Two Optimization Problems for Unit Disks

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- 5 2 FMF and FRI, University of Ljubljana, Slovenia
- Abstract
- 7 To be done

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Introduction

In this paper we consider two geometric optimization problems in the plane where unit disks play a prominent role. For both problems we discuss efficient algorithms to solve them, provide an implementation of these algorithms, and present experimental results on the implementation.

The first problem we consider is computing a shortest-path tree in the (unweighted) intersection graph of unit disks. The input to the problem is a set \mathcal{D} of n disks of the same size, each disk represented by its center. The corresponding unit disk (intersection) graph has a vertex for each disk, and an edge connecting two disks D and D' of \mathcal{D} whenever D and D' intersect. An alternative, more convenient point of view, is to take as vertex set the set of centers of the disks, denoted by P, and connecting two points p and q of P whenever the Euclidean length |pq| is at most the diameter of a disk. Given a root $r \in P$, the task is to compute a shortest-path tree from r in this graph.

The second problem we consider is the *minimum-separation problem*. The input is a set \mathcal{D} of n unit disks in the plane and two points s and t not covered by any disks of \mathcal{D} . We say that \mathcal{D} separates s and t if each curve in the plane from s to t intersects some disk of \mathcal{D} . The task is to find the minimum cardinality subset of \mathcal{D} that separates s and t. Formally, we want to solve

 $\begin{aligned} & \text{min} & |\mathcal{D}'| \\ & \text{s.t.} & \mathcal{D}' \subset \mathcal{D} \text{ and } \mathcal{D}' \text{ separates } s \text{ and } t. \end{aligned}$

Unit disks are the most standard model used for wireless sensor networks; see for example [6,8,16]. Often the model is referred as UDG. This model provides an appropriate trade off between simplicity and accuracy. Other models are more accurate, as for example discussed in [10,13], but obtaining efficient algorithm for them is much more difficult.

While unit disks give a simple model, exploiting the geometric features of the model is often challenging. Shortest paths in unit disk graphs are essential for routing and are a basic subroutine for several other more complex tasks. A somehow unexpected application of shortest paths in unit-disk graphs to boundary recognition is given in [15]. The minimum-separation problem and variants thereof have been considered in [2,7,14]. The problem is dual to the barrier-resilience problem considered in [1,9,11,12]. It is not obvious that the minimum-separation problem can be solved optimally in polynomial time, and the known algorithm for this uses as a subroutine shortest paths in unit disk graphs. Thus, both problems considered in this paper are related and it is worth to consider them together.

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Our contribution. We are aware of two algorithms to compute shortest-path trees in unit disk graphs in $O(n \log n)$ worst-case time: one by Cabello and Jejčič [3] and one Efrat, Itai and Katz [5]. Here we report on an implementation of a modification of the algorithm in [3], and compare it against two obvious alternatives. The only complex ingredient in the algorithm is computing the Delaunay triangulation, but efficient libraries are available for this. The algorithm of [5] would be substantially harder to implement and it has worse constants hidden in the O-notation.

As mentioned before, it is not obvious that the minimum-separation problem can be solved in polynomial time. A 2-approximation algorithm is given by Gibson, Kanade, and Varadarajan [7]. Cabello and Giannopoulos [2] provide an exact algorithm that takes $O(n^3)$ wosrt-case time and works for arbitrary shapes, not just disks. In this paper we improve this last algorithm to near-quadratic time for the special case of unit disks. The basic principle of the algorithm is the same, but several additional tools from Computational Geometry have to be employed to reduce the worst-case running time. We implement a variant of the new, near-quadratic-time algorithm and report on the experiments.

Assumptions. We will assume that $unit\ disk$ means that it has radius 1/2. Up to scaling the input data, this choice is irrelevant. However, it is convenient for the exposition because then the disks intersect whenever the distance between their centers is 1. The implementation and the experiments also make this assumption.

Henceforth P will be the set of centers of \mathcal{D} . All the computation will be concentrated on P. In particular, we assume that P is known. (For the shortest path problem, one could possibly consider weaker models based on adjacencies.)

We will work with the graph $G_{\leq 1}(P)$ with vertex set P and an edge between two points $p,q \in P$ whenever their Euclidean distance |pq| is at most 1. In the notation we remove the dependency on P and on the distance. Thus we just use G instead of $G_{\leq 1}(P)$. For simplicity of the theoretical exposition we will sometimes assume that G is connected. It is trivial to adapt to the general case, for example treating each connected component separately. The implementation does not make this assumption.

In the minimum-separation problem we will use s and t for the points to separate. We will assume that s is the origin and t is the point $(0, \tau)$, with $\tau \ge 0$. Thus, the segment st is vertical and t is above s. The implementation also makes this assumption. A simple rigid transformation can be applied to the input to get to this setting.

Organization of the paper. In Section 2 we discuss the theoretical algorithms for both problems and their guarantees. In Section 3 we discuss the implementations and in Section 4 we present our experimental results.

2 Description of algorithms

2.1 Shortest-path tree in unit-disk graphs

We describe here the algorithm of Cabello and Jejčič [3] to compute a shortest path tree in G from a given root point $r \in P$. As it is usually done for shortest path algorithms we use tables $dist[\cdot]$ and $\pi[\cdot]$ indexed by the points of P to record, for each point $p \in P$, the distance $d_G(s, p)$ and the ancestor of p in a shortest (s, p)-path.

The pseudocode of the algorithm, which we call UNWEIGHTEDSHORTESTPATH, is in Figure 1. We explain the intuition, taken almost verbatim from [3]. We start by computing the Delaunay triangulation DT(P) of P. We then proceed in rounds for increasing values of i, where at round i we find the set W_i of points at distance exactly i in G from the source r.

We start with $W_0 = \{r\}$. At round i, we use DT(P) to grow a neighbourhood around the points of W_{i-1} that contains W_i . More precisely, we consider the points adjacent to W_{i-1} in DT(P) as candidate points for W_i . For each candidate point that is found to lie in W_i , 81 we also take its adjacent vertices in DT(P) as new candidates to be included in W_i . For checking whether a candidate point p lies in W_i we use a data structure to find a nearest neighbour of p in W_{i-1} . If the distance from p to its nearest neighbour w in W_{i-1} is smaller than 1, then the shortest path tree is extended by connecting p to w.

```
UNWEIGHTEDSHORTESTPATH(P, r)
    build the Delaunay triangulation DT(P)
     for p \in P
 3
           dist[p] = \infty
          \pi[p] = \text{NIL}
     dist[r] = 0
     W_0 = \{r\}
 6
 7
     i = 1
     while W_{i-1} \neq \emptyset
 8
 9
          build data structure for nearest neighbour queries in W_{i-1}
10
           Q = W_{i-1}
                            // candidate points
           W_i = \emptyset
11
           while Q \neq \emptyset
12
13
                q an arbitrary point of Q
14
                remove q from Q
15
                for qp edge in DT(P)
                      w = \text{nearest neighbour of } p \text{ in } W_{i-1}
16
17
                      if dist[p] = \infty and |pw| \leq 1
                            dist[p] = i
18
                            \pi[p] = w
19
20
                            add p to Q
21
                            add p to W_i
22
          i = i + 1
     return dist[\cdot] and \pi[\cdot]
23
```

Figure 1 Algorithm from [3] to compute a shortest path tree in the unweighted case.

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Cabello and Jejčič [3] show that the algorithm correctly computes the shortest-path tree 86 from r. If for nearest neighbors we use a data structure that, for n points, has construction 87 time $T_c(n)$ and query time $T_q(n)$, and the Delaunay triangulation is computed in $T_{DT}(n)$ time, 88 then the algorithm takes $O(T_{DT}(n) + T_c(n) + nT_q(n))$ time. Standards tools in Computational Geometry imply that $T_{DT}(n) = O(n \log n)$, $T_c(n) = O(nq \log n)$ and $T_q(n) = O(\log n)$. This leads to the following.

▶ Theorem 2.1 (Cabello and Jejčič [3]). Let P be a set of n points in the plane and let r be a point from P. In time $\mathcal{O}(n \log n)$ we can compute a shortest path tree from s in the unweighted graph $G_{\leq 1}(P)$.

It is clear that, when computing the shortest path tree from several sources, we only need to compute the Delaunay triangulation once.

2.2 Minimum separation with unit-disk

Cabello and Giannopouolos [2] present an algorithm for the minimum separation problem that in the worst-case runs in cubic-time. The algorithm has one feature that is both an advantage and a disadvantage: it works for any reasonable shapes, like segments or ellipses, and not just unit disks. This means that it is very generic, which is good, but it cannot exploit any properties of unit disks.

In this section we are going to describe an algorithm to solve the minimum separation problem for unit disks in roughly quadratic time. The improvement is based on 3 ingredients. The first ingredient is a reinterpretation of the algorithm of [2] for disks. In the original algorithm, we had to select a point inside each shape. For disks there is a natural, obvious choice, the center of the disk. This allows for a simpler description and interpretation of the algorithm. We provide the description in Section 2.2.1

The second ingredient is the efficient algorithm for shortest-path trees for the graph G. The third ingredient is a compact treatment of the edges of G using a few tools from Computational Geometry, namely range trees, point-line duality, and nearest-neighbour searches. This is explained in Section 2.2.2.

2.2.1 Generic algorithm specialized for unit disks

Let us first introduce some notation. Each walk W in the graph $G = G_{\leq 1}(P)$ defines a planar curve in the obvious way: we connect the points of P with segments in the order given by W. We will relax the notation slightly and denote also by W the curve itself. For any spanning tree T of G and any edge $e \in E(G) \setminus E(T)$, let cycle(T,e) be the unique cycle in T + e. Finally, for any walk in G(P), let $\operatorname{cr}_2(st,W)$ be the modulo 2 value of the number of crossings between the segment st and (the curve defined by) W.

The following property is implicit in [2] and explicit in [4]:

Let T be any spanning tree of G. The set of unit disks with centers in P separate s and t if and only if there exists some edge $e \in E(G) \setminus E(T)$ such that $cr_2(st, cycle(T, e)) = 1$.

A consequence of this is that finding a minimum separation amounts to finding a shortest cycle G that crosses the segment st an odd number of times. Moreover, one can show that we can restrict our search to a very concrete family cycles, as follows. For each root r let us fix a shortest-path tree T_r from r in G. When there are many, the choice of T_r is irrelevant. Then, we can restrict our search to

$$\{cycle(T_r, e) \mid r \in P, e \in E(G) \setminus E(T_r)\}.$$

This follows from the co-called 3-path condition; see [2] for the ideas, and appendix ?? for a self-contained proof.

The values $\operatorname{cr}_2(st, cycle(T_r, e))$ can be computed in constant amortized time per edge with some easy bookkeeping. Consider a fixed tree T_r . For each point $p \in P$ we store N[p] as the parity of the number of crossings of the path in T_r from r to p. When p is not the root, the value N[p] can be computed from the value of its parent $\pi[p]$ in T_r using that $N[p] = N[\pi[p]] + \operatorname{cr}_2(st, p\pi[p])$. In the algorithm we have written it this way (lines 4–6), but one can of course compute the values at the time of computing the shortest path tree T_r .

We then have for each T_r

```
\forall pq \in E(G) \setminus E(T_r): \quad \operatorname{cr}_2(st, cycle(T_r, pq)) = N[p] + N[q] + \operatorname{cr}_2(st, pq) \pmod{2}\forall pq \in E(T_r): \quad 0 = N[p] + N[q] + \operatorname{cr}_2(st, pq) \pmod{2}
```

APPENDIX

because crossings that are counted twice cancel out modulo 2. (In particular, the path in T_r from r to the lowest common ancestor of p and q is counted twice.) This implies that we can just check for all edges pq of G whether the sum $N[p] + N[q] + \operatorname{cr}_2(st, pq)$ is 0 modulo 2. The final resulting algorithm, denoted as Generic Minimum Separation, is given in Figure 2.

```
GENERICMINIMUMSEPARATION(P, s, t)
     best = \infty // length of the best separation so far
 2
     for r \in P
           (dist[], \pi[]) = shortest path tree from r in G(P)
 3
           /\!\!/ Compute N[]
           N[r] = 0
 4
 5
           for p \in P \setminus \{r\} in non-decreasing values of dist[p]
 6
                N[p] = N[\pi[p]] + \operatorname{cr}_2(st, p\pi[p]) \pmod{2}
 7
           for pq \in E(G(P))
 8
                if N[p] + N[q] + \operatorname{cr}_2(pq, st) \pmod{2} = 1
 9
                      best = min\{best, dist[p] + dist[q] + 1
10
     return best
```

Figure 2 Adaptation of the generic algorithm to compute the minimum separation for unit disks.

Let us look into the time complexity of the algorithm. For each point $r \in P$ we have to compute a shortest-path tree in G. This can be done in $O(n \log n)$ in our case, as discussed in Section 2.1. Then, for each edge pq of G some constant amount of work is done. Thus for each point r we spend $O(n \log n + |E(G)|)$. This is cubic in the worst-case. We could get an improved running time if we can treat all the edges of G compactly. This is what we explain next.

2.2.2 Compact treatment of edges

Without loss of generality, in our theoretical discussion we will assume that z = (0,0) and z' = (0,s). Therefore, zz' is a segment, henceforth denoted by σ , contained in the y-axis.

3 Implementation

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3.1 Shortest-path tree in unit-disk graphs

Alternative constructions. Let us mention two obvious alternatives that we use in our comparison.

The first alternative is to build the graph $G = G_{\leq 1}(P)$ explicitly. Thus, for each pair of points p,q we check whether their distance is at most once and add an edge to a graph data structure. We can then use breadth-first-search (BFS) from the given root r. The preprocessing is quadratic, and the time spent to compute a shortest-path tree depends on the density of the graph G.

The second option we consider is to use a unit-length grid. Two points (x, y) and (x', y') are in the same grid cell if and only if $(\lfloor x \rfloor, \lfloor y \rfloor) = (\lfloor x' \rfloor, \lfloor y' \rfloor)$. We store all the points of a grid cell c in a list $\ell(c)$. The non-empty lists $\ell(c)$ are stored in a dictionary, where the bottom-left corner of the cell is used as key. We can then run some sort of BFS using this

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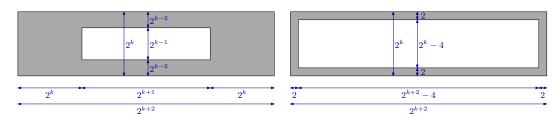


Figure 3 Data generation with a small hole (left) and a large hole (right).

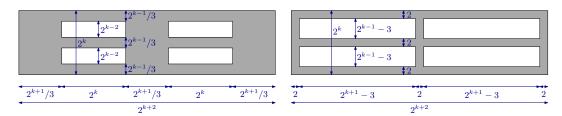


Figure 4 Data generation with four small holes (right) and four large holes (right).

data. When processing a point p in a cell c, we have to treat all the points in c and its adjacent cells as candidate points. The preprocessing is linear, and the time spent to compute a shortest-path tree depends on the density of the graph G. The number of candidates that are checked is proportional to the number of edges of the graph.

Are perhaps 16 points deleted from the list 16 once they are assigned a 16 level?

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3.2 Minimum separation with unit-disk

4 Experimental results

Data was generated uniformly at random in polygons...

4.1 Shortest-path tree in unit-disk graphs

4.2 Minimum separation with unit-disk

5 Conclusions

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A Building block

The following ideas are probably folklore. See [?].

▶ **Lemma 1.1.** Let P be a set of n weighted points in the plane. In $O(n \log n)$ time we can construct a data structure that, for a query point q, it finds in time $O(\log^2 n)$ a point in

```
\arg\min\{w_p \mid p \in P, |pq| \leqslant 1\}.
```

Proof. We sort the points of P by non-decreasing weight, build a balanced binary search tree \mathcal{T} with n leaves, and place the points of P in the leaves of \mathcal{T} such that the order arising from the search tree and the sorting of P match.

For every node ν of \mathcal{T} :

- Let $P(\nu)$ denote the set of points stored at the subtree rooted at ν . We refer to $P(\nu)$ as a canonical subset.
- Let $U(\nu)$ be the union of unit disks centered at the points of $P(\nu)$.
- We build a data structure to decide whether a given query point is contained in $U(\nu)$. A way to do this is to construct a point-location data structure for the Voronoi diagram of $P(\nu)$. Given a query point q, we find a closest neighbor from $P(\nu)$ by locating a cell c of the Voronoi diagram that contains q. We then check whether q is at most at unit distance from the site defining c. If the points of $P(\nu)$ are sorted (lexicographically), then the Voronoi diagram of $P(\nu)$ can be build in linear time ?? The preprocessing time for node ν is $O(|P(\nu)|)$ and a query can be answered in $O(\log |P(\nu)|) = O(\log n)$ time.

This finishes the description of the construction of the data structure. Let us analyze the building time. We sort the points of P lexicographically at the root and for each node ν we get the points $P(\nu)$ sorted from its parent. Since the canonical subsets at each level of the tree are disjoint, at each level of the tree we spend O(n) time. Since \mathcal{T} is balanced, there are $O(\log n)$ levels and we spend in total $O(n \log n)$ time to build the data structure.

Given a point q we first check whether q lies in U(r) for the root r of \mathcal{T} . If q is not in U(r), then no point of P is close enough to q. Otherwise we set $\nu = r$ and proceed with a top-to-down search in the tree \mathcal{T} until we reach a leaf. When ν is not a leaf, let ν_{ℓ} and ν_{r} be the left and right children of ν , respectively. If q is in $U(\nu_{\ell})$, we then proceed with the search in the subtree at ν_{ℓ} by setting $\nu = \nu_{\ell}$, otherwise we proceed the search in the right subtree setting $\nu = \nu_{r}$. At each point of the search, we maintain the following invariant:

$$P(\nu) \cap \arg\min\{w_p \mid p \in P, |pq| \le 1\} \ne \emptyset.$$

The search path in \mathcal{T} has $O(\log n)$ nodes and, for each node in the path, spend $O(\log n)$ time to decide whether q lies in $U(\nu)$. The total query time is thus $O(\log^2 n)$.

Consider the following problem for two weighted point sets A and B.

$$\Phi(A, B) := \min \quad w_a + w_b$$
s.t. $a \in A, b \in B$

$$|ab| \leq 1.$$

Lemma 1.2. Let A and B be sets of at most n weighted points in the plane. In $O(n \log^2 n)$ time we can compute $\Phi(A, B)$.

Proof. We build the data structure of the previous Lemma for B. For each $a \in A$, we query the data structure to obtain

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b^*(a) \in \arg\min\{w_b \mid b \in B, |ab| \le 1\}.
```

Then we find the $a \in A$ minimizing the sum $w_a + w_{b^*(a)}$.

We need $O(n \log n)$ time to construct the data structure and $O(\log^2 n)$ time per query. The result follows.

Let σ be a segment in the plane. Without loss of generality, we will assume that s lies on the y-axis with endpoints (0,0) and (0,s). Let A be a set of points with negative x-coordinates and let B be a set of points with positive x-coordinates. Each point p of $A \cup B$ has a weight w_p . We are interested in minimizing $w_a + w_b$ over all pairs $(a,b) \in A \times B$ whose segment ab have length at most 1 and intersects s. Thus, we want to find

$$\Phi_{\sigma}(A,B) := \min \quad w_a + w_b$$
 s.t. $a \in A, b \in B$
$$|ab| \leqslant 1$$

$$ab \text{ intersects } \sigma.$$

For every point $a \in A$ we define the sets

$$B(a) = \{b \in B \mid ab \text{ intersects } \sigma\},$$

$$B_{\leqslant 1}(a) = \{b \in B \mid ab \text{ intersects } \sigma \text{ and } |ab| \leqslant 1\} = \{b \in B(a) \mid |ab| \leqslant 1\}$$

and the optimization problem

$$\Phi_{\sigma}(a, B) = w_a + \min\{w_b \mid b \in B_{\leq 1}(a)\}.$$

We thus have

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$$\Phi_{\sigma}(A,B) = \min_{a \in A} \Phi_{\sigma}(a,B).$$

We first provide a data structure to obtain the sets $B(\cdot)$ compactly.

Lemma 1.3. There is a family $\{B_1, \ldots, B_t\}$ of subsets of B and a data structure $\mathcal{D}(B)$ with the following properties

```
\sum_{i=1}^{t} |B_i| = O(n \log n);
```

 $\mathcal{D}(B)$ has size $O(n \log n)$ and can be constructed in $O(n \log n)$ time;

for each point a with negative x-coordinate, there is a subset of indices $I(a) \subset \{1, \ldots, t\}$ such that $|I(a)| = O(\log^2 n)$ and B(a) is the disjoint union of $\{B_i\}_{i \in I(a)}$;

for each query point a with $a_x < 0$, the data structure $\mathcal{D}(B)$ returns I(a) in $O(\log^2 n)$ time.

Proof. It is convenient to use point-line duality. We use the precise duality described in BKOS Chapter XXX: the non-vertical line $\ell \equiv y = mx + c$ is mapped to the point $\ell^* = (m, -c)$.

Let $\mathbb L$ be the set of non-vertical lines. Let σ^* be the set of points dual to non-vertical lines that intersect s. Thus

$$\sigma^* = \{l^* \mid \ell \in \mathbb{L}, \ell \cap \sigma \neq \emptyset\}$$

In the dual space, the set σ^* is the horizontal slab

$$\sigma^* = \{ (m, -c) \in \mathbb{R}^2 \mid 0 \leqslant c \leqslant s \}.$$

For every point $b \in B$, let L_b^* be the set of points dual to the lines through b that intersect σ :

$$L_b^* = \{ \ell^* \mid \ell \in \mathbb{L}, b \in \ell, \text{ and } \sigma \cap \ell \neq \emptyset \}.$$

In the dual space, L_b^* is a segment with endpoints $(\varphi_1(b), 0)$ and $(\varphi_2(b), s)$, for some values $\varphi_1(b)$ and $\varphi_2(b)$ that are easily computable. In particular, L_b^* is contained in the slab σ^* and has the endpoints on different boundaries of σ^* . Finally, define the mapping point $\varphi(b) = (\varphi_1(b), \varphi_2(b))$. Thus, φ maps points to the right of the y-axis to points in the plane. Namely, $\varphi_1(b)$ is the slope of the line through b and (0,0) while $\varphi_2(b)$ is the slope of the line through b and (0,s).

Let a be any point to the left of the y-axis and let $b \in B$. The segment ab intersects σ if and only if L_a^* intersects L_b^* . Namely, an intersection of L_a^* and L_b^* is dual to the line through a and b. The segments L_a^* and L_b^* intersect if and only if the order of their endpoints on boundaries of σ^* are reversed. Thus we have the following property:

$$ab \cap \sigma \neq \emptyset \iff (\varphi_1(a) - \varphi_1(b)) \cdot (\varphi_2(a) - \varphi_2(b)) < 0.$$

Given a point a, the set of points $b \in B$ with the property that ab intersects σ corresponds to the points b with $\varphi(b)$ in 2 quadrants with apex $\varphi(a)$.

We can use a 2-dimensional range tree to store the point set $\varphi(B)$, where each point $b \in B$ is identified with its image $\varphi(b)$. For any query $a \in A$, the points $b \in B$ such that ab intersects σ are obtained by querying the 2-dimensional range tree for the points of $\varphi(B)$ contained in the two quadrants

$$\{(x,y) \mid \varphi_1(a) > x \text{ and } \varphi_2(a) < y\} \cup \{(x,y) \mid \varphi_1(a) < x \text{ and } \varphi_2(a) > y\}.$$

The details are standard in computational geometry; see Chapter XXX of BKOS.

Lemma 1.4. We can compute $\Phi_{\sigma}(A, B)$ in $O(n \log^4 n)$ time.

Proof. We construct the sets $\{B_1, \ldots, B_t\}$ and the data structure $\mathcal{D}(B)$ described in Lemma 1.3. For each B_j , where $j = 1, \ldots, t$, we build the data structure of Lemma 1.1. Since $\sum_{i=1}^{t} |B_i| = O(n \log n)$, we need

$$\sum_{i=1}^{t} O(|B_i| \log |B_i|) = O(n \log^2 n)$$

time for this.

Consider a point $a \in A$. We query $\mathcal{D}(B)$ and obtain the set I(a) of indices such that

$$B(a) = \bigcup_{i \in I(a)} B_i.$$

Now, for each $i \in I(a)$, we query with a the data structure associated to B_i to obtain

$$b_i^*(a) \in \arg\min\{w_b \mid b \in B_i, |ab| \leq 1\}.$$

We then have that

$$\min\{w_b \mid b \in B_{\leq 1}(a)\} = \min\{w_{b_i^*(a)} \mid i \in I(a)\},\$$

so we can obtain $\Phi_{\sigma}(a, B)$ from the points $b_i^*(a)$, $i \in I(a)$.

For a point a, we need $O(\log^2 n)$ time to obtain I(a), and then we need $O(\log^2 n)$ time for each index i in I(a). This means that we spend time $O(\log^4 n)$ for each point $a \in A$.

B Crossing a fixed segment

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We will drop in the notation the dependency on P and just write G instead of G(P). We will regard G as an unweighted graph and denote by $d_G(r,p)$ the shortest-path distance in G between r and p.

For each point $r \in P$, let let T_r be a shortest-path tree from r in G. If there are several possible candidates, we fix T_r . We use $T_r[p]$ for the path in T_r from r to p. Let E_r be the set of edges of G not appearing in T_r . Thus $E_r = E(G) \setminus E(T_r)$.

For each r and p from P, let $\gamma(r,p)$ be the polygonal path defined by $T_r[p]$. For each $r \in P$ and each edge $pq \in E_r$, let $\gamma(r,pq)$ be the polygonal closed path obtained be concatenating $\gamma(r,p)$, pq, and the reverse of $\gamma(r,q)$. We define len(r,pq) as the number of edges in $\gamma(r,pq)$. Thus we have $len(r,pq) = d_G(r,p) + d_G(r,q) + 1$.

Let σ be a segment. For simplicity we assume that σ does not contain any point of P. For each polygonal path γ , let $X(\gamma, \sigma)$ be the modulo two value of the number of crossings between γ and σ . For each $r, p \in P$ we define $N(r, p) = X(\gamma(r, p), \sigma)$, and for each $r \in P$ and $pq \in E_r$ we define $N(r, pq) = X(\gamma(r, pq), \sigma)$.

$$\Psi_{\sigma}(P) := \min \quad len(r,pq)$$
 s.t. $r \in P$
$$pq \in E_r$$

$$N(r,pq) = 1.$$

The results of Cabello and Giannopoulos [?] imply the following result. (Should we provide a self-contained proof adapted to this scenario?)

Theorem 2.1. Let (r^*, p^*q^*) be an optimal solution to $\Psi_{\sigma}(P)$. Then the points in $T_{r^*}[p] \cup T_{r^*}[q]$ are an optimal solution to the separation problem.

To compute $\Psi_{\sigma}(P)$ we will iterate over the points r, so we define

$$\Psi_{\sigma}(r, P) := \min \ d_G(r, p) + d_G(r, q)$$

s.t. $pq \in E_r$
 $N(r, pq) = 1.$

We readily have

$$\Psi_{\sigma}(P) = 1 + \min_{r \in P} \Psi_{\sigma}(r, P)$$

Let us fix some $r \in P$ and discuss how to solve $\Psi_{\sigma}(r, P)$. We compute a shortest path tree T_r from r. Together with each point $p \in P$ we store $d_G(r, p)$.

We compute the values N(r,p), $p \in P$, with a simple bottom-up traversal of T_r . We start setting $N_r[r] := 0$. For each point $p \in P$ whose parent in T_r is p', we obtain $N_r[p]$ using that $N_r[p] = N_r[p'] + |\sigma \cap pp'| \pmod{2}$. This is, if pp' crosses σ , then $N_r[p] = N_r[p'] + 1 \pmod{2}$, otherwise $N_r[p] = N_r[p']$.

Assume that σ is a vertical segment with endpoints (0,0) and (0,s).

We split the search for the optimal edge pq into cases:

pq has both endpoints on the same side of the y-axis and connects points with $N(r,p) \neq N(r,p)$, or

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pq has endpoints in both sides of the y-axis, $N(r,p) \neq N(r,p)$, and pq does not intersect σ , or

pq has endpoints in both sides of the y-axis, $N(r,p) \neq N(r,p)$, and pq intersects σ .

▶ **Lemma 2.2.** We can compute $\Psi_{\sigma}(r, P)$ in $O(n \log^4 n)$ time.

Proof. For $i \in \{0,1\}$, let L_i be the subset of points of P to the left of the y-axis with $N(r,p) = i \pmod{2}$. For $i \in \{0,1\}$, let R_i be the subset of points of P to the right of the y-axis and $N(r,p) = i \pmod{2}$. Clearly

$$P = L_0 \cup L_1 \cup R_0 \cup R_1.$$

Let σ_+ and σ_- be the rays contained in the y-axis after deleting σ . Using $d_G(r, p)$ as the weight w_p of point $p \in P$, we then have

$$\begin{split} \Psi_{\sigma}(r,P) \ = \ & \min\{\Phi(L_0,L_1), \Phi(R_0,R_1), \Phi_{\sigma_+}(L_0,R_1), \Phi_{\sigma_+}(L_1,R_0), \\ & \Phi_{\sigma_-}(L_0,R_1), \Phi_{\sigma_-}(L_1,R_0), \Phi_{\sigma}(L_0,R_0), \Phi_{\sigma}(L_1,R_1)\}. \end{split}$$

Each of these instances can be solved using Lemmas 1.2 and 1.4 in $O(n \log^4 n)$ time.

▶ **Theorem 2.3.** We can solve the minimum separation problem using $O(n^2 \log^4 n)$ time.

Proof. For each $r \in P$ we compute the shortest path tree T_r from r in T, the distances $d_G(r,p), p \in P$, and the parities N(r,p). We can then compute $\Psi_{\sigma}(r,P)$ using Lemma 2.2. Repeat for each $r \in P$. Time bound.

Correctness.

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