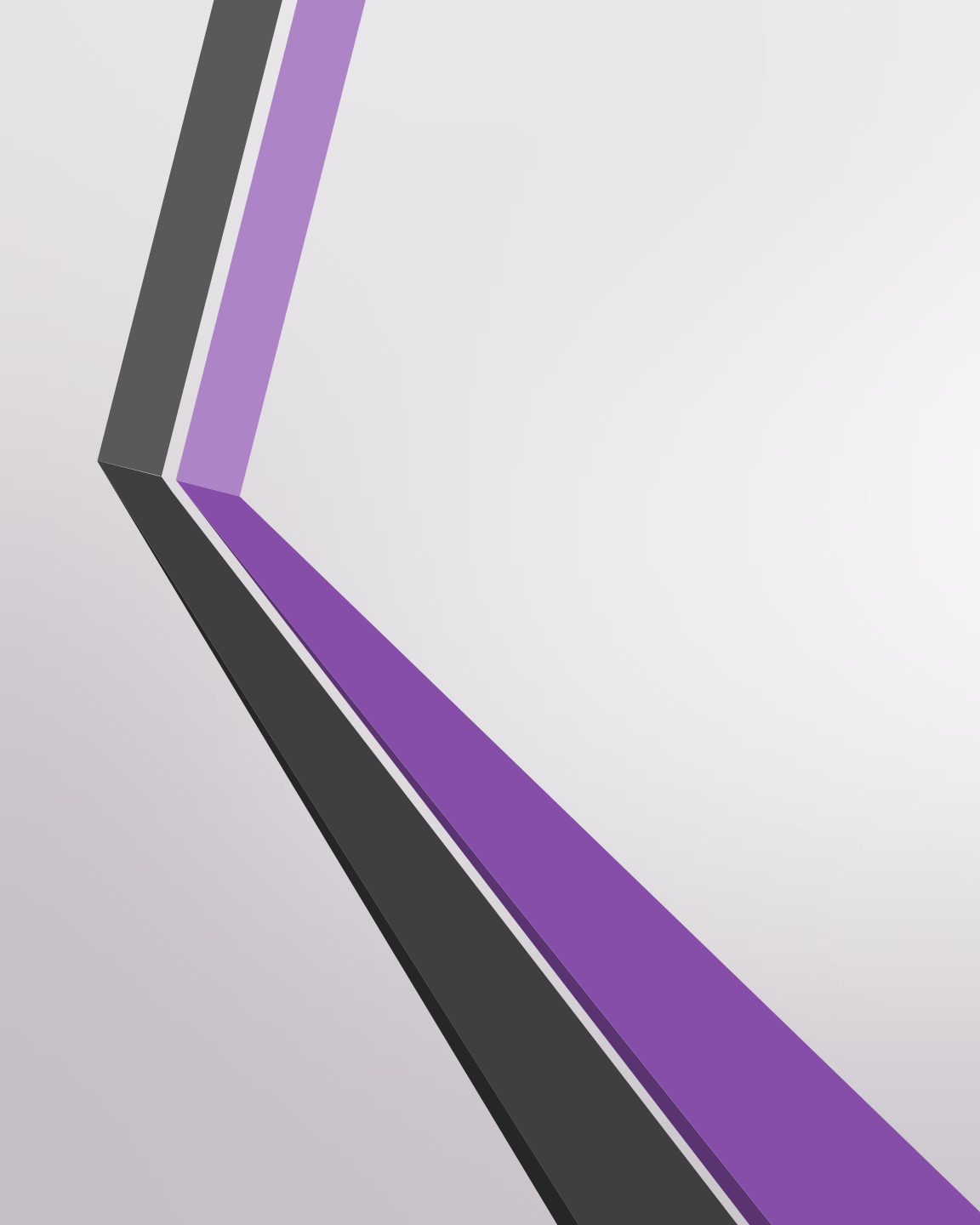




FastTree

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Instalación

- Se accede a: <http://www.microbesonline.org/fasttree/#Install>
- Windows y Linux: descargar ejecutable.
- Mac OS X: descargar fichero fuente FastTree.c (es necesario tener instalado xcode.gcc) y compilar con

`gcc -o3 -finline-functions -funroll-loops -Wall -o FastTree FastTree.c -lm`

FastTree

Also see [Fast Tree-Comparison Tools](#)

FastTree infers approximately-maximum-likelihood phylogenetic trees from alignments of nucleotide or protein sequences. FastTree can handle alignments with up to a million of sequences in a reasonable amount of time and memory. For large alignments, FastTree is 100-1,000 times faster than [PhyML 3.0](#) or [RAxML 7](#). FastTree is open-source software -- you can download the code [below](#).

FastTree is more accurate than PhyML 3 with default settings, and much more accurate than the distance-matrix methods that are traditionally used for large alignments. FastTree uses the Jukes-Cantor or [generalized time-reversible](#) (GTR) models of nucleotide evolution and the JTT ([Jones-Taylor-Thornton 1992](#)), WAG ([Whelan & Goldman 2001](#)), or LG ([Le and Gascuel 2008](#)) models of amino acid evolution. To account for the varying rates of evolution across sites, FastTree uses a single rate for each site (the "CAT" approximation). To quickly estimate the reliability of each split in the tree, FastTree computes [local support values](#) with the [Shimodaira-Hasegawa test](#) (these are the same as PhyML 3's "SH-like local supports").

More information:

- [How it Works](#)
- Downloads:
 - [Install FastTree](#)
 - [16S ribosomal RNA trees and alignments](#)
- [Running FastTree](#)
- [Speed & Accuracy](#)
 - [Simulated alignments and trees](#)
- [Special features:](#)
 - [Constrained topology search](#)
 - [Pseudocounts for highly fragmentary sequences](#)
 - [Multi-threaded FastTree with OpenMP](#)
 - [CAT/Gamma20 branch lengths and likelihoods](#)
- [Frequently Asked Questions](#)
- [Contact us](#)

Downloading and Installing FastTree

FastTree 2.1.10 is available as

- [Linux 64-bit executable \(+SSE\)](#)
 - [Double-precision executable](#) for nearly-identical sequences (see [explanation](#))
 - [Multi-threaded executable \(+SSE +OpenMP\)](#) (see [usage guide](#))
- [Windows 32-bit command-line executable \(no SSE\)](#)
- [C code](#)

If you use a Mac or other platform not included above, download [FastTree.c](#) and run

```
gcc -O3 -finline-functions -funroll-loops -Wall -o FastTree FastTree.c -lm
```

Ejecución

- Se ejecuta con: FastTree [parámetros] archivo alineado > archivo de salida
- Acepta ficheros en formato FASTA o PHYLIP.
- Parámetros:
 - **-nt**: para nucleótidos.
 - **-quiet**: suprime mensajes de salida.
 - **-log**: guarda árboles intermedios y detalles del modelo.
 - **-n <number>**: analiza múltiples alineamientos (solo formato PHYLIP).
 - **-intrees in_tree**: establece el árbol de partida.
 - **-quote**: obtiene número entero de cada secuencia.
 - **-gamma**: reescala ramas del árbol y calcula la probabilidad gammazo.
 - **-expert**: muestra más información.

Usage for FastTree version 2.1.10 No SSE3:

```
FastTree protein_alignment > tree
FastTree < protein_alignment > tree
FastTree -out tree protein_alignment
FastTree -nt nucleotide_alignment > tree
FastTree -nt -gtr < nucleotide_alignment > tree
FastTree < nucleotide_alignment > tree
```

FastTree accepts alignments in fasta or phylip interleaved formats

Common options (must be before the alignment file):

```
-quiet to suppress reporting information
-nopr to suppress progress indicator
-log logfile -- save intermediate trees, settings, and model details
-fastest -- speed up the neighbor joining phase & reduce memory usage
            (recommended for >50,000 sequences)
-n <number> to analyze multiple alignments (phylip format only)
            (use for global bootstrap, with seqboot and CompareToBootstrap.pl)
-nosupport to not compute support values
-intree newick_file to set the starting tree(s)
-intree1 newick_file to use this starting tree for all the alignments
            (for faster global bootstrap on huge alignments)
-pseudo to use pseudocounts (recommended for highly gapped sequences)
-gtr -- generalized time-reversible model (nucleotide alignments only)
-lg -- Le-Gascuel 2008 model (amino acid alignments only)
-wag -- Whelan-And-Goldman 2001 model (amino acid alignments only)
-quote -- allow spaces and other restricted characters (but not ' ) in
         sequence names and quote names in the output tree (fasta input only;
         FastTree will not be able to read these trees back in)
-noml to turn off maximum-likelihood
-nome to turn off minimum-evolution NNIs and SPRs
            (recommended if running additional ML NNIs with -intree)
-nome -mlen with -intree to optimize branch lengths for a fixed topology
-cat # to specify the number of rate categories of sites (default 20)
      or -nocat to use constant rates
-gamma -- after optimizing the tree under the CAT approximation,
          rescale the lengths to optimize the Gamma20 likelihood
-constraints constraintAlignment to constrain the topology search
          constraintAlignment should have 1s or 0s to indicates splits
-expert -- see more options
```

For more information, see <http://www.microbesonline.org/fasttree/>

Windows users: Please remember to run this inside a command shell

Hit return to continue

Modelos evolutivos

- Nucleótidos
 - Jukes-Cantor
 - GTR: usando -gtr
- Aminoácidos:
 - JTT
 - WAG: usando -wag
 - LG: usando -lg

Resultados

- Ignora las letras desconocidas e informa
- Muestra el número de cadenas únicas
- Muestra el tiempo consumido
- Muestra la plausibilidad (likelihood) de cada rama que calcula. Para ser comparada es necesario usar la opción -gamma para que la plausibilidad que devuelve sea Gamma20.
- El resultado es un árbol en formato Network
- La raíz del árbol es asignada arbitrariamente

Ejemplo

```
pendiente@DESKTOP-OQOS191:/mnt/c/Users/Guillermo/Documentos/Trabajo/Bioinformatica/P5$ ./FastTree -nt alineado.fasta>t.nwk
FastTree Version 2.1.10 SSE3
Alignment: alineado.fasta
Nucleotide distances: Jukes-Cantor Joins: balanced Support: SH-like 1000
Search: Normal +NNI +SPR (2 rounds range 10) +ML-NNI opt-each=1
TopHits: 1.00*sqrtN close=default refresh=0.80
ML Model: Jukes-Cantor, CAT approximation with 20 rate categories
Ignored unknown character k (seen 12 times)
Ignored unknown character m (seen 12 times)
Ignored unknown character r (seen 85 times)
Ignored unknown character s (seen 2 times)
Ignored unknown character w (seen 14 times)
Ignored unknown character y (seen 54 times)
Initial topology in 0.03 seconds
Refining topology: 21 rounds ME-NNIs, 2 rounds ME-SPRs, 11 rounds ML-NNIs
Total branch-length 0.488 after 0.31 sec1, 1 of 39 splits
ML-NNI round 1: LogLk = -3496.923 NNIs 9 max delta 4.31 Time 0.53
Switched to using 20 rate categories (CAT approximation)1 of 20
Rate categories were divided by 0.748 so that average rate = 1.0
CAT-based log-likelihoods may not be comparable across runs
Use -gamma for approximate but comparable Gamma(20) log-likelihoods
ML-NNI round 2: LogLk = -3283.944 NNIs 5 max delta 0.45 Time 0.75
ML-NNI round 3: LogLk = -3283.909 NNIs 0 max delta 0.00 Time 0.86
Turning off heuristics for final round of ML NNIs (converged)
ML-NNI round 4: LogLk = -3283.855 NNIs 1 max delta 0.05 Time 1.02 (final)
Optimize all lengths: LogLk = -3283.855 Time 1.07
Total time: 1.30 seconds Unique: 41/42 Bad splits: 0/38
pendiente@DESKTOP-OQOS191:/mnt/c/Users/Guillermo/Documentos/Trabajo/Bioinformatica/P5$ _
```


Salida

