

# The Viterbi Algorithm

In this section we will describe the Viterbi algorithm in more detail. The Viterbi algorithm provides an efficient way of finding the most likely state sequence in the maximum *a posteriori* probability sense of a process assumed to be a **finite-state discrete-time** Markov process. Such processes can be subsumed under the general statistical framework of **compound decision theory**. We proceed to explain this theory in the context of Text Recognition as defined in [3], since this is the application that we will focus on later on.

## Compound decision theory

Suppose we have a text of  $n$  characteres. Each character yields a feature vector  $z_i$ ,  $i=1,2,...,n$ . Let  $p(\mathbf{Z}|\mathbf{C})$  denote the probability density function of the vector sequence  $\mathbf{Z}=z_1,z_2,...,z_n$  conditioned on the sequence of identities  $\mathbf{C}=c_1,c_2,...,c_n$ , where  $z_k$  is the feature vector for the  $k$ -th character, and where  $c_k$  takes on  $M$  values (number of letters in the alphabet) for  $k=1,2,...,n$ . Also, let  $P(\mathbf{C})$  be the *a priori* probability of the sequence of values  $\mathbf{C}$ . In other words  $P(\mathbf{C})$  is the *a priori* probability distribution of all sequences of  $n$  characters. The probability of correctly classifying the text is maximized by choosing that sequence of characters that has a maximum posterior probability or the so called: **maximum a posteriori (MAP)** probability, given by  $P(\mathbf{C}|\mathbf{Z})$ .

From **Bayes' rule** we obtain

$$P(\mathbf{C}|\mathbf{Z}) = \frac{p(\mathbf{Z}|\mathbf{C}) P(\mathbf{C})}{p(\mathbf{Z})}. \quad (1)$$

Since  $p(\mathbf{Z})$  is independent of the sequence  $\mathbf{C}$  (it is just a scale factor) we need only maximize the discriminant function

$$g_c(\mathbf{Z}) = p(\mathbf{Z}|\mathbf{C}) P(\mathbf{C}). \quad (2)$$

The amount of storage required for these probabilities is huge in practice, for that reason, assumptions are made in order to reduce the problem down to manageable size. These assumptions are:

- ⚡ **The size of the sequence of observations is not very large.** Let  $n$  be the size of a word. Then  $P(\mathbf{C})$  is the frequency of occurrence of words.
- ⚡ **Conditional independence among the features vectors.** The shape of a character, which generates a given feature vector, is independent of the shapes of neighbouring characters and is, therefore, dependent only on the character in question.

Under these assumptions, and taking logarithms, Eq. 2 reduces to

$$g_c(\mathbf{Z}) = \sum_{i=1}^n \log p(z_i | c_i) + \log P(c_1, ..., c_n). \quad (3)$$

For the case of the Viterbi algorithm, if we assume that the process is first-order Markov, then Eq. 3 reduces to:

$$g_c(\mathbf{Z}) = \sum_{i=1}^n \log p(z_i | c_i) + \log [P(c_1 | c_0) + P(c_2 | c_1) + ... + P(c_n | c_{n+1})]. \quad (4)$$

The **MAP** sequence estimation problem previously stated can also be viewed as the problem of finding the **shortest route** [1] through a certain graph. By thinking in this similarity one can see the natural recursivity of the Viterbi algorithm.

To illustrate how the Viterbi algorithm obtains this shortest path, we need to represent the Markov process in an easy way. A **state diagram**, like one shown in Fig. 2, is often used. In this state diagram, the **nodes** (circles) represent **states**, **arrows** represent **transitions**, and over the course of time the process traces some path from state to state through the state diagram.

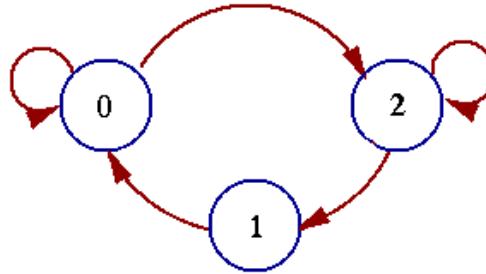


Fig. 2 State diagram of a three-state process.

A more redundant description of the same process is shown in Fig. 3, this description is called **trellis**. In a trellis, each node corresponds to a distinct state at a given time, and each arrow represents a transition to some new state at the next instant of time. The trellis begins and ends at the known states  $c_0$  and  $c_n$ . Its most important property is that to every possible state sequence  $\mathbf{C}$  there corresponds a unique path through the trellis, and vice versa.

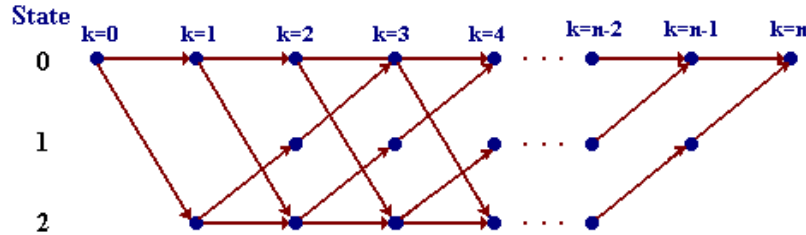


Fig. 3 Trellis for the three-state process of Fig. 1.

Now, suppose we assign to every path a length proportional to  $-\log [p(\mathbf{Z}|\mathbf{C})P(\mathbf{C})]$ . Since  $\log()$  is a **monotonic** function and there is a **one-to-one** correspondence between paths and sequences, we only need to find the path whose  $-\log [p(\mathbf{Z}|\mathbf{C})P(\mathbf{C})]$  is **minimum**, this will give us the state sequence for which  $p(\mathbf{Z}|\mathbf{C})P(\mathbf{C})$  is **maximum**, in other words, the state sequence with the maximum *a posteriori* (MAP) probability, which take us back to the original problem we want to solve. The **total length** of the path corresponding to some state sequence  $\mathbf{C}$  is

$$-\log [p(\mathbf{Z}|\mathbf{C})P(\mathbf{C})] = \sum_{k=1}^n l(t_k),$$

where  $l(t_k)$  is the associated length to each transition  $t_k$  from  $c_k$  to  $c_{k+1}$ . The shortest such path segment is called the **survivor** corresponding to the node  $c_k$ , and is denoted  $\mathbf{S}(c_k)$ . For any time  $k > 0$ , there are  $M$  survivors in all, one for each  $c_k$ . The observation is this: the shortest complete path  $\mathbf{S}$  must begin with one of these survivors. Thus for any time  $k$  we need to remember only the  $M$  survivors  $\mathbf{S}(c_k)$  and their correspondent lengths. To get to time  $k+1$ , we need only extend all time- $k$  survivors by one time unit, compute the lengths of the extended path segments, and for each

node  $c_{k+1}$  as the corresponding time-( $k+1$ ) survivor. Recursion proceeds indefinitely without the number of survivors ever exceeding  $M$ . This algorithm is a simple version of forward dynamic programming.

We illustrate this with an example taken from [1] that involves a four-state trellis covering five time units. The complete trellis with each branch labeled with a length is shown in Fig. 4.

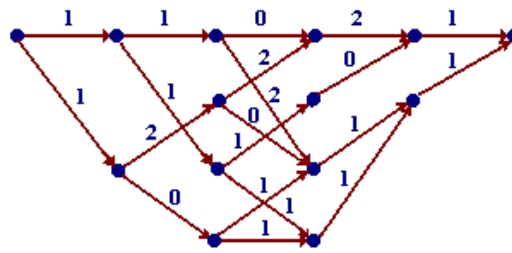


Fig. 4 Trellis labeled with branch lengths;  $M=4$ ,  $n=5$

The 5 recursive steps by which the algorithm determines the shortest path from the initial to the final node are shown in Fig. 5. At each step only the 4 (or fewer) survivors are shown, along with their lengths.

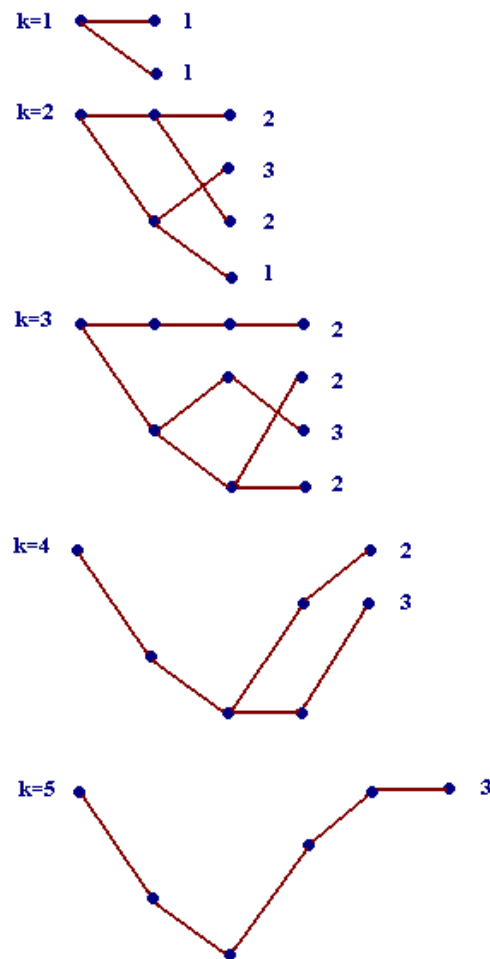


Fig. 5 Recursive determination of the shortest path via the Viterbi algorithm.

The Viterbi algorithm seen as finding the shortest route through a graph is:

**Input:**

$\mathbf{Z}=z_1, z_2, \dots, z_n$  the input observed sequence

**Initialization:**

$k=1$  time index

$S(c_1)=c_1$

$L(c_1)=0$  this is a variable that accumulate the lengths, the initial length is 0

**Recursion:**

For all transitions  $t_k=(c_k, c_{k+1})$

compute:  $L(c_k, c_{k+1}) = L(c_k) + l[t_k = (c_k, c_{k+1})]$  among all  $c_k$ .

Find  $L(c_{k+1}) = \min L(c_k, c_{k+1})$

For each  $c_{k+1}$

store  $L(c_{k+1})$  and the corresponding survivor  $S(c_{k+1})$ .

$k=k+1$

Repeat until  $k=n$

With finite state sequences  $\mathbf{C}$  the algorithm terminates at time  $n$  with the shortest complete path stored as the survivor  $S(c_K)$ .

The complexity of the algorithm is easily estimated:

**Memory:** the algorithm requires  $M$  storage locations, one for each state, where each location must be capable of storing a length  $L(m)$  and a truncated survivor listing  $S(m)$  of the symbols.

**Computation:** in each unit of time the algorithm must make  $M^2$  additions at most, one for each existing transition, and  $M$  comparisons among the  $M^2$  results.

Thus, the amount of storage is proportional to the number of states, and the amount of computation to the number of transitions.

