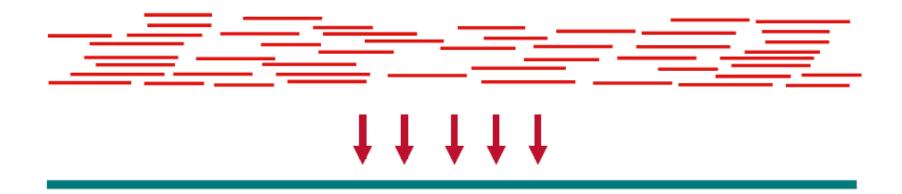
Read Mapping and Assembly

Gail Rosen

Read Mapping Leads to Assembly



- Modern fast read aligners: BWT, Bowtie, SOAP
 - Based on Burrows-Wheeler transform

Outline

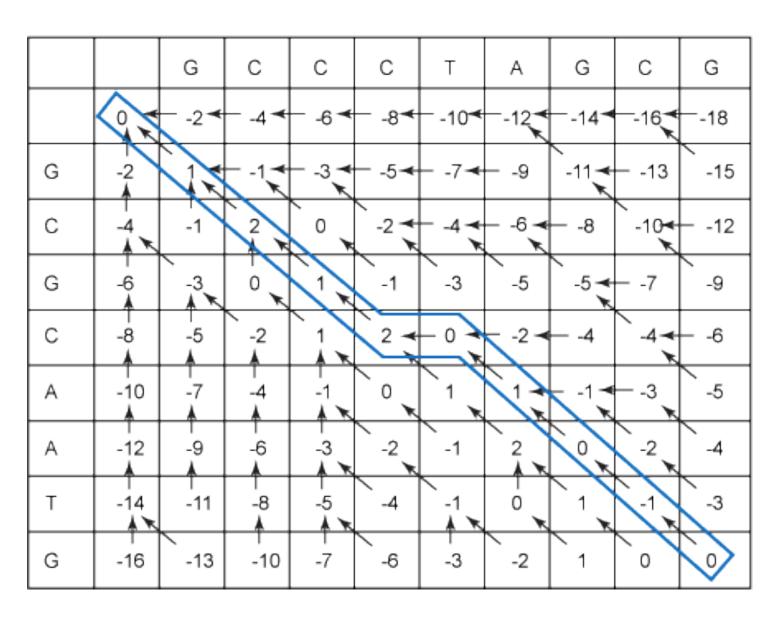
- Traditional Alignment approaches
- Read Mapping
 - Hashing
 - BWT
- Assembly
 - Difficult for metagenomics
- Binning Contigs

Alignment

- Needleman-Wunsch
- Smith-Waterman

Uses Dynamic Programming Approaches

Needleman Wunsch



Smith Waterman Alignment

		G	С	С	С	Т	А	G	С	G
	0 *	0	0	0	0	0	0 x	0	0 *	0
G	0	1 1	0 💌	0 *	0	0	0	1 x	0	1
С	0 1	0	2	1	1	0	0 *	0	2	0
G	0	1 *	0 🔭	1	0	0	0	1	0	3
С	0	0	2	1	2	0	0	0	2	1
А	0	0	0	1	0	1 1	1	0	0	1
А	0	0	0	0	0	0	2	0	0	0
Т	0 *	0	0	0	0	1	0 🔭	1	0 🗶	0
G	0	1	0	0	0	0	0	1	0	1

Approaches to Short Read Alignment

Two main approaches

- Hash-Based mapping:
 - Hashing of reads (E.g. Maq, Eland, SHRiMP)
 - Hashing of genome (E.g. novoalign, SHRiMP2)
- Indexing of tree-like structures:
 - Bowtie (ungapped)
 - Bowtie2 (gapped)
 - Bwa (gapped)
 - these all use Suffix Arrays/Burrows-Wheeler Transform (BWT), coupled with FM index

Hashing

 A hash function simply converts a string ("key") to an integer ("value").

• The integer is then used as an index in an array, for fast

look up.

Search can Still be O(nm) Time if hashing not good hash function keys buckets 00 521-8976 01 John Smith 02 521-1234 03 Lisa Smith 13 Sandra Dee 14 521-9655 15

MAQ

- First widely used open source short read aligner
- Very fast, but at the cost of accuracy (ungapped)
- Uses hashing to index sequence reads, then scans with reference sequence
- Guaranteed to find alignments with up to 2 mismatches
- Can take advantage of paired end reads
- Uses quality scores to determine best alignments
- Generally no longer used, as has been superceded by newer aligners

http://sourceforge.net/projects/maq/

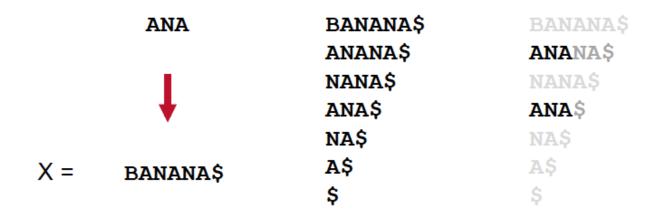
Bowtie

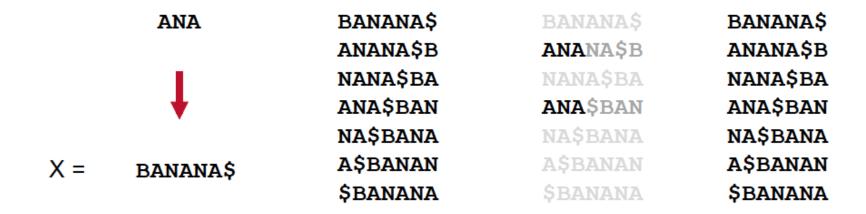
- Similar to MAQ, in that it uses quality scores to find best alignments.
- Uses "Burrows-Wheeler index" to keep its memory footprint small.
- Can find alignments with up to 3 mismatches in the first L bases of the read.
- Only ungapped alignments
- Also supports paired end reads. http://bowtie-bio.sourceforge.net/
- Bowtie2 supports gapped alignments too.

Burrows Wheeler Transform (BWT)

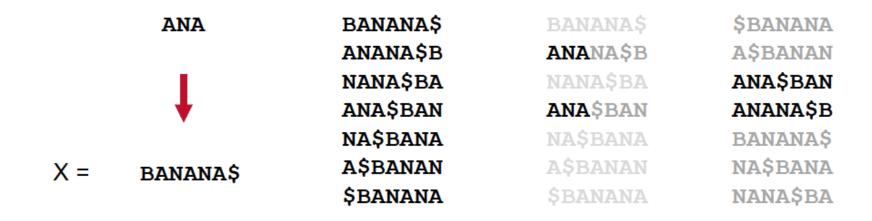
The BWT of a string S is a reversible permutation of S that enables the search for a pattern P in S to take linear time with respect to the length of P (O(|P|) time, independent of the length of S)

> suffixes of BANANA

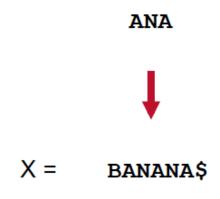


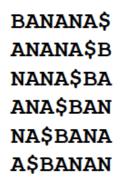


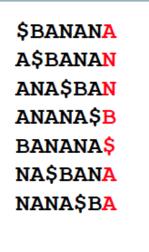
Suffixes sharing a common prefix are grouped together with a BWT



Tends to put runs of same character together (good for compression)







BWT matrix of string 'BANANA'

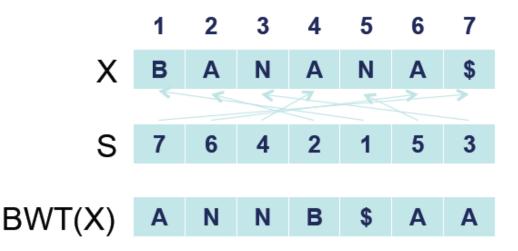
BWT(BANANA) = ANNB\$AA

Suffix Array

\$BANANA 1 \$BANANA A\$BANAN 2 A\$BANAN ANA\$BAN 3 ANA\$BAN ANANA\$B 4 ANANA\$B BANANA\$ 5 BANANA\$ NA\$BANA 6 NA\$BANA NANA\$BA 7 NANA\$BA

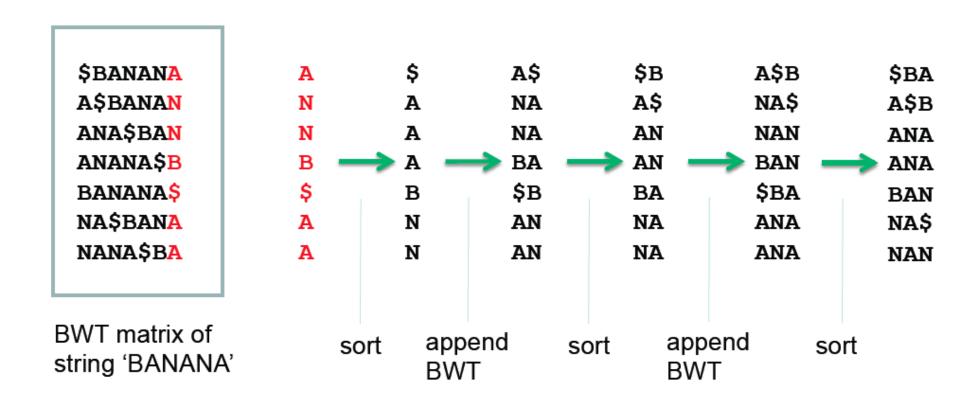
Suffixes are sorted in the BWT matrix

S(i) = j, where $X_j ... X_n$ is the i-th suffix lexicographically



BWT(X) constructed from S: At each position, take the letter to the left of the one pointed by S

Reconstructing BANANA



Reconstructing BANANA - faster

\$BANANA

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

Lemma. The i-th occurrence of character c in last column is the same text character as the i-th occurrence of c in the first column

\$BANANA

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

Reconstructing BANANA - faster

\$BANANA

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

Lemma. The i-th occurrence of character c in last column is the same text character as the i-th occurrence of c in the first column

A \$BANAN

NA\$BANA

NANA\$BA

BANANA\$

\$BANANA

ANA\$BAN

ANANA\$B

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

Lemma. The i-th occurrence of character c in last column is the same text character as the i-th occurrence of c in the first column

A \$BANAN

NA\$BANA

NANA\$BA

BANANA\$

\$BANANA

ANA\$BAN

ANANA\$B

A\$BANAN ANA\$BAN ANANA\$B

Same words, same sorted order

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

Lemma. The i-th occurrence of character 'a' in last column is the same text character as the i-th occurrence of 'a' in the first column

LF(): Map the i-th occurrence of character 'a' in last column to the first column

LF(r): Let row r contain the i-th occurrence of 'a' in last column

Then, LF(r) = r'; r': i-th row starting with 'a'

A\$BANAN

ANA\$BAN

ANANA\$B

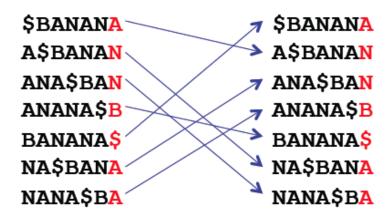
BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

LF(r): Let row r be the i-th occurrence of 'a' in last column Then, LF(r) = r'; r': i-th row starting with 'a'

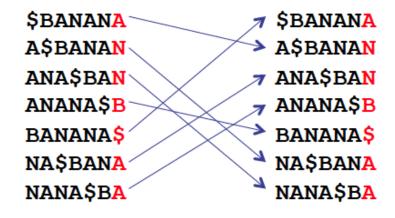


LF[] = [2, 6, 7, 5, 1, 3, 4]

Row LF(r) is obtained by rotating row r one position to the right

\$BANANA
A\$BANAN
ANA\$BAN
ANANA\$B
BANANA\$
NA\$BANA

BWT matrix of string 'BANANA'



$$LF[] = [2, 6, 7, 5, 1, 3, 4]$$

Computing LF() is easy:

Let C(a): # of characters smaller than 'a' Example: C(\$) = 0; C(A) = 1; C(B) = 4; C(N) = 5

Let row r end with the i-th occurrence of 'a' in last column

Then,
$$LF(r) = C(a) + i$$

\$BANANA
A\$BANAN
ANA\$BAN
ANANA\$B
BANANA\$
NA\$BANA

BWT matrix of string 'BANANA'

```
Reconstruct BANANA:

S := ""; r := 1; c := BWT[r];

UNTIL c = '$' {

S := cS;

r := LF(r);

c := BWT(r); }
```

Credit: Ben Langmead thesis

A\$BANAN

ANA\$BAN

ANANA\$B

BANANA\$

NA\$BANA

NANA\$BA

BWT matrix of string 'BANANA'

L(W): lowest index in BWT matrix where W is prefix U(W): highest index in BWT matrix where W is prefix

Example:

$$L("NA") = 6$$

 $U("NA") = 7$

Lemma (prove as exercise)

$$L(aW) = C(a) + i + 1,$$

$$where i = \# 'a's up to L(W) - 1 in BWT(X)$$

$$U(aW) = C(a) + j,$$

$$where j = \# 'a's up to U(W) in BWT(X)$$

Example:

Summary of BWT algorithm

Suffix array of string X:

S(i) = j, where $X_j ... X_n$ is the j-th suffix lexicographically

- BWT follows immediately from suffix array
 - Suffix array construction possible in O(n), many good O(n log n) algorithms
- Reconstruct X from BWT(X) in time O(n)
- Search for all exact occurrences of W in time O(|W|)
- BWT(X) is easier to compress than X

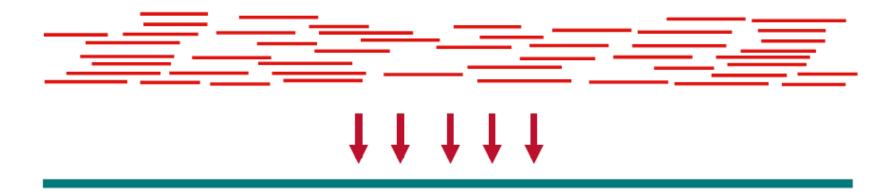
BWT-based Aligners In Practice

- Inexact matching: allow mismatches and gaps during alignment
 - => keep track of a set of SA intervals
- Heuristics: seeds, bounds on number of allowed differences, scoring (gap open/extend, mismatches)
- Memory considerations: sampling the suffix array and the occurrence array, compression
- Typical aligner phases
 - stage 1: BWT index construction
 - stage 2: short-Read Mapping
 - stage 3: alignment results reporting/evaluation

BWA

- From the author of Maq, but now also uses
 Burrows-Wheeler transform to significantly speed
 it up, and use less memory.
- Can also find small indels, in contrast to both Maq and Bowtie.
- Is slightly slower than bowtie, but ability to find indels make it more useful if SNVs are important to you.

Back to Assembly



CATCGACCGAGCGCGATGCTAGCTAGGTGATCGT.....

TGCCGCATCGACCGAGCGCGATGCTAGCTAGGTGATCGT...

GCATGCCGCATCGACCGAGCGCGATGCTAGCTAGGTGATCGT

GTGCATGCCGCCATCGACCGAGCGCGATGCTAGCTAGCTCATCCT

k-mer

"k-mer" is a substring of length k

```
S: GGCGATTCATCG mer: from Greek meaning "part"
```

A 4-mer of S: ATTC

```
All 3-mers of S: GGC
GCG
CGA
GAT
ATT
TTC
TCA
CAT
ATC
```

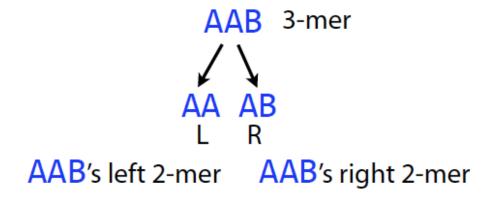
I'll use "k-1-mer" to refer to a substring of length k - 1

TCG

As usual, we start with a collection of reads, which are substrings of the reference genome.

AAA, AAB, ABB, BBB, BBA

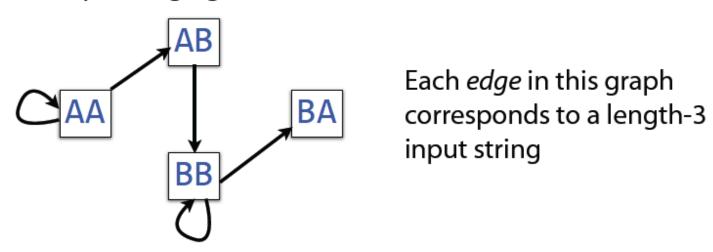
AAB is a k-mer (k = 3). AA is its left k-1-mer, and AB is its right k-1-mer.

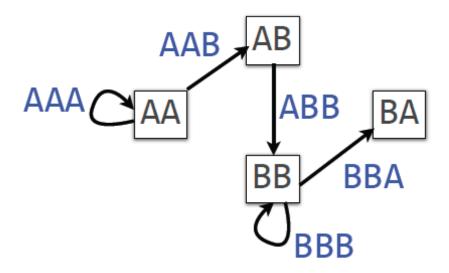


Take each length-3 input string and split it into two overlapping substrings of length 2. Call these the *left* and *right 2-mers*.

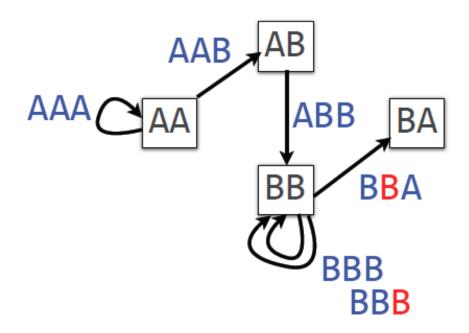
AAABBBA take all 3-mers: AAA, AAB, ABB, BBB, BBA

Let 2-mers be nodes in a new graph. Draw a directed edge from each left 2-mer to corresponding right 2-mer:





An edge corresponds to an overlap (of length k-2) between two k-1 mers. More precisely, it corresponds to a k-mer from the input.



If we add one more B to our input string: AAABBBBA, and rebuild the De Bruijn graph accordingly, we get a *multiedge*.

A procedure for making a De Bruijn graph for a genome

Assume *perfect sequencing* where each length-*k* substring is sequenced exactly once with no errors

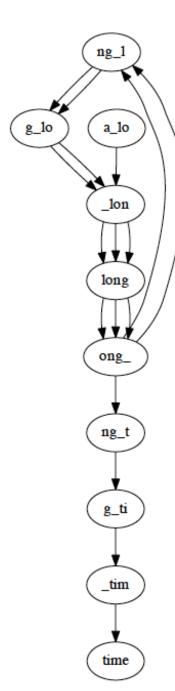
Pick a substring length k: 5

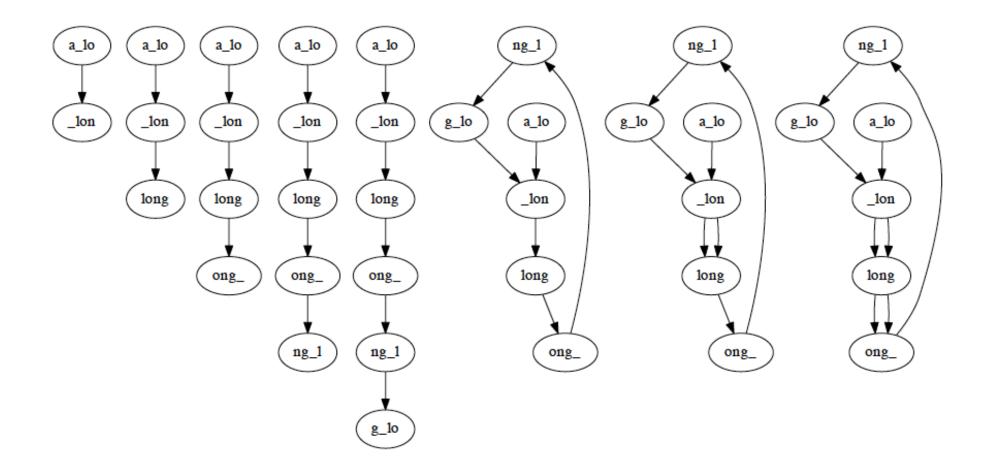
Start with an input string: a_long_long_time

Take each *k* mer and split into left and right *k*-1 mers

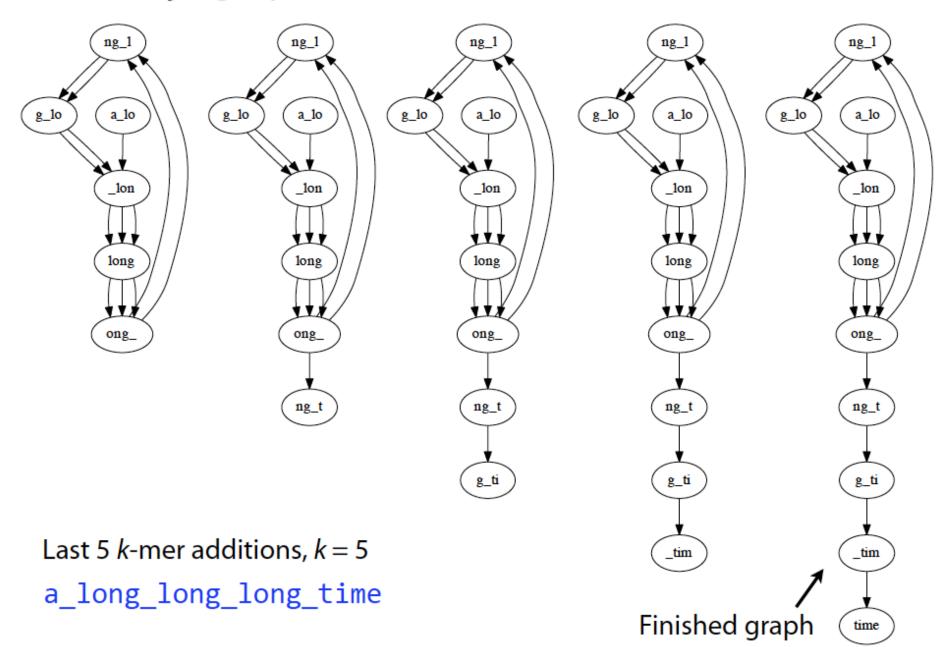
long_ long ong_

Add k-1 mers as nodes to De Bruijn graph (if not already there), add edge from left k-1 mer to right k-1 mer





First 8 k-mer additions, k = 5a_long_long_time



In practice, De Bruijn graph-based tools give up on unresolvable repeats and yield fragmented assemblies, just like OLC tools.

But first we note that using the De Bruijn graph representation has other advantages...

genome of length \mathbf{m} $\mathbf{m} = 3 \times 10^9 \text{ nt} \approx \text{human}$

Say a sequencer produces
$$d = 6 \times 10^9 \text{ reads}$$
 $\approx 1 \text{ sequencing run}$ d reads of length $m = 100 \text{ nt}$ $m = 3 \times 10^9 \text{ nt} \approx 1 \text{ sequencing run}$

To build a De Bruijn graph in practice:

Pick k. Assume $k \le$ shortest read length (k = 30 to 50 is common).

For each read:

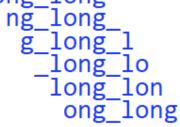
For each *k*-mer:

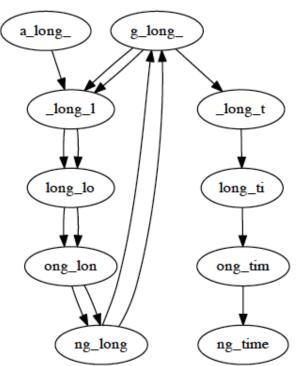
Add k-mer's left and right k-1-mers to graph if not there already. Draw an edge from left to right k-1-mer.

```
Pick k = 8 Genome: a_long_long_time
```

Reads: a_long_long_long, ng_long_l, g_long_time

k-mers: a_long_l ng_long_ g_long_t _long_lo g_long_l _long_ti _long_lon ong long ong time





Given n (# reads), N (total length of all reads) and k, and assuming k < length of shortest read:

Exact number of k-mers: N - n(k-1) O(N)

This is also the number of edges, |E|

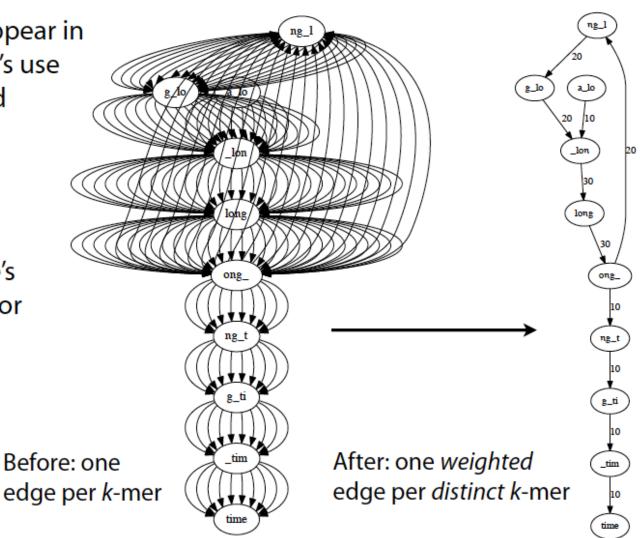
Number of nodes |V| is at most $2 \cdot |E|$, but typically much smaller due to repeated k-1-mers

In typical assembly projects, average coverage is ~ 30 - 50

Same edge might appear in dozens of copies; let's use edge weights instead

Weight = # times k-mer occurs

Using weights, there's one weighted edge for each distinct k-mer



What did we give up?

Reads are immediately split into shorter k-mers; can't resolve repeats as well as overlap graph

Only a very specific type of "overlap" is considered, which makes dealing with errors more complicated, as we'll see

Read coherence is lost. Some paths through De Bruijn graph are inconsistent with respect to input reads.

This is the OLC \leftrightarrow DBG tradeoff

Single most important benefit of De Bruijn graph is the $O(\min(G, N))$ space bound, though we'll see this comes with large caveats

GroopM

