

## ON THE WORST CASE OF A MINIMAL SPANNING TREE ALGORITHM FOR EUCLIDEAN SPACE

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### Abstract.

This paper concerns the worst case running time of the minimal spanning tree algorithm presented by Bentley and Friedman.

For a set of  $N$  points in  $k$ -dimensional Euclidean space the worst case performance of the algorithm is shown to be  $\Theta(N^2 \log N)$ , for  $k \geq 2$  and  $\Theta(N^2)$ , for  $k = 1$ .

*Keywords:* minimal spanning tree, Euclidean space,  $k$ - $d$  tree.

### 1. Introduction.

We shall study the special minimal spanning tree (MST) problem where a set of  $N$  points in  $k$ -dimensional Euclidean space is to be connected in such a way that the sum of the lengths of the  $N-1$  straight line segments is minimal. The problem can be solved by the method of Bentley and Friedman [1]. We shall discuss the running time of this algorithm in the worst case. Nevalainen, Ernvall and Katajainen [2] gave a simplified version of the algorithm. It is easy to see that the worst case of these MST-algorithms is the same. So we shall consider only the simplified version.

The algorithm of [2] functions as follows. At first  $N$  single point subtrees called fragments are formed. After this, at each step, the fragment of *minimal size* is joined to the nearest fragment with a minimal line segment. The process continues until there is only one fragment left which is the MST. To facilitate the choice of the minimal distance a priority queue that stores the distance to the nearest neighbour for each point of a fragment is maintained for each fragment. Because points are added to the fragment in the course of the process, for some points their nearest neighbour is no longer outside the fragment. If this kind of distance is met on the top of the priority queue we must search the nearest neighbour of the point outside the fragment by means of a  $k$ - $d$  tree [1] and update the queue.

The  $k$ - $d$  tree is a binary tree in which each node represents a subcollection of the points. In an internal node it is told how the subcollection is partitioned into two disjoint subsets. At each node the partitioning is accomplished by dividing the subcollection at the median value of one of the coordinates. The root of the tree represents the entire point set and the leaf nodes contain the actual points. With each node a set of bounds is associated, which jointly form a rectilinearly oriented hyperrectangle in  $k$ -space, within which all the points of the particular subcollection must lie.

The nearest neighbour of a point is searched among the points outside the particular fragment by using the  $k$ - $d$  tree and in this paper this is called the *nearest outside neighbour search*. The search can be described recursively. For an internal node we first search the subtree in which the discriminator coordinate of the points is on the same side of the partition value as the coordinate of the given point. If any part of the hyperrectangle of the opposite subtree does intersect a ball centered at the given point, with radius equal to the current nearest outside neighbour distance, then that subtree must also be searched. Generally, for a leaf node, the distance from the given point is calculated but in case the point in the leaf node belongs to the same fragment as the given point the leaf node is rejected at once.

Zolnowsky [3] considered three different choices for the selection of discriminator coordinates. He showed how the performance of the  $k$ - $d$  tree depends on the choice in context of the usual nearest neighbour search. We shall apply his ideas and show that if the discriminator is the coordinate in which the points have the *greatest spread* in values, then the running time of our MST-algorithm is  $\Theta(N^2 \log N)$ <sup>1)</sup> in the worst case for  $k \geq 2$ . In [2] a one-dimensional point set was given for which the running time is  $\Omega(N^2)$ . We shall show that  $O(N^2)$  is the worst case upper bound for  $k = 1$ .

## 2. The worst case study.

### 2.1 The upper and lower bounds for $k \geq 2$ .

In the MST-algorithm the number of points to be merged into bigger fragments is at most  $(N/2) \log N$  [2]. In the worst case, for all of these points the nearest outside neighbour has to be searched. The worst thing that could

<sup>1)</sup> The symbols  $O$ ,  $\Omega$ ,  $\Theta$  for functions  $f$  and  $g$  denote:

- (i)  $f(x) = O(g(x))$  iff there exists a positive constant  $C$  such that for every  $x$   $f(x) \leq Cg(x)$  (upper bound).
- (ii)  $f(x) = \Omega(g(x))$  iff there exists a positive constant  $D$  such that for every  $x$   $f(x) \geq Dg(x)$  (lower bound).
- (iii)  $f(x) = \Theta(g(x))$  iff  $f(x) = O(g(x))$  and  $f(x) = \Omega(g(x))$  (tight bound).

happen is that the entire  $k$ - $d$  tree must be searched each time. So we have a straightforward upper bound  $O(N^2 \log N)$  for the worst case complexity of the algorithm.

Now let us give a point set in the two-dimensional space for which the running time of the MST-algorithm reaches the above upper bound. Let us assume the number of the points is a power of 2, say  $N = 2^s$  ( $s \geq 2$ ). We choose the first  $N/2$  points as

$$\left( s + \sum_{j=0}^{s-2} \left\lfloor \frac{i-1}{2^j} \right\rfloor, 0 \right), \quad \text{for } 1 \leq i \leq N/2$$

and the next  $N/2 - 1$  points as  $\left( 0, \frac{1}{2^{N-i}} \right)$ , for  $N/2 < i < N$ . Finally we choose the  $N$ th point as  $(0, N)$  (see Figure 1).

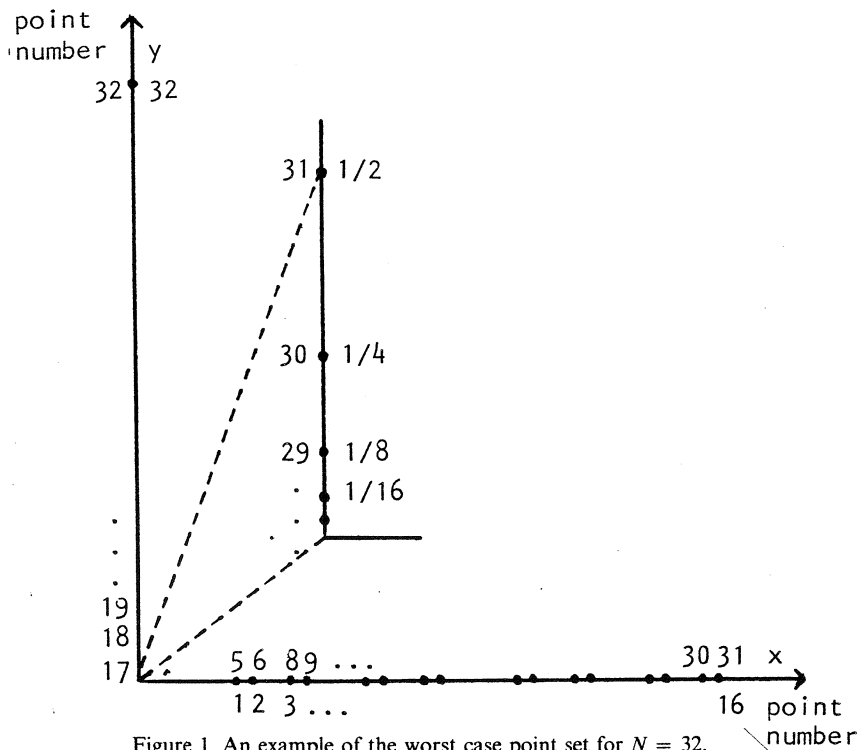


Figure 1. An example of the worst case point set for  $N = 32$ .

Because the point set on the  $y$ -axis is expressed with decreasing density all the points belong to the same fragment after the fragments of size one have been considered. On the  $x$ -axis we get first  $N/4$  fragments of equal size, then

$N/8$  fragments of equal size etc. We know that the number of the nearest outside neighbour searches is totally  $(N/4) \log(N/2) + N/2$  [2] because every priority in a priority queue of a fragment on the  $x$ -axis must be updated.

For these points the spread in the  $y$ -coordinate is one unit greater than that in the  $x$ -coordinate. Thus the discriminator of the root of the  $k$ - $d$  tree is the  $y$ -coordinate. The points on the  $y$ -axis and those on the  $x$ -axis are put into different subcollections. The points on the  $y$ -axis have zero spread in the  $x$ -coordinate. So the discriminator coordinate is the  $y$ -coordinate for each node in the subtree representing the points on the  $y$ -axis. Hence the search ball of each point on the  $x$ -axis must intersect the rectangle of each point on the  $y$ -axis.

The number of the nearest outside neighbour searches for this point set is  $\Omega(N \log N)$  each demanding  $\Omega(N)$  operations. So, the running time of the MST-algorithm is  $\Omega(N^2 \log N)$ . To sum up we know the algorithm has a worst case time complexity of  $\Theta(N^2 \log N)$  for  $k \geq 2$ .

## 2.2 The upper bound for $k = 1$ .

In one-dimensional space points are on a line. The fragments consist of adjacent points on the line. In the worst case during the search of the nearest neighbour of a fragment for each point of that fragment the nearest outside neighbour must be determined. It is further possible that in connection with each search all the leaf nodes containing a point of the fragment and two successive leaf nodes must be visited in the  $k$ - $d$  tree. Thus in the worst case,  $m+2$  adjacent leaf nodes are visited when the size of a fragment is  $m$ .

We now count the number of nodes visited in the  $k$ - $d$  tree during the search procedure. Since a constant time is spent in each node when searching the nearest neighbour, the number of nodes counted in this way indicates the complexity of the algorithm.

We denote the *closest common ancestor* of the fragment  $F$  ( $cca(F)$ ) by a node which is the root of the smallest possible subtree containing all of the adjacent  $m+2$  leaf nodes. Let  $F' \subseteq F$  be the subset which contains the points in the leaf nodes of the left subtree of  $cca(F)$  and  $F''$  the corresponding subset for the right subtree of  $cca(F)$ . The complete subtree, whose root is  $cca(F')$  is called  $A$  and the subtree whose root is  $cca(F'')$  is called  $B$ . In Figure 2 the situation is described in the case where  $cca(F)$  is the root of the whole tree.

The nodes visited during the search are on the path from the root of the  $k$ - $d$  tree to  $cca(F')$  and to  $cca(F'')$  and in the subtrees  $A$  and  $B$ . We shall first count the number of nodes on the path. This number is at most  $2h$  where  $h$  is the height of the  $k$ - $d$  tree. At most  $O(N \log N)$  nearest outside neighbour searches have to be done [2]. Thus the total number of nodes visited in order to reach both these two nodes is at most  $O(N \log^2 N)$ .

When we count the number of nodes in the subtrees  $A$  and  $B$  we get a rough



### 3. Refinements to the algorithm.

Instead of using the greatest spread partition criteria in the construction of the  $k$ - $d$  tree the so-called *square* criteria could give better running times in the worst case as seen in [3]. The idea is to choose the discriminator as the coordinate in which the separation of the bounds is currently greatest. However, in one-dimensional space the two  $k$ - $d$  tree structures are the same. Thus the point set presented in [2] is still valid in order to get  $\Omega(N^2)$  performance. This point set has just the property that during the nearest outside neighbour searching in the  $k$ - $d$  tree all the leaf nodes containing a point of that fragment must be visited. Additionally this search must be carried out for every point of the fragment.

An essential refinement is achieved for  $k = 1$  if during the search of the nearest outside neighbour an entrance to the branches for which every point belonging to the same fragment as the given point is not allowed. Bentley and Friedman called this "poisoning". Then a new label must be added to the nodes of a  $k$ - $d$  tree. The label tells the name of the fragment to which all the points of the corresponding subtree belong. If the left subtree and the right subtree of a node contain points from different fragments the label is undefined.

The updating of the labels takes  $O(N \log^2 N)$  time. For each point of a smaller fragment we must first search the leaf node of that point and then backtrack towards the root until the updating is no more possible. At most  $O(N \log N)$  update operations are performed each demanding  $O(\log N)$  time. After this refinement each nearest outside neighbour search takes no longer than  $O(2 \log N)$  time because of the special form of the fragments. The number of the nearest outside neighbour searches is at most  $O(N \log N)$  and so the algorithm takes  $O(N \log^2 N)$  time in the worst case. It is an open question whether the method could be applied successfully even to  $k > 1$ .

It should be pointed out that after this refinement the algorithm is not yet optimal, since  $O(N \log N)$  time is possible by sorting. Perhaps, it is not sensible to replace the search of the nearest neighbour of a fragment with many nearest outside neighbour searches. During the algorithm the nearest neighbour of a fragment must be searched  $N-1$  times and in the worst case this has to be replaced with  $(N/2) \log N$  nearest outside neighbour searches. It is worth studying if there exists a data structure which supports an efficient search of the nearest neighbour of a fragment.

We have seen that the priority queues became useless because it is possible to construct point sets for which the nearest outside neighbour search is needed to each point of the fragment until it is certain which point is the nearest. When an algorithm with a good performance in the worst case is formed, priority queues need not be used this way. However, the use of the priority queues is fundamental in the algorithm of Bentley and Friedman where the fragment is grown from the area with the lowest local density of points towards the points closer to each other.

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