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# **Background**

### 1.1 Introduction

#### 1.1.1 Motivation

This work has been motivated by recent advances of molecular genetics. The human genome has been sequenced in 2001. Also mouse, drosophila, etc. Nowadays # reference model genomes are available in genbank.

Next-generation sequencing has been the second revolution. NGS produces billions of reads for 1000\$ dollars. Why should one re-sequence a known genome? Resequencing applications include variant calling, etc. So NGS impacts biomedicine.

Given a set of reads, two approaches are possible: assembly and mapping.

Assembly methods are based on overlaps, de brujin graphs, or...

Read mapping methods work on a previously assembled reference genome.

The typical SNPs analysis pipeline 1.1 consists of...

In this work we focus on read mapping, although many core algorithms considered are also applicable to assembly, as well as to later pipeline stages.

Figure 1.1: NGS pipeline.

### 1.1.2 Organization of this manuscript

## 1.2 Stringology preliminaries

We now introduce fundamental definitions and problems of stringology, in order to keep the manuscript self-contained. The reader familiar with basic stringology can skip this section and proceed to section ??.

#### 1.2.1 Definitions

Let us start by defining primitive objects of stringology: alphabets and strings. An alphabet is a finite set of symbols (or characters); a string (or word) over an alphabet is a finite sequence of symbols from that alphabet. We denote the length of a string s by |s|, and by

 $\epsilon$  the empty string s.t.  $|\epsilon| = 0$ . Given an alphabet  $\Sigma$ , we define  $\Sigma^0 = \{\epsilon\}$  as the set containing the empty string,  $\Sigma^n$  as the set of all strings over  $\Sigma$  of length n, and  $\Sigma^* = \bigcup_{n=0}^{\infty} \Sigma^n$  as the set of all strings over  $\Sigma$ . Finally, we call any subset of  $\Sigma^*$  a language over  $\Sigma$ .

We now define concatenation, the most fundamental operation on strings. The concatenation operator of two strings is denoted with  $\cdot$  and defined as  $\cdot: \Sigma^* \times \Sigma^* \to \Sigma^*$ . Given two strings,  $x \in \Sigma^n$  with  $x = x_1 x_2 \dots x_n$ , and  $y \in \Sigma^m$  with  $y = y_1 y_2 \dots y_m$ , their concatenation  $x \cdot y$  (or simply denoted xy) is the string  $z \in \Sigma^{n+m}$  consisting of the symbols  $x_1 x_2 \dots x_n y_1 y_2 \dots y_m$ .

From concatenation we can derive the notion of prefix, suffix, and substring. A string x is a prefix of y iff there is some string z s.t.  $y = x \cdot z$ . Analogously, x is a suffix of y iff there is some string z s.t.  $y = z \cdot x$ . Moreover, x is a substring of y iff there is some string y s.t.  $y = y \cdot x \cdot z$ , and then we say that y occurs within y at position y.

**Example 1.1.** These definitions allow us to model basic biological sequences. Let us consider the alphabet consisting of DNA bases:  $\Sigma = \{A, C, G, T\}$ . Examples of strings over  $\Sigma$  are x = A, y = AGGTAC, z = TA. For instance,  $y \in \Sigma^6$  and |y| = 6. Moreover, the concatenation  $x \cdot z$  produces ATA. The string x is a prefix of y, and the string z is a substring of y occurring at position 4 in y.

### 1.2.2 Transformations

The next step is to define the minimal set of edit operations to transform one string into another: substitutions, insertions and deletions. Given two strings x, y of equal length n, the string x can be transformed into the string y by substituting (or replacing) all symbols  $x_i$  s.t.  $x_i \neq y_i$  into  $y_i$ , for  $1 \leq i \leq n$ . If the given strings have different lengths, insertion and deletion of symbols from x become necessary to transform it into y. Therefore, given any two strings x, y, we define as edit transcript for x, y any finite sequence of substitutions, insertions and deletions transforming x into y.

#### **Example 1.2.** TODO: example of edit transcript.

Edit transcripts lead us to the definition of distance functions between strings. The Hamming distance between two strings  $x,y\in \Sigma^n$  is defined as the function  $d_H:\Sigma^n\times \Sigma^n\to \mathbb{N}$  counting the number of substitutions necessary to transform x into y. More generally, the edit (or Levenshtein) distance between two strings  $x,y\in \Sigma^*$  is defined as the function  $d_E:\Sigma^*\times \Sigma^*\to \mathbb{N}$  counting the minimum number of edit operation necessary to transform x into y.

**Example 1.3.** TODO: example of edit and hamming distance.

### 1.2.3 Edit distance computation

The edit distance problem is to compute the edit distance between two given strings, along with an optimal edit transcript that describes the transformation? The edit distance problem is a minimization problem and can be efficiently computed via dynamic

programming (DP). Below we describe the three essential components of the DP approach: the recurrence relation, the DP table, and the traceback.

Given two strings x, y, for all  $1 \le i \le |x|$  and  $1 \le j \le |y|$  we define with  $d_E(x_{1..i}, y_{1..j})$  the edit distance between their prefixes  $x_{1..i}$  and  $y_{1..j}$ . The base conditions of the recurrence relation are:

$$d_E(\epsilon, \epsilon) = 0 \tag{1.1}$$

$$d_E(x_{1.i}, \epsilon) = i \text{ for } 1 \le i \le |x| \tag{1.2}$$

$$d_E(\epsilon, y_{1..j}) = j \text{ for } 1 \le j \le |y| \tag{1.3}$$

and the recursive case is defined as follows:

$$d_E(x_{1,i}, y_{1,i}) = \min\{d_E(x_{1,i-1}, y_{1,i}) + 1, d_E(x_{1,i}, y_{1,i-1}) + 1, d_E(x_{1,i-1}, x_{1,i-1}) + \delta(x_i, y_i)\}$$
(1.4)

The recurrence relation can be computed in time  $\mathcal{O}(|x| \cdot |y|)$  using a table of  $(n+1) \times (m+1)$  cells. However only  $\mathcal{O}(\min\{n,m\})$  space is required.

The table can be filled in four different ways: column-wise, row-wise, diagonal-wise or antidiagonal-wise.

**Figure 1.2:** DP table representing the computation of the edit distance  $d_E(x_{1..5}, y_{1..4})$ .

An optimal alignment can be computed in time O(n + m).

### 1.2.4 Alignments

An alignment is a way of visualizing a transformation between two strings.

The problem of finding the optimal alignment between two strings is the dual of the edit distance problem.

**Example 1.4.** TODO: example of alignment.

## 1.3 Overview of string matching

#### 1.3.1 Problem definition

We can now define *exact string matching*, perhaps the most fundamental problem in stringology. Given a string p (with |p| = m) called the *pattern* and a longer string t (with |t| = n) called the *text*, the exact string matching problem is to find all occurrences, if any, of pattern p into text t?. This problem has been extensively studied from the theoretical standpoint and is well solved in practice.

The definition of distance functions between strings let us generalize exact string matching into a more challenging problem: *approximate string matching*. Given a text

t, a pattern p, and a *distance threshold*  $k \in \mathbb{N}$ , the approximate string matching (a.s.m.) problem is to find all occurrences of p into t within distance k. The a.s.m. problem under the Hamming distance is commonly referred as the k-mismatches problem and under the edit distance as the k-differences problem. For k-mismatches and k-differences, it must hold k>0 as the case k=0 corresponds to exact string matching, and k< m as a pattern occurs at any position in the text if we substitute all its m characters. Under these distances, we define the *error rate* as  $\epsilon=\frac{k}{m}$ , with  $0<\epsilon<1$  given the above conditions, and we alternatively refer to these a.s.m. problems as  $\epsilon$ -mismatches and  $\epsilon$ -differences.

We can classify string matching problems in two categories, *online* and *offline*, depending on which string, the pattern or the text, is given first. Algorithms for online string matching work by preprocessing the pattern and scanning the text from left to right (or right to left); their worst-case runtime complexity ranges from  $\mathcal{O}(nm)$  to  $\mathcal{O}(n)$  while their worst-case memory complexity ranges from  $\mathcal{O}(\sigma^k m^k)$  to  $\mathcal{O}(m)$ . Algorithms for offline string matching are instead allowed to preprocess the text; their worst-case runtime complexity ranges from  $\mathcal{O}(m)$  to  $\mathcal{O}(\sigma^k m^k)$  while their worst-case memory complexity is usually  $\mathcal{O}(n)$ . In practice, if the text is long, static and searched frequently, offline methods largely outperform online methods in terms of runtime, provided the necessary amount of memory. Therefore, we concentrate on offline algorithms.

We can subdivide algorithms for offline string matching in two categories: *fully-indexed* and *filtering*. Fully-indexed algorithms work solely on the index of the text, while filtering methods first use the index to discard uninteresting portions of the text and subsequently use an online method to verify narrow areas of the text. Filtering methods outperform fully-indexed methods for a vast range of inputs<sup>1</sup> and are thus very interesting from a practical standpoint. Nonetheless, filtering methods are just opportunistic combinations of online and fully-indexed methods.

In the following of this section we thus give a brief and non-exhaustive overview of the fundamental techniques for online and (both fully-indexed and filtering) offline string matching. This overview serves as an introduction to the more involved algorithms presented in chapter ?? and directly used in applications of part ??. For an extensive treatment of the subject, we refer the reader to complete surveys on exact ? and approximate ? online string matching methods, as well as to a succint survey on indexed methods ?.

#### 1.3.2 Online methods

We consider two classes of algorithms for online string matching, those based on automata and those based on dynamic programming.

#### **Automata**

Exact search of one pattern. Boyer-moore automaton.

Exact search of multiple patterns. Aho-corasick automaton.

Approximate search of one pattern. Ukkonen automaton.

<sup>&</sup>lt;sup>1</sup> When the error rate is low.

#### **Dynamic programming**

The dynamic programming algorithm ?? to compute the edit distance of two strings can be easily turned into a string matching algorithm. Since an occurrence of the pattern can start and end anywhere in the text, a.s.m. consists of computing the edit distance between the pattern and all substrings of the text. The problem can be thus solved by computing the edit distance between the text and the pattern without penalizing leading and trailing deletions in the text.

Let pose x = t and y = p and consider equations ??. Since an occurrence of the pattern can start anywhere in the text, we change the initialization of the top row as:

$$d_E(\epsilon, y_{1..j}) = 0 \text{ for } 1 \le j \le |y| \tag{1.5}$$

and since an occurrence of the pattern can end anywhere in the text, we check every cell of the bottom row for the condition:

$$d_E(x_{1..m}, y_{1..j}) \le k \text{ for } 1 \le j \le |y|. \tag{1.6}$$

**Figure 1.3:** DP table representing the match of p = ... in t = ...

### 1.3.3 Indexed methods

No matter how fast online methods can be, these approaches quickly become impractical when the text is long and searched frequently. If the text is static and given in advance, we are allowed to preprocess it. We build an index of the text beforehand and use it to speed up subsequent searches. To this intent, we first introduce *suffix trees*, optimal data structures to index strings. Later on, we consider algorithms solving string matching problems on indices.

The suffix tree is a idealized data structure to elegantly expose our algorithms. In chapter ?? we consider other substring indices.

### Suffix tree and suffix trie

The suffix tree [?] is a lexicographically ordered tree data structure representing all suffixes of a string. Assume w.l.o.g. a string s of length n, padded with a *terminator symbol* s not being part of the string alphabet s. The suffix tree s of the string s has one node designated as the root and s leaves, each one pointing to a distinct suffix of s, denoted as s larger s larger

In the following of this manuscript we consider w.l.o.g. *suffix tries* instead of suffix trees. On suffix tries, internal nodes can have only one child and each edge is labeled by

 $<sup>^2</sup>$  The terminator symbol is necessary to ensure that no suffix  $s_{i..n}$  is a prefix of another suffix  $s_{j..n}$ .

one single character (see Figure ??). This fact simplifies the exposition of all given algorithms without affecting their runtime complexity nor their result. However, we remark that all given algorithms can be generalized to work on trees.

Figure 1.4: Suffix tree and suffix trie.

Therefore, from now on we assume the text t to be indexed using a suffix trie  $\mathcal{T}$ . Given a node x of  $\mathcal{T}$ , we denote with label(x) the label of the edge entering into x, with  $\mathbb{C}(x)$  the set of children of x being internal nodes, with  $\mathbb{E}(x)$  the set of children of x being leaves, and with  $\mathbb{L}(x)$  the set of all leaves of the subtree rooted in x. We remark that entering edges of internal nodes in  $\mathbb{C}(x)$  are always labeled with symbols in  $\Sigma$ , while entering edges of leaves in  $\mathbb{L}(x)$  and  $\mathbb{E}(x)$  are always labeled with terminator symbols.

#### **Exact string matching**

Using the suffix trie  $\mathcal T$  of the text t, we can find all occurrences of a pattern p into t in optimal time  $\mathcal O(m)$ , thus independently of n. Algorithm 1.1 searches a pattern p by starting in the root node of  $\mathcal T$  and following the path spelling the pattern. If the search ends up in a node x, each leaf  $l_i \in \mathbb L(x)$  points to a distinct suffix  $t_{i..n}$  such that  $t_{i..i+m} = p$ . Algorithm 1.1 is correct since each path from the root to any internal node of the suffix trie  $\mathcal T$  spells a different unique substring of t; consequently all equal substrings of t are represented by a single common path.

### **Algorithm 1.1** Exact string matching on a suffix trie.

```
1: procedure ExactSearch(x, p)
2: if p = \epsilon then
3: report \mathbb{L}(x)
4: else if \exists c_x \in \mathbb{C}(x) : label(c_x) = p_1 then
5: ExactSearch(c_x, p_{2..|p|})
6: end if
7: end procedure
```

Figure 1.5: Exact string matching on a suffix tree.

### Backtracking k-mismatches

We can solve k-mismatches by backtracking [??] on the suffix trie  $\mathcal{T}$ , in average time sublinear in n [?]. A top-down traversal on  $\mathcal{T}$  spells incrementally all distinct substrings of t. While traversing each branch of the trie, we incrementally compute the distance between the query and the spelled string. If the computed distance exceeds k, we stop the traversal and proceed on the next branch. Conversely, if we completely spelled the pattern p, and we ended up in a node x, each leaf  $l_i \in \mathbb{L}(x)$  points to a distinct suffix  $t_{i..n}$  such that  $d_H(t_{i..i+m}, p) \leq k$ . See algorithm 1.2.

### **Algorithm 1.2** *k*-mismatches on a suffix trie.

```
1: procedure KMISMATCHES(x, p, e)
        if e = 0 then
 2:
            EXACTSEARCH(x, p)
 3:
 4:
            for all c_x \in \mathbb{C}(x) do
 5:
                if label(c_x) = p_1 then
 6:
 7:
                    KMISMATCHES(c_x, p_{2..|p|}, e)
 8:
                else
 9:
                    KMISMATCHES(c_x, p_{2..|p|}, e-1)
10:
        end if
11:
12: end procedure
```

### Backtracking k-differences

We can compute k-differences on a suffix tree in two different ways. Algorithm 1.3 explicitly enumerates errors by recursing on the suffix trie. Algorithm 1.4 computes the edit distance on the suffix trie.

### **Algorithm 1.3** k-differences on a suffix trie.

```
1: procedure KDifferences(x, p, e)
 2:
        if e = 0 then
            EXACTSEARCH(x, p)
 3:
 4:
        else
            KDifferences(x, p_{2..|p|}, e - 1)
 5:
            for all c_x \in \mathbb{C}(x) do
 6:
 7:
                KDifferences(c_x, p, e-1)
 8:
                if label(c_x) = p_1 then
                    KDifferences(c_x, p_{2..|p|}, e)
 9:
10:
                else
                   KDifferences(c_x, p_{2..|p|}, e-1)
11:
                end if
12:
        end if
13:
14: end procedure
```

#### **Algorithm 1.4** *k*-difference on a suffix trie.

```
    procedure KDIFFERENCES(x, p, e)
    for all c<sub>x</sub> ∈ ℂ(x) do
    KDIFFERENCES(c<sub>x</sub>, p<sub>2..|p|</sub>, e − d<sub>E</sub>(repr(x), p))
    end procedure
```

In algorithm 1.4, any online method can be used to compute the edit distance. However, for theoretical considerations, it is important to consider an algorithm which computes in  $\mathcal{O}(1)$  per node. Note that it is sufficient to have an algorithm capable of checking whether the current edit distance is within the imposed threshold k.

Algorithm 1.3 reports more occurrences than algorithm 1.4. Discuss neighborhood, condensed neighborhood, and super-condensed neighborhood.

### 1.3.4 Filtering methods

Filtering methods work under the assumption that given patterns occurs in the text with a *low average probability*. Such probability is a function of the error rate  $\epsilon$ , in addition to the alphabet size  $\sigma$ , and can be computed or estimated under the assumption of the text being generated by a specific random source, e.g. the uniform Bernoulli model where each symbol of  $\Sigma$  occurs with probability  $\frac{1}{\sigma}$ . Under such assumptions, [?] estimates that  $\epsilon < 1 - \frac{1}{\sigma}$  is a conservative bound on the error rate which ensures few matches, and for which filtering algorithms are effective. For higher error rates, non-filtering online and indexed methods work better.

The goal of filtering methods is thus to obtain algorithms that have favorable average running times. The principle under which they work is that large and uninteresting portions of the text can be quickly discarded, while narrow and highly similar portions can be verified with a conventional online method. We remark that any filtering method can either work in a online fashion or take advantage of an index of the text to speed up the filtration phase. Here we always consider the filtration phase to be indexed.

We call a filter *lossless* or *full-sensitive* if it guarantees not to discard any occurrence of the pattern, otherwise we call it *lossy*. Lossy filters can be designed to solve approximately  $\epsilon$ -differences. We focus our attention on lossless filters.

We now consider two classes of filtering methods: those based on q-grams and those based on seeds. Filters based on q-grams consider all overlapping substrings of the pattern having length q, the so-called q-grams. Eventually, simple lemmas gives us lower bounds on the number of q-grams that must be present in a narrow window of the text as necessary condition for an occurrence of the pattern. Instead, seed based filters partition the pattern into non-overlapping factors called seeds. We can derive full-sensitive partitioning strategies by application of the pigeonhole principle.

### q-Gram filters

q-Gram filters are based on the so-called q-gram similarity measure  $\tau_q: \Sigma^* \times \Sigma^* \to \mathbb{N}$ , defined as the number of substrings of length q common to two given strings. The following lemma relates the q-gram similarity to edit distance<sup>3</sup>. It gives a lower bound on the q-gram similarity  $\tau_q(x,y)$  for any two strings having  $d_E(x,y)=k$ . This means that  $\tau_q(x,y) \geq k$  is a necessary but not sufficient condition for  $d_E(x,y) \leq k$ .

**Lemma 1.1.** ? Let x, y be two strings with edit distance k with  $min\{|x|, |y|\} = m$ , then x and y have q-gram similarity  $\tau_q(m, k) \ge m - q + 1 - kq$ .

<sup>&</sup>lt;sup>3</sup> Thus also to Hamming distance.

*Proof.* By induction on k.

### Example 1.5.

How can we use this result to solve a.s.m. problems? The lemma itself does not give us the direct solution, indeed it considers the edit distance between two arbitrary strings, while in a.s.m. problems the pattern can match any substring of the text. In the case of Hamming distance, if the pattern matches any substring s of t, then s must have length m. In the case of edit distance, it must hold  $m-k \le |s| \le m+k$ . The dot-plot representation helps us to visualize this concept. Hamming distance occurrences cover one single diagonal of the dot-plot, while edit distance occurrences are enclosed inside a parallelogram of side 2k+1.

Overlapping parallelograms?

We can design a filtration algorithm that scans the text and counts how many q-grams of the pattern falls into each parallelograms. Only the parallelograms exceeding the threshold  $\tau_q(m,k)$  have to be verified with an online method, e.g. standard DP. To speed up the filtration phase, the q-grams can be counted with the help of a substring index.

Which length of q makes filtration lossless?  $q = \left\lfloor \frac{m}{k+1} \right\rfloor$ .

*Figure 1.6:* Filtration with q-grams.

### **Seed filters**

Again, we start by considering the case of two arbitrary strings x, y s.t.  $d_E(x, y) \le k$ . If we partition w.l.o.g. y into k+1 non-overlapping seeds, then at least one seed  $y^i$  will occur as a factor of x.

**Lemma 1.2.** ? If  $y = y^1 y^2 ... y^{k+1}$  then  $x = a y^i b$  for some a, b.

Proof. Again, by induction on k?

Thus, we solve k-differences by partitioning p into k+1 seeds and searching them into t, e.g. with the help of a substring index. Note that we are reducing one approximate search into many smaller exact searches. Again, this lemma gives us a necessary but not sufficient condition, consequently we must verify whether any candidate location induced by an occurrence of some seed corresponds to an approximate occurrence of p in t. Thus, we verify any substring s of the text of length  $m-k \le |s| \le m+k$  containing one seed of p.

Which length of *q* makes filtration lossless?

Figure 1.7: Filtration with exact seeds.

### **Analysis of filters**

## 1.4 Related problems

### 1.4.1 Dictionary search

Dictionary search is a restriction of string matching. Given a set of database strings  $\mathbb D$  and a query string q find all strings in  $\mathbb D$  within distance k from q. Note that strings in  $\mathbb D$  usually have length similar to |q|, as  $||d|-|q||\leq k$  is a necessary condition for  $d_E(d,q)\leq k$ .

### Online methods

The problem can be solved by checking whether  $d_E(d,q) \leq k$  for all  $d \in \mathbb{D}$ . Answering the question whether the distance  $d_E(d,q) \leq k$  is an easier problem than computing the edit distance  $d_E(d,q)$ : a band of size k+1 is sufficient.

#### Indexed methods

Using a radix tree  $\mathcal{D}$  we can find all strings in  $\mathbb{D}$  equal to a query string q, in optimal time  $\mathcal{O}(|q|)$  and independently of  $||\mathbb{D}||$ .

**Algorithm 1.5** Exact dictionary search on a radix trie.

**Algorithm 1.6** Approximate dictionary search on a radix trie.

### Filtering methods

### 1.4.2 Overlaps computation

Define problem.

### Online methods

DP solution.

#### **Indexed methods**

Indexed solution, exact and approximate.

### 1.4.3 Local similarity search

Define score and scoring scheme.

Define local similarity.

### Online methods

Give dynamic programming solution.

### **Indexed methods**

Backtracking over substring index. BWT-SW.

## Filtering methods

SWIFT/Stellar is based on the q-gram lemma.