Dynamic Programming for Reduced NFAs for Approximate String and Sequence Matching[1]

Jan Holub

Department of Computer Science and Engineering
Czech Technical University
Karlovo nám. 13, 121 35 Prague 2, Czech Republic
phone: (++420 2) 2435 7287, fax: (++420 2) 298098

e-mail: `holub@cs.felk.cvut.cz`

Abstract. We present a new simulation method for the reduced nondeterministic finite automata (*NFA*s) for the approximate string and sequence matching using the Levenshtein and generalized Levenshtein distances. These reduced *NFA*s are used in case that we are interested only in all occurrences of a pattern in an input text such that the edit distance between the pattern and the found strings is less or equal to a given k and we are not interested in the values of these edit distances. The presented simulation method is based on the dynamic programming.

Key words: approximate string and sequence matching, simulation of nondeterministic finite automata, Levenshtein distance, generalized Levenshtein distance, dynamic programming

1 Introduction

Given a string $T = t_1 t_2 \ldots t_n$ over an alphabet Σ, a pattern $P = p_1 p_2 \ldots p_m$ over the alphabet Σ, and an integer k, $k \leq m \leq n$. The approximate string matching is defined as a searching for all occurrences of pattern P in text T such that edit distance $D(P, X)$ between pattern P and string $X = t_i t_{i+1} \ldots t_j$, $0 < i \leq j \leq n$, found in the text is less than or equal to k. The approximate sequence matching is defined in the same way as the approximate string matching, but any number of symbols can be located between the occurrences of two adjacent symbols of the pattern in the text. In this paper we consider two types of distances called the Levenshtein distance and the generalized Levenshtein distance.

The Levenshtein distance $D_L(P, X)$ between strings P and X not necessarily of the same length is the minimum number of edit operations *replace* (one character is replaced by another), *insert* (one character is inserted), and *delete* (one character is removed) needed to convert string P to string X. The generalized Levenshtein distance $D_G(P, X)$ between strings P and X not necessarily of the same length is the minimum number of edit operations *replace*, *insert*, *delete*, and *transpose* (two adjacent characters are exchanged) needed to convert string P to string X.
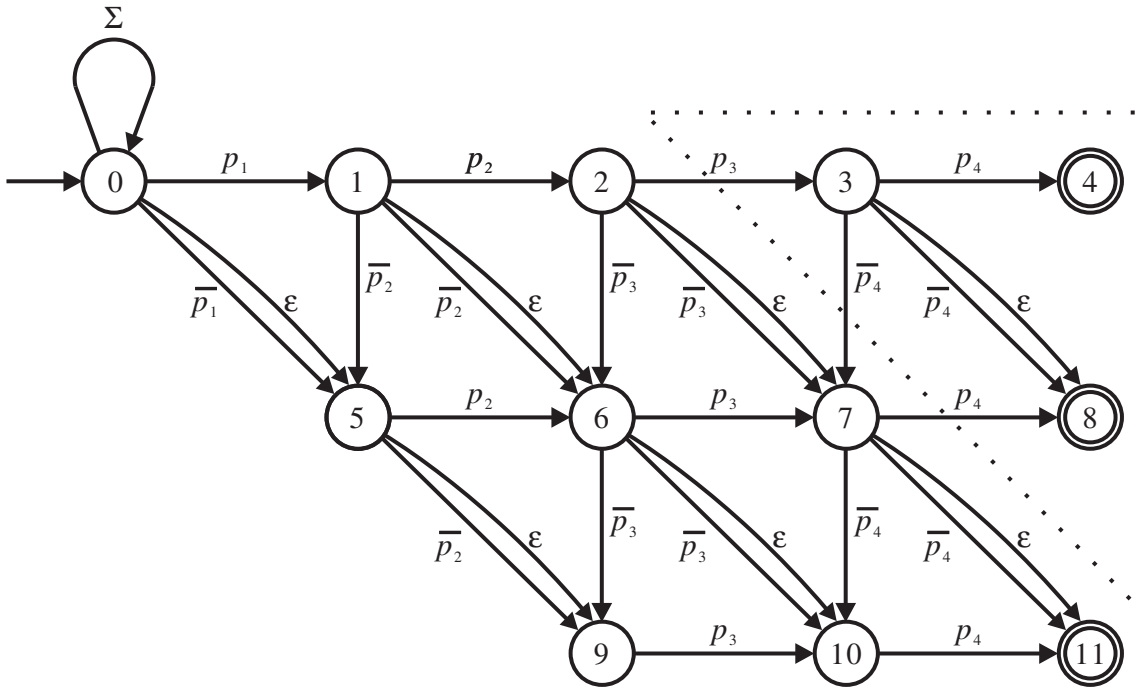

Figure 1: *NFA* for the approximate string matching using the Levenshtein distance $(m = 4, k = 2)$.

The nondeterministic finite automaton (*NFA*) for the approximate string matching using the Levenshtein distance has been presented in [Mel96, Hol96]. In the *NFA* there is for each edit distance l, $0 \leq l \leq k$, one level of states. An example of such *NFA* for $m = 4$ and $k = 2$ is shown in Figure 1[2].

There are known two algorithms for the approximate string matching for which there was shown [Mel96, Hol97] that they simulate the run of the *NFA* for the approximate string matching. The first method is Shift-Or algorithm [BYG92] and its variations — Shift-Add [BYG92] and Shift-And [WM92]. The second method is the dynamic programming [Sel80, Ukk85].

2   Reduced *NFA*s

If we are interested only in all occurrences of the pattern in the text with the edit distance less or equal to k, and we do not want to know the edit distance between the found string and the pattern, we can remove such states from the *NFA* for the approximate string matching that are needed only to determine the edit distance of the found string [Hol96]. Such states are bordered by the dotted line in Figure 1. The resulting *NFA* is shown in Figure 2 and has only one final state that represents that the pattern has been found with the edit distance less or equal to k.

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[2]Symbol $\overline{p}_j$, $0 < j \leq m$, represents $\Sigma - \{p_j\}$ in figures.
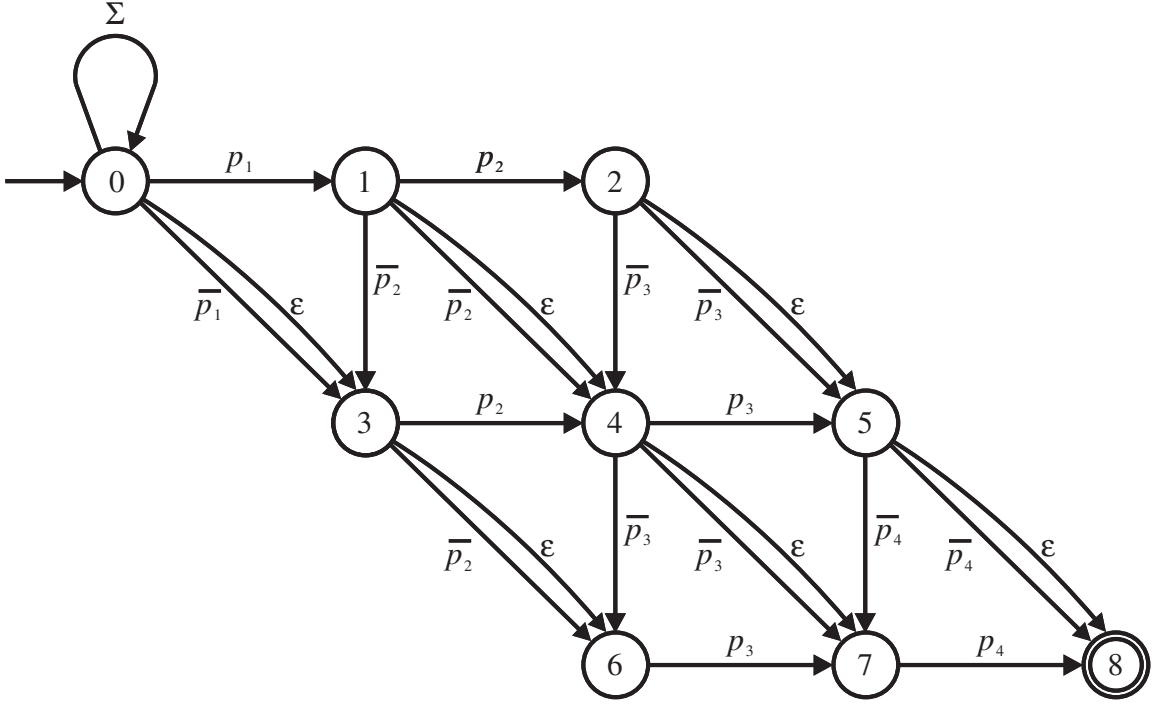
Figure 2: Reduced *NFA* for the approximate string matching using the Levenshtein distance ($m = 4$, $k = 2$).

The modification of Shift-Or algorithm for the reduced *NFA*s was presented in [Hol96] and the modification of the dynamic programming is discussed in the following sections.

3   Dynamic Programming

The dynamic programming [Sel80, Ukk85] computes in each step i of the run of the *NFA* ith column of matrix D which is of size $(m \times n)$; one element of the column is for each depth of the *NFA* and contains the number of level of the highest active state of this depth. If there is no active state in this depth, then the element contains the number of the level not existing in this depth. Since each *NFA* for the approximate string matching has $m + 1$ depths, it needs space $\mathcal{O}(m)$ and runs in time $\mathcal{O}(mn)$.

Since last k depths of the *NFA* do not have states on all $k + 1$ levels of the *NFA*, this method is not suitable for the reduced *NFA*s for the approximate string matching. Instead of having one element of the column for each depth of the *NFA* we have one element for each diagonal of the *NFA*; these diagonals are formed by the ε-transitions and are of the same length. If any state on a diagonal is active, then all states located lower on this diagonal are also active because of ε-transitions. Therefore in the element for each diagonal l, $0 \leq l \leq m - k$, we store only the number of the level of the highest active state on diagonal l. In this way we get for each step i, $0 \leq i \leq n$, of the run of the *NFA* the column $D_i = d_{0,i}, d_{1,i}, \ldots, d_{m-k,i}$ of length $m - k + 1$. Each element of the column can contain a value ranging from 0 to $k + 1$, where value $k + 1$

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represents that there is no active state on the corresponding diagonal. The formula for computing columns D_i is as follows:

$$
\begin{aligned}
d_{0,i} \;&:=\; 0, && 0 \le i \le n \\
d_{j,0} \;&:=\; k+1, && 0 < j \le m-k \\
d_{j,i} \;&:=\; \min(k+1, \\
& \qquad g_{d_{j-1,i-1}+j,t_i} + d_{j-1,i-1}, && \text{delete \& match} \\
& \qquad \textbf{if } p_{d_{j,i-1}+j+1} \ne t_i \\
& \qquad \textbf{then } d_{j,i-1}+1 && \text{replace} \\
& \qquad \textbf{else } k+1, \\
& \qquad \textbf{if } p_{d_{j+1,i-1}+j+2} \ne t_i \\
& \qquad \textbf{then } d_{j+1,i-1}+1 && \text{insert} \\
& \qquad \textbf{else } k+1), && 0 < j < m-k, 0 < i \le n \\
d_{j,i} \;&:=\; \min(k+1, \\
& \qquad g_{d_{j-1,i-1}+j,t_i} + d_{j-1,i-1}, && \text{delete \& match} \\
& \qquad \textbf{if } p_{d_{j,i-1}+j+1} \ne t_i \\
& \qquad \textbf{then } d_{j,i-1}+1 && \text{replace} \\
& \qquad \textbf{else } k+1), && j = m-k, 0 < i \le n
\end{aligned}
\tag{3}
$$

The first line in the formula says that the initial state lying on the 0th diagonal of the *NFA* is always active because of its self-loop.

The second one says that at the beginning of the searching there is no active state on diagonals l, $0 < l \le m-k$, because there is no initial state on such diagonals.

Part $g_{d_{j-1,i-1}+j,t_i} + d_{j-1,i-1}$ represents *match* and *delete* transitions. The *match* is represented by the horizontal transitions and edit operation *delete* is represented by the diagonal ε-transitions in Figure 2. An implementation of *match* transition is simple — if the state on diagonal $j-1$ and on level $d_{j-1,i-1}$ is active and horizontal transition leading from this state is labeled by symbol t_i, then the state on diagonal j and on level $d_{j-1,i-1}$ becomes active. For an implementation of *delete* transition we have to search for the state on diagonal $j-1$ and on level l, $d_{j-1,i-1} \le l \le m-k$, such that there is a *match* transition labeled by input symbol t_i leading from this state. In order to find such state in the constant time we have to use auxiliary matrix G in which there is for each position r in pattern P and input symbol t_i the number r', $0 \le r'$, such that $p_{r+r'} = t_i$ where r' is the lowest possible. If there is no such position, then $r' = k+1$. Since the value of $d_{j-1,i-1}$ can be $k+1$ and the maximum number of diagonal, into which there lead *match* transitions, is $m-k$, the maximum position for which a value of matrix G is required is $m-k+k+1 = m+1$. Therefore the matrix has to be of size $(m+2) \times |\Sigma'|$ where $\Sigma' \subseteq \Sigma$ is the alphabet used in pattern P. The formula for computation of matrix G is as follows:

$$
\begin{aligned}
g_{j,a} \;&:=\; \min(\{k+1\} \cup \{(l \mid p_{j+l} = a, 0 \le l) \text{ or} \\
& \qquad (k+1 \mid \text{if there is no such } l)\}), && 0 < j \le m, a \in \Sigma \\
g_{m+1,a} \;&:=\; k+1, && a \in \Sigma
\end{aligned}
\tag{4}
$$

Number $d_{j-1,i-1}+j$ gives the position of symbol $p_{d_{j-1,i-1}+j}$ in the pattern which is used as a label of the *match* transition leading from the highest active state on

diagonal $j - 1$ to a state on diagonal j. Therefore $g_{d_{j-1,i-1}+j,t_i} + d_{j-1,i-1}$ gives the level of the highest active state on diagonal j that has arisen by using *match* transition to each active state on diagonal $j - 1$.

Part $d_{j,i-1} + 1$ represents *replace* transition. In Figure 2, edit operation *replace* is represented by the diagonal transition labeled by symbol $\overline{p}_{d_{j,i-1}+j+1}$ mismatching symbol $p_{d_{j,i-1}+j+1}$. To implement *replace* transition it is only needed to move the highest active state on diagonal j to the next lower position on the same diagonal. Since $d_{j,i-1}$ can reach $k + 1$ the value of expression $d_{j,i-1} + j + 1$ can be greater than m and in that case $p_{d_{j,i-1}+j+1}$ would give undefined value. To solve this problem we can add some **if** statements but it increases the time of the computation. The better solution is to put some symbols, that are not in input alphabet Σ, at positions $m + 1$ and $m + 2$ of the pattern — for example symbol $\langle end\ of\ string \rangle$.

Part $d_{j+1,i-1} + 1$ represents *insert* transition. In Figure 2, edit operation *insert* is represented by the vertical transition also labeled by mismatching symbol $\overline{p}_{d_{j+1,i-1}+j+2}$. The active state on diagonal $j + 1$ and on level $d_{j+1,i-1}$ moves to level $d_{j+1,i-1} + 1$ on diagonal j.

From these transitions we get minimum in order to obtain the highest active state on each diagonal. An example of matrix G for pattern $P = adbbca$ is shown in Table 1 and the process of searching for pattern $P = adbbca$ with at most $k = 3$ errors in text $T = adcabcaabadbbca$ is shown in Table 2.

G	a	b	c	d	$\Sigma - \{a, b, c, d\}$
1	0	2	4	1	4
2	4	1	3	0	4
3	3	0	2	4	4
4	2	0	1	4	4
5	1	4	0	4	4
6	0	4	4	4	4
7	4	4	4	4	4

Table 1: Matrix G for pattern $P = adbbca$ and $k = 3$.

D	-	a	d	c	a	b	c	a	a	b	a	d	b	b	c	a
0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
1	4	0	1	1	0	1	2	0	0	1	0	1	2	2	3	0
2	4	4	0	1	2	1	2	3	4	1	2	0	1	2	3	4
3	4	4	4	2	3	4	2	3	3	4	3	4	0	1	2	3

Table 2: Matrix D for pattern $P = adbbca$, text $T = adcabcaabadbbca$, and $k = 3$.

Below we also present an algorithm that uses the dynamic programming for the reduced *NFA* for the approximate string matching using the Levenshtein distance. While in Formula 3 there were evaluated the transitions incoming to the diagonals,

in this algorithm there are evaluated the outgoing transitions. It simplifies the computation because then there is only one test whether input symbol t_i is a matching symbol. This test is necessary for deciding whether to use only *match* transition or to use *replace*, *insert*, and *delete* transitions.

Algorithm 1
DP for the reduced *NFA* for the approximate string matching using the Levenshtein distance
Input: Pattern $P = p_1 p_2 \ldots p_m$, text $T = t_1 t_2 \ldots t_n$, maximum number of differences allowed k.
Output: Matrix D of size $(m - k + 1) \times (n + 1)$.
Method:

$$d_{0,0} := 0$$
$$d_{j,0} := k + 1, 0 < j \le m - k$$
for $i := 1, 2, \ldots, n$ **do**
$\quad d_{0,i} := 0$ $\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* $j = 0$ */
$\quad d_{1,i} := g_{1,t_i}$ $\qquad\qquad\qquad$ /* *delete* & *match* from the initial state *** */
$\quad$ **if** $p_{d_{1,i-1}+2} = t_i$ **then** $\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* $j = 1$ */
$\quad\quad d_{2,i} := d_{1,i-1}$ $\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* *match* */
$\quad$ **else**
$\quad\quad d_{2,i} := \min(g_{d_{1,i-1}+2,t_i} + d_{1,i-1}, k + 1)$ $\qquad\qquad$ /* *delete* & *match* */
$\quad\quad d_{1,i} := \min(d_{1,i-1} + 1, d_{1,i})$ $\qquad\qquad\qquad\qquad$ /* *replace* */
$\quad$ **endif** $\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* *** */
$\quad$ **for** $j := 2, 3, \ldots, m - k - 1$ **do**
$\quad\quad$ **if** $p_{d_{j,i-1}+j+1} = t_i$ **then**
$\quad\quad\quad d_{j+1,i} := d_{j,i-1}$ $\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* *match* */
$\quad\quad$ **else**
$\quad\quad\quad d_{j+1,i} := \min(g_{d_{j,i-1}+j+1,t_i} + d_{j,i-1}, k + 1)$ $\qquad$ /* *delete* & *match* */
$\quad\quad\quad d_{j,i} := \min(d_{j,i-1} + 1, d_{j,i})$ $\qquad\qquad\qquad\qquad$ /* *replace* */
$\quad\quad\quad d_{j-1,i} := \min(d_{j,i-1} + 1, d_{j-1,i})$ $\qquad\qquad\qquad$ /* *insert* */
$\quad\quad$ **endif** $\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* *** */
$\quad$ **endfor**
$\quad j := m - k$ $\qquad\qquad\qquad\qquad\qquad\qquad\qquad$ /* the last diagonal */
$\quad$ **if** $p_{d_{j,i-1}+j+1} \ne t_i$ **then**
$\quad\quad d_{j,i} := \min(d_{j,i-1} + 1, d_{j,i})$ $\qquad\qquad\qquad\qquad$ /* *replace* */
$\quad\quad d_{j-1,i} := \min(d_{j,i-1} + 1, d_{j-1,i})$ $\qquad\qquad\qquad$ /* *insert* */
$\quad$ **endif**
$\quad$ **if** $d_{m-k,i} < k + 1$ **then**
$\quad\quad$ **write**("pattern found at position i")
$\quad$ **endif**
endfor

The first command in the first **for** cycle in the algorithm ($d_{0,0} := 0$) represents the self-loop of the initial state — the highest active state in 0th diagonal is always in level 0 and this is the initial state.

The second command ($d_{1,i} := g_{1,t_i}$) represents the only transition that leads from 0th diagonal which is *match* transition. g_{1,t_i} gives the position l of the pattern, on which t_i is located, or $k + 1$ if t_i is not in the pattern. If $l < k + 1$, then this position l

is equal to the level of 1st diagonal in which there is the active state that arose by using *match* transition for t_i going from 0th diagonal.

The first **if** statement represents transitions leading from the highest active state on the 1st diagonal. In this case we do not evaluate *insert* transitions because they lead always to 0th diagonal where the initial state is always active. If input symbol t_i is the same as the symbol $p_{d_{1,i-1}+2}$ used as a label of *match* transition leading from the highest active state in 1st diagonal, then we evaluate only this *match* transition ($d_{2,i} := d_{1,i-1}$). If the symbols are different, then we evaluate *delete* and *replace* transitions. For *delete* transition we search for the next occurrence of input symbol t_i in the pattern behind position $d_{j,i-1} + j + 1$ (the number of the diagonal plus the number of the level gives the position in the pattern corresponding to the state on that level of that diagonal). At first we perform *delete* transition (we move the highest active state down in the diagonal) and then we perform *match* transition for input symbol t_i. For *replace* transition we move the highest active state in the diagonal to the next lower position in the diagonal.

In the next **for** cycle the transitions leading from the highest active state of the next diagonals except the last one are evaluated. It is done in the same way as described in the previous paragraph but in addition *insert* transition is evaluated. For this *insert* transition we put the level of diagonal j increased by one to the previous diagonal $j - 1$.

In the last diagonal we evaluate only *replace* and *insert* transitions because *match* transition has no diagonal into which it could lead.
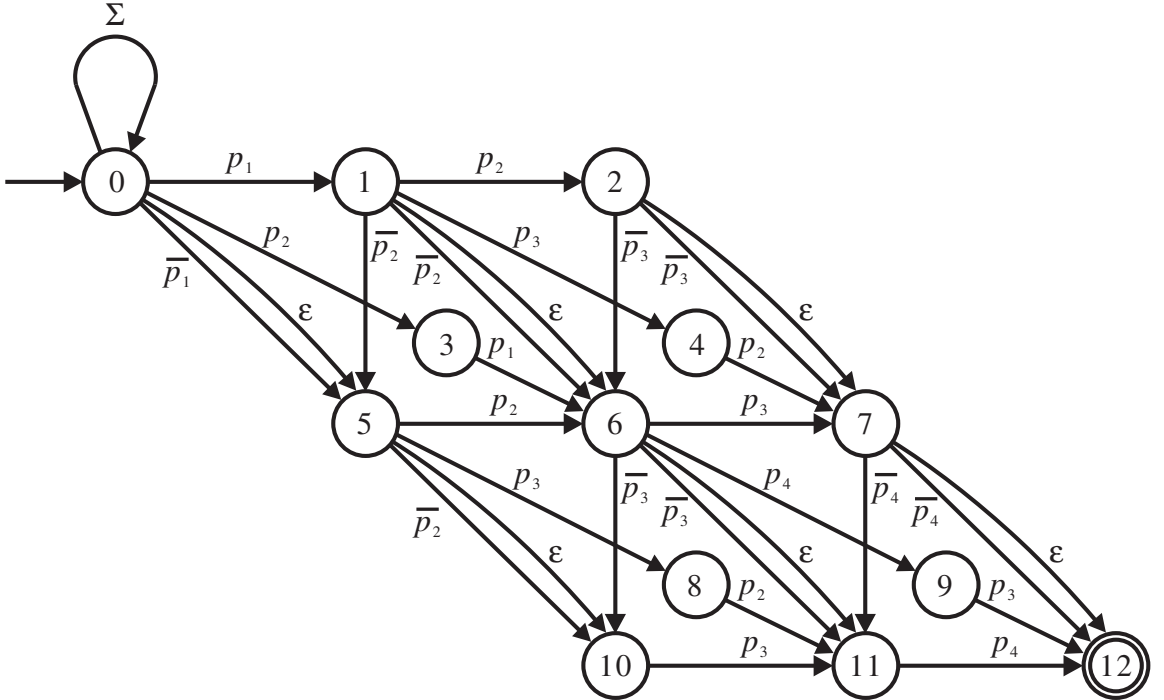

Figure 3: Reduced *NFA* for the approximate string matching using the generalized Levenshtein distance ($m = 4$, $k = 2$).

This method can be also used for the simulation of the run of the reduced *NFA*

for the approximate string matching using the generalized Levenshtein distance. An example of such reduced *NFA* for $m = 4$ and $k = 2$ is shown in Figure 3. We have only to add the part representing edit operation *transpose*. In Formula (3), the added part is as follows:

$$\begin{array}{ll} \textbf{if } p_{d_{j-1,i-2}+j+1} = t_{i-1} \textbf{ and } p_{d_{j-1,i-2}+j} = t_i & \\ \textbf{then } d_{j-1,i-2} + 1 & transpose \\ \textbf{else } k + 1, & 0 < j \le m - k, 1 < i \le n \end{array} \qquad (5)$$

And in Algorithm 1, the added part is as follows:

$$\begin{array}{ll} \textbf{if } p_{d_{j,i-2}+j+2} = t_{i-1} \textbf{ and } p_{d_{j,i-2}+j+1} = t_i \textbf{ then} & \\ \quad d_{j+1,i} := \min(d_{j,i-2} + 1, d_{j+1,i}) & /* \; transpose \; */ \\ \textbf{endif} & \end{array}$$

This part should be inserted into each part of Algorithm 1 where $0 \le j < m - k$ and $1 < i \le n$. Such places are behind the lines marked by '***'.
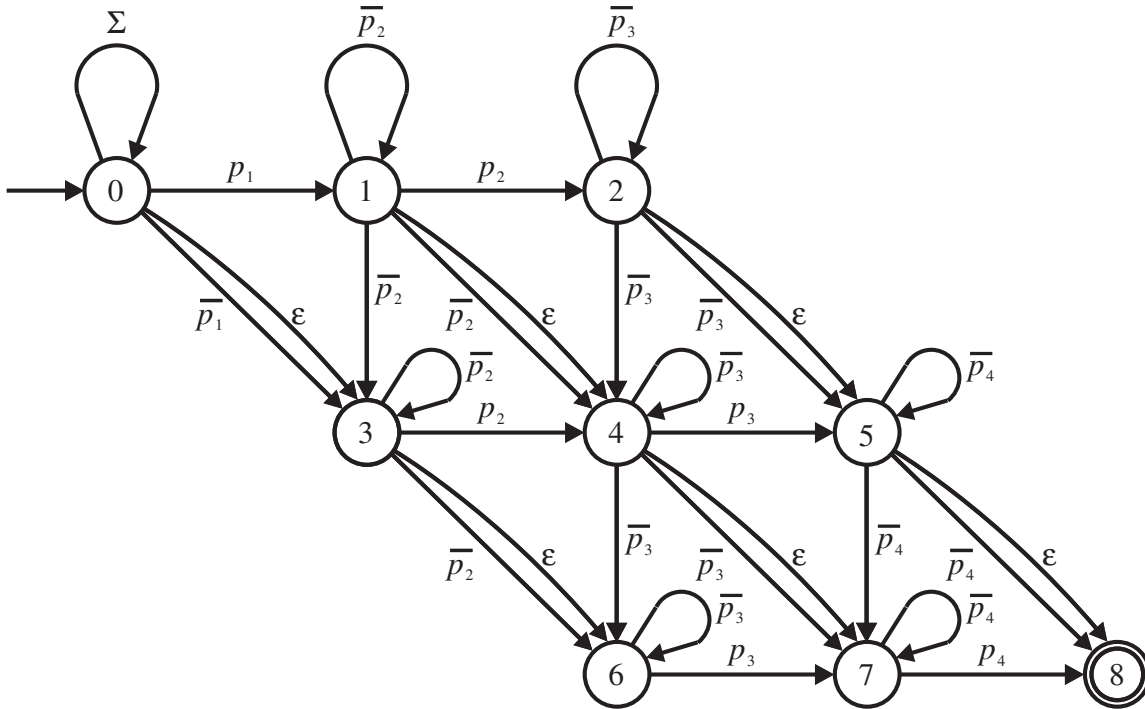

Figure 4: Reduced *NFA* for the approximate sequence matching using the Levenshtein distance ($m = 4$, $k = 2$).

This type of simulation of the reduced *NFA*s can be also used for the reduced *NFA*s for the approximate sequence matching using the Levenshtein and generalized Levenshtein distances [Hol97]. An example of the reduced *NFA* for the approximate

sequence matching using the Levenshtein distance for $m = 4$ and $k = 2$ is shown in Figure 4.

To modify the presented algorithm so that it could simulate this reduced *NFA* we have to implement the self-loops in each nonfinal and noninitial state. It can be performed by inserting the following part into Formulae (3) and (3+5) for the approximate string matching.

$$
\begin{array}{lll}
\textbf{if } p_{d_{j,i-1}+j+1} \neq t_i & & \\
\textbf{then } d_{j,i-1} & self-loop & \\
\textbf{else } k+1, & 0 < j < m - k, 0 < i \leq n & \\
\textbf{if } p_{d_{j,i-1}+j+1} \neq t_i \textbf{ and } d_{j,i-1} < k & & (6)\\
\textbf{then } d_{j,i-1} & self-loop & \\
\textbf{else } k+1, & j = m - k, 0 < i \leq n &
\end{array}
$$

The presented formulae and algorithm compute whole matrix D but in the practice only two (three for the generalized Levenshtein distance) columns from this matrix are used in each step of the computation.

4   Conclusion

The resulting simulation runs in time $\mathcal{O}((m - k)n + m\mu)$ and needs space $\mathcal{O}(m\mu)$, where μ is the number of different symbols used in the pattern. We can decrease the space complexity by using another implementation of auxiliary matrix G but it increases the time complexity. Our algorithm also uses only one input symbol in each step of computation in case of the Levenshtein distance and two input symbols in case of the generalized Levenshtein distance.

The resulting algorithm has the time bound better than [Sel80, Ukk85] which runs in time $\mathcal{O}(mn)$ and for $k > \frac{m}{2}$ it has also the time bound better (not considering the preprocessing time) than [GP89] which runs in time $\mathcal{O}(kn + m \log \tilde{m})$ where $\tilde{m} = \min(m, |\Sigma|)$.

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