Some of these slides have been borrowed from Dr. Paul Lewis, Dr. Joe Felsenstein. Thanks!

Paul has many great tools for teaching phylogenetics at his web site:

http://hydrodictyon.eeb.uconn.edu/people/plewis
We estimate branch lengths in terms of expected number of changes *per site*. To do this we standardize the total rate of divergence in the Q matrix and estimate $\nu = \mu t = 3\alpha t$ for each branch.

<table>
<thead>
<tr>
<th>From State</th>
<th>A</th>
<th>C</th>
<th>G</th>
<th>T</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>$-1$</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
</tr>
<tr>
<td>C</td>
<td>$\frac{1}{3}$</td>
<td>$-1$</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
</tr>
<tr>
<td>G</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
<td>$-1$</td>
<td>$\frac{1}{3}$</td>
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<tr>
<td>T</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
<td>$\frac{1}{3}$</td>
<td>$-1$</td>
</tr>
</tbody>
</table>
**JC “state-comparison” probabilities**

$i$ and $j$ refer to states (A, C, G, T). $i \neq j$:

\[
\Pr(i \rightarrow i | \nu) = \frac{1}{4} + \frac{3}{4} e^{-4 \nu / 3}
\]

\[
\Pr(i \rightarrow j | \nu) = \frac{1}{4} - \frac{1}{4} e^{-4 \nu / 3}
\]
CFN transition probabilities

\[
\begin{align*}
\Pr(0 \to 0|\nu) &= \Pr(1 \to 1|\nu) = \frac{1}{2} + \frac{1}{2}e^{-2\nu} \\
\Pr(0 \to 1|\nu) &= \Pr(1 \to 0|\nu) = \frac{1}{2} - \frac{1}{2}e^{-2\nu}
\end{align*}
\]
“$M_k$” transition probabilities

$k$-state version of the one-rate model.

$$\Pr(i \rightarrow i | \nu) = \frac{1}{k} + \frac{(k - 1)e^{-\left(\frac{k}{k-1}\right)\nu}}{k}$$

$$\Pr(i \rightarrow j | \nu) = \frac{1}{k} - \frac{e^{-\left(\frac{k}{k-1}\right)\nu}}{k}$$
**Kimura (1980) model or “the K80 model”**

Transitions and transversions occur at different rates:

<table>
<thead>
<tr>
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</thead>
<tbody>
<tr>
<td>A</td>
<td>$-2\beta - \alpha$</td>
<td>$\beta$</td>
<td>$\alpha$</td>
<td>$\beta$</td>
</tr>
<tr>
<td>C</td>
<td>$\beta$</td>
<td>$-2\beta - \alpha$</td>
<td>$\beta$</td>
<td>$\alpha$</td>
</tr>
<tr>
<td>G</td>
<td>$\alpha$</td>
<td>$\beta$</td>
<td>$-2\beta - \alpha$</td>
<td>$\beta$</td>
</tr>
<tr>
<td>T</td>
<td>$\beta$</td>
<td>$\alpha$</td>
<td>$\beta$</td>
<td>$-2\beta - \alpha$</td>
</tr>
</tbody>
</table>
Kimura (1980) model or “the K80 model”. Reparameterized.

We only care about the relative rates, so we can choose one rate to be frame of reference. This turns the 2 parameter model into a 1 parameter form:

<table>
<thead>
<tr>
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<th>A</th>
<th>C</th>
<th>G</th>
<th>T</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>$-(2+\kappa)\beta$</td>
<td>$\beta$</td>
<td>$\kappa\beta$</td>
<td>$\beta$</td>
</tr>
<tr>
<td>C</td>
<td>$\beta$</td>
<td>$-(2+\kappa)\beta$</td>
<td>$\beta$</td>
<td>$\kappa\beta$</td>
</tr>
<tr>
<td>G</td>
<td>$\kappa\beta$</td>
<td>$\beta$</td>
<td>$-(2+\kappa)\beta$</td>
<td>$\beta$</td>
</tr>
<tr>
<td>T</td>
<td>$\beta$</td>
<td>$\kappa\beta$</td>
<td>$\beta$</td>
<td>$-(2+\kappa)\beta$</td>
</tr>
</tbody>
</table>
Kimura (1980) model or “the K80 model”. Reparameterized again.

<table>
<thead>
<tr>
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<th>C</th>
<th>G</th>
<th>T</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>$-2 - \kappa$</td>
<td>1</td>
<td>$\kappa$</td>
<td>1</td>
</tr>
<tr>
<td>C</td>
<td>1</td>
<td>$-2 - \kappa$</td>
<td>1</td>
<td>$\kappa$</td>
</tr>
<tr>
<td>G</td>
<td>$\kappa$</td>
<td>1</td>
<td>$-2 - \kappa$</td>
<td>1</td>
</tr>
<tr>
<td>T</td>
<td>1</td>
<td>$\kappa$</td>
<td>1</td>
<td>$-2 - \kappa$</td>
</tr>
</tbody>
</table>
Kappa is the transition/transversion rate ratio:

\[ \kappa = \frac{\alpha}{\beta} \]

(if \( \kappa = 1 \) then we are back to JC).
What is the instantaneous probability of an particular transversion?

\[ \text{Pr}(A \rightarrow C) = \text{Pr}(A) \text{Pr(\text{change to } C)} = \frac{1}{4} (\beta dt) \]
What is the instantaneous probability of an particular transition?

\[ \Pr(A \rightarrow G) = \Pr(A) \Pr(\text{change to } G) = \frac{1}{4} (\kappa \beta dt) \]
There are four types of transitions:

\[ A \rightarrow G, G \rightarrow A, C \rightarrow T, T \rightarrow C \]

and eight types of transversions:

\[ A \rightarrow C, A \rightarrow T, G \rightarrow C, G \rightarrow T, C \rightarrow A, C \rightarrow G, T \rightarrow A, T \rightarrow G \]

\[
\text{Ti/Tv ratio} = \frac{\Pr(\text{any transition})}{\Pr(\text{any transversion})} = \frac{4 \left( \frac{1}{4} (\kappa \beta dt) \right)}{8 \left( \frac{1}{4} (\beta dt) \right)} = \frac{\kappa}{2}
\]

For K2P instantaneous transition/transversion ratio is one-half the instantaneous transition/transversion rate ratio
Kimura model change probabilities

\[
\text{Pr}(A \rightarrow A|\nu) = \frac{1}{4} \left( 1 + e^{-\left(\frac{4}{2+\kappa}\right)\nu} + 2e^{-\left(\frac{2+2\kappa}{2+\kappa}\right)\nu} \right)
\]

\[
\text{Pr}(A \rightarrow G|\nu) = \frac{1}{4} \left( 1 + e^{-\left(\frac{4}{2+\kappa}\right)\nu} - 2e^{-\left(\frac{2+2\kappa}{2+\kappa}\right)\nu} \right)
\]

\[
\text{Pr}(A \rightarrow C|\nu) = \frac{1}{4} \left( 1 - e^{-\left(\frac{4}{2+\kappa}\right)\nu} \right)
\]
### Felsenstein 1981 model or “F81 model”

<table>
<thead>
<tr>
<th>From State</th>
<th>A</th>
<th>C</th>
<th>G</th>
<th>T</th>
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</thead>
<tbody>
<tr>
<td>A</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\pi_T$</td>
</tr>
<tr>
<td>C</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\pi_T$</td>
</tr>
<tr>
<td>G</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\pi_T$</td>
</tr>
<tr>
<td>T</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\pi_T$</td>
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</tbody>
</table>
HKY 1985 model

<table>
<thead>
<tr>
<th>From State</th>
<th>A</th>
<th>C</th>
<th>G</th>
<th>T</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\kappa\pi_G$</td>
<td>$\pi_T$</td>
</tr>
<tr>
<td>C</td>
<td>$\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\kappa\pi_T$</td>
</tr>
<tr>
<td>G</td>
<td>$\kappa\pi_A$</td>
<td>$\pi_C$</td>
<td>$\pi_G$</td>
<td>$\pi_T$</td>
</tr>
<tr>
<td>T</td>
<td>$\pi_A$</td>
<td>$\kappa\pi_C$</td>
<td>$\pi_G$</td>
<td>$-$</td>
</tr>
</tbody>
</table>
F84* vs. HKY85

F84 model:

\[ \begin{align*}
\mu & \quad \text{rate of process generating all types of substitutions} \\
\kappa \mu & \quad \text{rate of process generating only transitions}
\end{align*} \]

Becomes F81 model if \( \kappa = 0 \)

HKY85 model:

\[ \begin{align*}
\beta & \quad \text{rate of process generating only transversions} \\
\kappa \beta & \quad \text{rate of process generating only transitions}
\end{align*} \]

Becomes F81 model if \( \kappa = 1 \)

### General Time Reversible – GTR model

<table>
<thead>
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<th>A</th>
<th>C</th>
<th>G</th>
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</thead>
<tbody>
<tr>
<td>A</td>
<td>−</td>
<td>$a \pi_C$</td>
<td>$b \pi_G$</td>
<td>$c \pi_T$</td>
</tr>
<tr>
<td>C</td>
<td>$a \pi_A$</td>
<td>−</td>
<td>$d \pi_G$</td>
<td>$e \pi_T$</td>
</tr>
<tr>
<td>G</td>
<td>$b \pi_A$</td>
<td>$d \pi_C$</td>
<td>−</td>
<td>$f \pi_T$</td>
</tr>
<tr>
<td>T</td>
<td>$c \pi_A$</td>
<td>$e \pi_C$</td>
<td>$f \pi_G$</td>
<td>−</td>
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</tbody>
</table>

In PAUP, $f = 1$ indicating that $G \rightarrow T$ is the reference rate.
Likelihood of a single sequence

First 32 nucleotides of the \( \psi \eta \)-globin gene of gorilla:

\[
\text{GAAGTCTTGGAGAAATAAACTGCACACACTGG}
\]

\[
L = \pi_G \pi_A \pi_A \pi_G \pi_T \pi_T \pi_G \pi_C \pi_C \pi_T \pi_G \pi_A \pi_A \pi_T \pi_A \pi_A \pi_C \pi_C \pi_T \pi_G \pi_C \pi_C \pi_A \pi_C \pi_A \pi_C \pi_T \pi_G \pi_G
\]

\[
= \pi_A^{12} \pi_C^7 \pi_G^7 \pi_T^6
\]

\[
\ln L = 12 \ln \left( \pi_A \right) + 7 \ln \left( \pi_C \right) + 7 \ln \left( \pi_G \right) + 6 \ln \left( \pi_T \right)
\]

We can already see by eye-balling this that the F81 model (which allows unequal base frequencies) will fit better than the JC69 model (which assumes equal base frequencies) because there are about twice as many As as there are Cs, Gs and Ts.
How can we calculate the likelihood score

Under the JC (or K2P) model:

\[
\ln L = 12 \ln \pi_A + 7 \ln \pi_C + 7 \ln \pi_G + 6 \ln \pi_T
\]
\[
= 12 \ln 0.25 + 7 \ln 0.25 + 7 \ln 0.25 + 6 \ln 0.25
\]
\[
= -44.361
\]
How can we calculate the likelihood score

Under the F81 (or HKY or GTR) model:

$$\ln L = 12 \ln \pi_A + 7 \ln \pi_C + 7 \ln \pi_G + 6 \ln \pi_T$$

But what are the values for the parameters: $\pi_A, \pi_C, \pi_G, \pi_T$?

In many cases we refer to these parameters as “nuisance parameters.” They must be specified in order to calculate the likelihood, but we are not interested in them by themselves.
We can find the maximum likelihood estimates of the parameters to give us the ML score: the maximum likelihood obtainable under this model:

\[
\ln L = 12 \ln \pi_A + 7 \ln \pi_C + 7 \ln \pi_G + 6 \ln \pi_T
\]
\[
= 12 \ln \hat{\pi}_A + 7 \ln \hat{\pi}_C + 7 \ln \hat{\pi}_G + 6 \ln \hat{\pi}_T
\]
\[
= 12 \ln 0.375 + 7 \ln 0.21875 + 7 \ln 0.21875 + 6 \ln 0.1875
\]
\[
= -43.091
\]

But how did I get the numbers to fill for the parameters? How do we know that \( \hat{\pi}_A = 0.375 \) and \( \hat{\pi}_C = 0.21875 \)
We might guess that:

\[ \hat{\pi}_A = \frac{12}{32} = 0.375 \]
\[ \hat{\pi}_C = \frac{7}{32} = 0.21875 \]
\[ \hat{\pi}_G = \frac{7}{32} = 0.21875 \]
\[ \hat{\pi}_T = \frac{6}{32} = 0.1875 \]

but how do we prove it?
ML parameter estimates

For simple problems we solve for the point in parameter space for which derivatives with respect to all parameters are 0 (we also have to consider boundary points).

We would have to do constrained optimization because

$$\pi_A + \pi_C + \pi_G + \pi_T = 1$$

and that is a pain.
ML parameter estimates

We can reparameterize:

\[
\begin{align*}
  r &= \pi_A + \pi_G \\
a &= \frac{\pi_A}{\pi_A + \pi_G} \\
c &= \frac{\pi_C}{\pi_C + \pi_T}
\end{align*}
\]

and always recover the original parameters:

\[
\begin{align*}
  \pi_A &= ra \\
  \pi_G &= r(1 - a) \\
  \pi_C &= (1 - r)c \\
  \pi_T &= (1 - r)(1 - c)
\end{align*}
\]
ML parameter estimates

\[
\ln L = 12 \ln \pi_A + 7 \ln \pi_C + 7 \ln \pi_G + 6 \ln \pi_T \\
= 12 \ln [ra] + 7 \ln [(1 - r)c] + 7 \ln [r(1 - a)] + 6 \ln [(1 - r)(1 - c)]
\]

Recall that:

\[
\frac{\partial \ln f(x)}{\partial x} = \frac{\partial f(x)}{\partial x} \cdot \frac{1}{f(x)}
\]
\[
\ln L = 12 \ln [ra] + 7 \ln [(1 - r)c] + 7 \ln [r(1 - a)] + 6 \ln [(1 - r)(1 - c)]
\]
\[
\frac{\partial \ln L}{\partial a} = \frac{12r}{ra} + \frac{7(-r)}{r(1 - a)}
\]
\[
= \frac{12}{a} - \frac{7}{(1 - a)}
\]
\[
0 = \frac{12}{\hat{a}} - \frac{7}{(1 - \hat{a})}
\]
\[
\hat{a} = \frac{12}{19}
\]
\[
\begin{align*}
\ln L &= 12 \ln [ra] + 7 \ln [(1 - r)c] + 7 \ln [r(1 - a)] + 6 \ln [(1 - r)(1 - c)] \\
\frac{\partial \ln L}{\partial c} &= \frac{7(1 - r)}{(1 - r)c} + \frac{6 - (1 - r)}{(1 - r)(1 - a)} \\
&= \frac{7}{c} - \frac{6}{(1 - c)} \\
0 &= \frac{7}{\hat{c}} - \frac{6}{(1 - \hat{c})} \\
\hat{c} &= \frac{7}{13}
\end{align*}
\]
\[
\ln L = 12 \ln [ra] + 7 \ln [(1 - r)c] + 7 \ln [r(1 - a)] + 6 \ln [(1 - r)(1 - c)]
\]

\[
\frac{\partial \ln L}{\partial r} = \frac{12a}{ra} + \frac{7(-c)}{(1 - r)c} + \frac{7(1 - a)}{r(1 - a)} + \frac{6(-1 - c)}{(1 - r)(1 - c)}
\]

\[
= \frac{12}{r} - \frac{7}{(1 - r)} + \frac{7}{r} - \frac{6}{1 - r}
\]

\[
= \frac{19}{r} - \frac{13}{(1 - r)}
\]

\[
0 = \frac{19}{\hat{r}} - \frac{13}{(1 - \hat{r})}
\]

\[
\hat{r} = \frac{19}{32}
\]
ML inference displays “scale invariance” so we can just transform the ML estimates into our original parameters:

\[
\hat{\pi}_A = \hat{r} \hat{\alpha} = \left(\frac{19}{32}\right) \left(\frac{12}{19}\right) = \frac{12}{32}
\]

\[
\hat{\pi}_G = \hat{r} (1 - \hat{\alpha}) = \left(\frac{19}{32}\right) \left(\frac{7}{19}\right) = \frac{7}{32}
\]

\[
\hat{\pi}_C = (1 - \hat{r}) \hat{\kappa} = \left(\frac{13}{32}\right) \left(\frac{6}{13}\right) = \frac{7}{32}
\]

\[
\hat{\pi}_T = (1 - \hat{r})(1 - \hat{\kappa}) = \left(\frac{13}{32}\right) \left(\frac{6}{13}\right) = \frac{6}{32}
\]
Likelihood ratio testing

\[ \ln L_{JC} = -44.361 \]
\[ \ln L_{F81} = -43.091 \]

But the F81 model has 3 more free parameters than JC.

The likelihood ratio test is a hypothesis testing approach to model selection.
Likelihood ratio testing

$H_0$: the data were generated under the simpler model
$H_A$: the data were generated under the more complex model.

test statistic: $2 \left( \ln L_{\text{complex}} - \ln L_{\text{simple}} \right)$

The LRT only works if the simple model is nested inside the more complex model (if the free parameters for the simple model are a subset of the free parameters for the more complex model).
Likelihood ratio testing

Null distribution: \( \chi^2 \) distribution with 
\( d.f. = \) difference in the number of free parameters.

If the LR test statistic is > than the critical value from the appropriate chi-square table, then we reject the simple model and prefer the more complex model.
Likelihood ratio testing - example

\[ \ln L_{JC} = -44.361 \]
\[ \ln L_{F81} = -43.091 \]
\[ LRT = 2(−43.091 − (−44.361)) = 2.54 \]
\[ df = 3 − 0 = 3 \]
\[ \chi^2_3(\text{critical, } P = 0.05) = 7.815 \]

Not significant. Do not reject the JC model.

(if we look up the \( P \)-value for this test statistic it is 0.4681).
Likelihoods on the simplest possible tree

\[\text{GA} \rightarrow \text{GG}\]

\[L = L_1 L_2\]

\[= \Pr(G) \Pr(G \rightarrow G) \Pr(A) \Pr(A \rightarrow G)\]

\[= \Pr(G) \Pr(G \rightarrow G|\nu) \Pr(A) \Pr(A \rightarrow G|\nu)\]

\[= \left(\frac{1}{4}\right) \left(\frac{1}{4} + \frac{3}{4} e^{-\frac{4\nu}{3}}\right) \left(\frac{1}{4}\right) \left(\frac{1}{4} - \frac{1}{4} e^{-\frac{4\nu}{3}}\right)\]
\[ d = \frac{1}{4} - \frac{1}{4} e^{-\frac{4\nu}{3}} \]

\[
\left( \frac{1}{4} + \frac{3}{4} e^{-\frac{4\nu}{3}} \right) = 1 - 3d
\]

\[
L = \left( \frac{1}{4} \right) \left( \frac{1}{4} + \frac{3}{4} e^{-\frac{4\nu}{3}} \right) \left( \frac{1}{4} \right) \left( \frac{1}{4} - \frac{1}{4} e^{-\frac{4\nu}{3}} \right)
\]

\[
= \frac{(1 - 3d)d}{16}
\]

\[
\frac{\partial \ln L}{\partial d} = \frac{1 - 6d}{16}
\]

\[
0 = \frac{1 - 6\hat{d}}{16}
\]

\[
\hat{d} = \frac{1}{6}
\]

\[
\hat{\nu} = 0.82396
\]

\[
L = 0.005208
\]
You may recall that the JC distance correction from lecture 8 looked like this:

$$\nu = \frac{-3}{4} \ln \left( 1 - \frac{4p}{3} \right)$$

If you put in $p = 0.5$, because half the sites differ in our example then you the same branch length:

$$\nu = 0.82396$$

Our JC distance correction formula is actually an ML estimator of the branch length between a pair of taxa.
The first 30 nucleotides of the ψη-globin gene

gorilla  \text{GAAGTCCTTGAGAAATAAAACTGCACACTGG}

orangutan  \text{GGACTCCTTGAGAAATAAAACTGCACACTGG}

\[ L = \left[ \left( \frac{1}{4} \right) \left( \frac{1}{4} + \frac{3}{4} e^{-\frac{4\nu}{3}} \right) \right]^{28} \left[ \left( \frac{1}{4} \right) \left( \frac{1}{4} - \frac{1}{4} e^{-\frac{4\nu}{3}} \right) \right]^2 \]

\[ \hat{\nu} = 0.06982 \]
\[ \ln L = -51.13396 \]
Likelihood of a tree
(data for only one site shown)

Arbitrarily chosen to serve as the root node

Ancestral states like this are not really known - we will address this in a minute.

Arbitrarily chosen to serve as the root node
Likelihood for site $k$

$$L_k = \frac{1}{4} \left[ \frac{1}{4} + \frac{3}{4} e^{-4v_1/3} \right] \left[ \frac{1}{4} + \frac{3}{4} e^{-4v_2/3} \right] \left[ \frac{1}{4} - \frac{1}{4} e^{-4v_3/3} \right] \left[ \frac{1}{4} - \frac{1}{4} e^{-4v_4/3} \right] \left[ \frac{1}{4} + \frac{3}{4} e^{-4v_5/3} \right]$$

$\pi_A$

$v_5$ is the expected no. substitutions for just this segment of the tree

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Brute force approach would be to calculate $L_k$ for all 16 combinations of ancestral states and sum
Pruning algorithm*
(same result, much less time)

Many calculations can be done just once, and then reused many times

*The pruning algorithm was introduced by: Felsenstein, J. 1981. Evolutionary trees from DNA sequences: a maximum likelihood approach. *Journal of Molecular Evolution* 17:368-376
<table>
<thead>
<tr>
<th>Taxon</th>
<th>Character</th>
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</thead>
<tbody>
<tr>
<td>1</td>
<td>A</td>
</tr>
<tr>
<td>2</td>
<td>C</td>
</tr>
<tr>
<td>3</td>
<td>C</td>
</tr>
<tr>
<td>4</td>
<td>C</td>
</tr>
<tr>
<td>5</td>
<td>G</td>
</tr>
</tbody>
</table>

![Diagram](attachment:image.png)
\[ L = \sum \sum \sum \sum \sum \Pr(x, y, z, w, A, C, C, C, G|\nu) \]
\[ L = \sum_x \sum_y \sum_z \sum_w \Pr(x) \Pr(y|x, \nu_6) \Pr(A|y, \nu_1) \Pr(C|y, \nu_2) \cdots \]
\[ \Pr(z|x, \nu_8) \Pr(C|z, \nu_3) \Pr(w|z, \nu_7) \Pr(C|w, \nu_4) \Pr(G|w, \nu_5) \]
\[ L = \sum_{x} \sum_{y} \sum_{z} \Pr(x) \Pr(y|x, \nu_6) \Pr(A|y, \nu_1) \Pr(C|y, \nu_2) \cdots \]

\[ \Pr(z|x, \nu_8) \Pr(C|z, \nu_3) \left( \sum_w \Pr(w|z, \nu_7) \Pr(C|w, \nu_4) \Pr(G|w, \nu_5) \right) \]
\[ L = \sum \sum \Pr(x) \Pr(y|x, \nu_6) \Pr(A|y, \nu_1) \Pr(C|y, \nu_2) \cdots \]

\[
\left( \sum \Pr(z|x, \nu_8) \Pr(C|z, \nu_3) \left( \sum \Pr(w|z, \nu_7) \Pr(C|w, \nu_4) \Pr(G|w, \nu_5) \right) \right)
\]
\[ L = \sum_x \Pr(x) \left( \sum_y \Pr(y|x, \nu_6) \Pr(A|y, \nu_1) \Pr(C|y, \nu_2) \right) \cdots \]

\[
\left( \sum_z \Pr(z|x, \nu_8) \Pr(C|z, \nu_3) \left( \sum_w \Pr(w|z, \nu_7) \Pr(C|w, \nu_4) \Pr(G|w, \nu_5) \right) \right)
\]
Maximum likelihood is a lot of work

- Site likelihoods involve products of transition probabilities, summed over ancestral states
- Overall log-likelihood for a tree is sum of site log-likelihoods
- Overall log-likelihood must be maximized!
  - must find MLEs for all edge lengths and all model parameters
  - this involves computing the overall log-likelihood many, many times (try turning on logiter in PAUP to get a feel for how much work this involves)
- Maximized lnL can now be compared to maximized lnL from other trees
Is it worth it?

• Uses all information
  – Parsimony ignores constant and autapomorphic sites
  – Distance methods ignore information not captured in pairwise comparisons

• Model generality
  – Some models possible with distance methods, but some quantities cannot be estimated reliably (e.g. variation in rates across sites)
  – Many parsimony variants exist, but parsimony does not allow estimation of the step matrix entries, for example
  – Many complex models are only possible under likelihood or Bayesian methods (which have a likelihood foundation)