

Constructing the relative neighborhood graph in 3-dimensional Euclidean space

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Abstract

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The relative neighborhood graph for a finite set $S = \{p_1, \dots, p_N\}$ of points, briefly $\text{RNG}(S)$, is defined by the following formation rule: $\overline{p_i p_j}$ is an edge in $\text{RNG}(S)$ if and only if for all $p_k \in S - \{p_i, p_j\}$, $\text{dist}(p_i, p_j) \leq \max(\text{dist}(p_i, p_k), \text{dist}(p_j, p_k))$.

We show that RNG for point sets in \mathbb{R}^3 can be constructed in optimal space and $O(N^2 \log N)$ time. Also, combinatorial estimates on the size of RNG in \mathbb{R}^3 are given.

1. Introduction

Let $S = \{p_1, \dots, p_N\}$ be a set of points in \mathbb{R}^d . The relative neighborhood graph of S , briefly $\text{RNG}(S)$, is a graph with the set of vertices S and the set of edges defined by the following formation rule:

$\overline{p_i p_j}$ is an edge in $\text{RNG}(S)$ if and only if for all $p_k \in S - \{p_i, p_j\}$,
 $\text{dist}(p_i, p_j) \leq \max(\text{dist}(p_i, p_k), \text{dist}(p_j, p_k))$

where dist denotes the standard Euclidean metric. Further in the paper we will simply use $|\overline{p_i p_j}|$ for $\text{dist}(p_i, p_j)$.

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Relative neighborhood graphs have been defined by Toussaint [18] as an interesting variant of graphs introduced by Lankford [14]. The immediate interest in these graphs was induced by both their combinatorial flavor and numerous potential applications, e.g. in pattern recognition (see [19]). A number of results on relative neighborhood graphs, their variants, and generalizations have been published; see [7, 9–11, 13, 15–18].

The 2-dimensional case of relative neighborhood graphs has been particularly well studied. The first optimal $O(N \log N)$ algorithm was obtained by Supowit [17] who utilized a clever sweep line approach with sweeping conducted in 6 directions. Lee [15] and O'Rourke [16] addressed 2-dimensional RNG with the metrics L_1 and L_∞ . The problem of constructing RNG from the Delaunay triangulation has been investigated by Jaromczyk, Kowaluk and Yao [10], where an optimal $O(N)$ construction has been shown. In fact, they also showed that this result extends to the metric L_p , $1 < p < \infty$, and is valid for a wider class of graphs.

On the other hand, the problem of constructing relative neighborhood graphs in higher-dimensional spaces is still unexplored. For a long time no better result than an $O(n^3)$ algorithm presented by Toussaint [18] was known. Apparently the first subcubic algorithm for constructing relative neighborhood graphs in \mathbb{R}^d was given by Supowit [17] for a special class of input points. He designed an $O(N^2)$ algorithm for sets of input points where no three points form an isosceles triangle (or equivalently, all pairwise distances between points in the input set are distinct). Later this result was generalized to the relative neighborhood graphs with the metric L_p , $1 < p < \infty$; still, the absence of isosceles triangles was crucial (see Jaromczyk and Kowaluk [9]).

This paper gives an $O(N^2 \log N)$ time and optimal space algorithm for arbitrary point sets in \mathbb{R}^3 . More specifically we prove the following:

Theorem 1.1. *Given $S = \{p_1, \dots, p_N\}$, $p_i \in \mathbb{R}^3$, $i = 1, \dots, N$, the relative neighborhood graph of S can be constructed in $O(N^2 \log N)$ time and optimal $O(\mu_3(S))$ space; $\mu_3(S)$ denotes the size of RNG(S).*

We will also discuss in this paper an estimate for $\mu_3(N) = \max(\mu_3(S))$: S is an N point set in \mathbb{R}^3 .

This paper is divided into five sections. The next section gives geometric preliminaries that lay the groundwork for the main algorithm. Section 3 will present our algorithm, Section 4 gives an upper bound on the size of the relative neighborhood graphs in \mathbb{R}^3 , and finally Section 5 will provide concluding remarks.

2. Geometry of RNG in \mathbb{R}^3

The definition of relative neighborhood graphs directly implies a rule that can be used to identify edges that do not belong to RNG. This rule tells that an edge $\overline{p_i p_j}$

does not belong to RNG if and only if there is a point p_k which is closer to both of p_i , p_j than the distance between p_i and p_j . For the purpose of our algorithm, however, it is more convenient to have an elimination rule that is based on relations between the lengths of edges and angles between them rather than on relations between distances.

Such a rule is implied by the following lemma:

Lemma 2.1. $\overline{p_i p_k} \notin \text{RNG}(S)$ if and only if there exists p_j such that $|\overline{p_i p_j}| < |\overline{p_i p_k}|$ and

$$\angle(\overline{p_i p_j}, \overline{p_i p_k}) < \arccos \frac{|\overline{p_i p_j}|}{2|\overline{p_i p_k}|}.$$

Proof. We will show that under the assumptions $\overline{p_i p_k}$ is the longest edge in the triangle $\triangle p_i p_j p_k$; see Fig. 1. Let $\alpha = \angle(\overline{p_i p_j}, \overline{p_i p_k})$. By the law of cosines we have

$$\begin{aligned} |\overline{p_j p_k}|^2 &= |\overline{p_i p_k}|^2 + |\overline{p_i p_j}|^2 - 2|\overline{p_i p_j}| \times |\overline{p_i p_k}| \times \cos \alpha \\ &< |\overline{p_i p_k}|^2 + |\overline{p_i p_j}|^2 - 2 \times \frac{|\overline{p_i p_j}|}{2|\overline{p_i p_k}|} \times |\overline{p_i p_j}| \times |\overline{p_i p_k}| \\ &= |\overline{p_i p_k}|^2. \end{aligned}$$

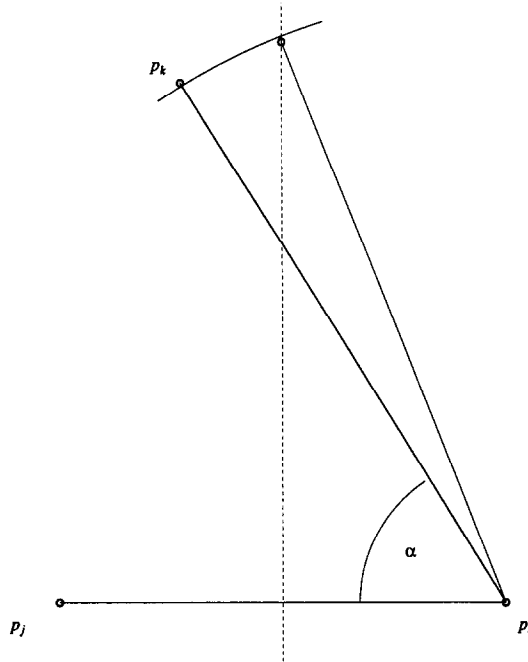


Fig. 1.

Hence $\overline{p_i p_k}$ is the longest edge in $\triangle p_i p_j p_k$ and therefore does not belong to $\text{RNG}(S)$. On the other hand, if $\overline{p_i p_k} \notin \text{RNG}(S)$, then there exists p_j such that $\overline{p_i p_k}$ is the (strictly) longest edge in $\triangle p_i p_j p_k$. By elementary geometry $\cos \alpha > |\overline{p_i p_j}|/(2|\overline{p_i p_k}|)$ which ends the proof. \square

Observe that in particular if the angle between $\overline{p_i p_j}$ and $\overline{p_i p_k}$ ($|\overline{p_i p_k}| > |\overline{p_i p_j}|$) is less than $\frac{1}{3}\pi$, then $\overline{p_i p_k}$ is the longest edge in $\triangle p_i p_j p_k$ and therefore cannot belong to $\text{RNG}(S)$. This simple fact which will be very useful later is formulated as the following lemma:

Lemma 2.2. *If $|\overline{p_i p_j}| < |\overline{p_i p_k}|$ and $\angle(\overline{p_i p_j}, \overline{p_i p_k}) < \frac{1}{3}\pi$, then $\overline{p_i p_k} \notin \text{RNG}(S)$.*

Remark that Lemma 2.1 has the following geometric interpretation: take two points p_i, p_j and the bisecting plane H perpendicular to $\overline{p_i p_j}$. If $|\overline{p_i p_k}| > |\overline{p_i p_j}|$ and $\angle(\overline{p_i p_j}, \overline{p_i p_k}) < \arccos(|\overline{p_i p_j}|/(2|\overline{p_i p_k}|))$, then $\overline{p_i p_k}$ intersects H . Lemma 2.1 states that no edge $\overline{p_i p_k}$ which is longer than $\overline{p_i p_j}$ and pierces H can belong to the relative neighborhood graph of S .

3. Algorithm to construct RNG in \mathbb{R}^3

The algorithm presented in the upcoming sections will consist of two elimination phases responsible for identifying those pairs of points that do not form edges in $\text{RNG}(S)$. The first phase, called a *coarse elimination*, will be based on the elimination rule implied by Lemma 2.2 from the previous section. The second phase, called a *fine elimination*, will utilize Lemma 2.1 to eliminate those edges not in $\text{RNG}(S)$ that remain after the coarse elimination. For the sake of clarify these two elimination phases, which can be interleaved, will be presented in sequence.

3.1. Coarse elimination

The objective of this phase is to construct a “small” supergraph of $\text{RNG}(S)$. This supergraph will have the property that the number of different lengths of edges adjacent to each point is bounded by a constant independent of N . This will be essential for a fast “cleaning-up” process which will be done in the fine elimination. Existence of such a supergraph is guaranteed by Lemma 2.2 which states that the longer of two edges (with a common endpoint) that form an angle not greater than $\frac{1}{3}\pi$ cannot belong to $\text{RNG}(S)$.

The coarse elimination will be carried out separately for each point $p_i \in S$. We start with the set $\text{adj}(p_i) = \{\overline{p_i p_j} : j = 1, \dots, N, i \neq j\}$ of all edges having p_i as an endpoint. This set will be subsequently modified by removing some of the edges that do not belong to $\text{RNG}(S)$. At each stage all of the (currently) shortest edges in this set are placed in a so-called *orbit*, and then they are removed from the $\text{adj}(p_i)$ set.

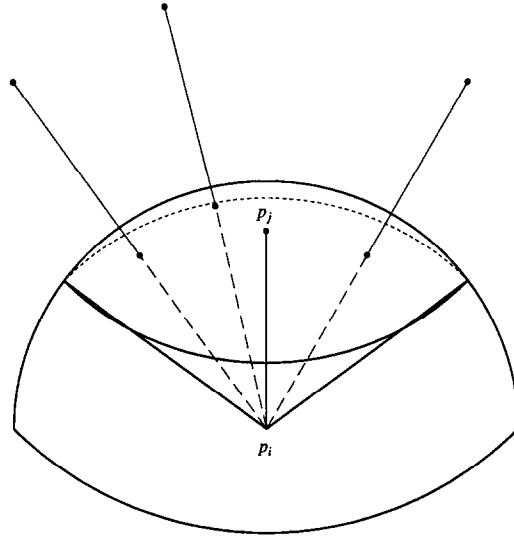


Fig. 2. $\overline{p_i p_j}$ is a central edge.

A representative edge from the orbit is designated as a *central* edge for this orbit and will serve as a basis for further eliminations.

Note that at this moment all edges in $\text{adj}(p_i)$ are longer than the central edge. Then those edges that form an angle not greater than $\frac{1}{3}\pi$ with central are removed from the set $\text{adj}(p_i)$; they do not belong to $\text{RNG}(S)$ by virtue of Lemma 2.2; see Fig. 2.

This process, which forms subsequent orbits and eliminates edges that form with the current central edge an angle less than $\frac{1}{3}\pi$, is repeated until $\text{adj}(p_i)$ is empty. Edges eliminated by this process will be placed in the set $\text{Coarse_Elim}(p_i)$. Let us note that in fact it is not necessary to store $\text{Coarse_Elim}(p_i)$ explicitly; this set is merely used to facilitate a proof of the algorithm correctness. The pseudo-code for the procedure $\text{Coarse_Elimination}(p_i)$ is given below.

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procedure Coarse_Elimination( $p_i$ );
   $\text{adj}(p_i) := \{ \overline{p_i p_j} : p_j \in S - \{p_i\} \}$ 
   $\text{number\_of\_orbits} := 0$ ;
   $\text{Coarse\_Elim}(p_i) := \emptyset$ ;
  while  $\text{adj}(p_i) \neq \emptyset$  do
    begin
       $\text{number\_of\_orbits} := \text{number\_of\_orbits} + 1$ ;
       $\overline{p_i p_l} := (\text{one of the}) \text{ shortest edge in } \text{adj}(p_i)$ ;
       $\text{central}(p_i, \text{number\_of\_orbits}) := \overline{p_i p_l}$ ;
       $\text{orbit}(p_i, \text{number\_of\_orbits}) := \{ \overline{p_i p_j} \in \text{adj}(p_i) : |\overline{p_i p_j}| = |\overline{p_i p_l}| \}$ ;
       $\text{adj}(p_i) := \text{adj}(p_i) - \text{orbit}(p_i, \text{number\_of\_orbits})$ ;
    end

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Coarse_Elim( $p_i$ ) := Coarse_Elim( $p_i$ )
                      $\cup \{ \overline{p_i p_j} \in \text{adj}(p_i) : \angle(\overline{p_i p_j}, \overline{p_i p_l}) < \frac{1}{3}\pi \}$ ;
adj( $p_i$ ) := adj( $p_i$ ) - Coarse_Elim( $p_i$ );
end;
number_of_orbits( $p_i$ ) := number_of_orbits;
end procedure;

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The major observation is that the number of the orbits for each point is bounded by a constant independent on N . Indeed, the following lemma holds.

Lemma 3.1. *For all $p_i \in S$, $\text{number_of_orbits}(p_i) \leq c_3$, where c_3 is a constant independent on N .*

Proof. For every $\text{central}(p_i, j)$, $j = 1, \dots, \text{number_of_orbits}(p_i)$, let cone_j denote an open circular cone with the apex p_i , symmetric with respect to the line passing through $\text{central}(p_i, j)$ and such that the angle at the apex (i.e., the angle in 2-dimensional plane passing through $\text{central}(p_i, j)$) is equal to $\frac{1}{3}\pi$. Let B be a unit ball centered at p_i . Note that $\angle(\text{central}(p_i, l), \text{central}(p_i, j)) \geq \frac{1}{3}\pi$, for $l \neq j$. Hence $\text{cone}_l \cap \text{cone}_j = \emptyset$ for $l \neq j$. Consequently, the volume of B is not smaller than the sum of the volumes of $\text{cone}_j \cap B$, $j = 1, \dots, \text{number_of_orbits}(p_i)$. Since the volume of each $\text{cone}_j \cap B$ is fixed the number of cones is bounded by a certain constant c_3 . Therefore, the number of central edges, and hence the number of orbits, is bounded by the constant c_3 which is independent on N . \square

Lemma 3.2. *For each $p_i \in S$, $\text{Coarse_Elim}(p_i) \cap \text{RNG}(S) = \emptyset$.*

Proof. This follows from Lemma 2.2 and the way $\text{Coarse_Elim}(p_i)$ is constructed. \square

Note that at this point we cannot claim that the set $\text{Coarse_Elim}(p_i)$ contains all the eliminated edges adjacent to p_i . In fact, the set $\text{orbit}(p_i, l)$ may contain some edges that do not belong to $\text{RNG}(S)$. However, if all the pairwise distances between points in S were distinct, then the orbits would exactly form $\text{RNG}(S)$.

By Lemma 3.1 the number of iterations of the while-loop in the $\text{Coarse_Elimination}$ procedure is bounded by a constant c_3 ; the cost of each iteration is $O(N)$. Hence, we have the following

Lemma 3.3. *The procedure $\text{Coarse_Elimination}$ constructs the family $\text{orbit}(p_i, l), \dots, \text{orbit}(p_i, \text{number_of_orbits})$ in $O(N)$ time.*

Because the coarse elimination is carried out for each point in S the overall cost of this phase is $O(N^2)$.

Clearly the above presented method is not the only way to construct a supergraph of $\text{RNG}(S)$ such that for each point edges adjacent to this point can be grouped into a constant number of orbits consisting of edges of the same length. For example, we can find such a supergraph starting from a general technique developed by Yao [20]. To this end we can think that each point in S is associated with a finite set of vectors (a frame) such that each pair of vectors form an angle of at most $\frac{1}{3}\pi$. Such a frame partitions the space around each point into a finite number of cones. Using Yao's solution to the general geographic neighbor problem, see [20], we can find for all points their nearest neighbors in each of the corresponding cones. The method guarantees that this can be done in a subquadratic time. The edges joining points with such neighbors play a role of our central edges. Clearly, in each cone no edge longer than the corresponding central edge can belong to $\text{RNG}(S)$ (because of the angles between vectors in the frame). This straightforward observation can be used to eliminate further edges in the way similar to the coarse elimination. The overall process takes $O(N^2)$ time and produces a supergraph with the desired properties. This graph is, in general, different than one produced by the Coarse_Elimination procedure.

3.2. Fine elimination

The family of orbits constructed by the Coarse_Elimination procedure may contain edges that do not belong to $\text{RNG}(S)$. The objective of the fine elimination is to discard those edges. For each edge $\overline{p_i p_j}$ in orbit and every shorter edge (than $\overline{p_i p_j}$) the angle between them will be measured and then compared to the threshold angle given in Lemma 2.1. To achieve an efficiency the process of comparing angles will be carried out simultaneously for all edges in a given orbit using an algorithm for the problem of point inclusion in the union of circles.

Before giving a more detailed description of the algorithm let us introduce some notation which will facilitate further presentation. Let V , and W be two sets of edges such that all edges in V are shorter than edges in W . We define $V \blacklozenge W = \{ \overline{p_i p_k} \in W : \angle(\overline{p_i p_j}, \overline{p_i p_k}) < \arccos(|\overline{p_i p_j}| / (2|\overline{p_i p_k}|)), \overline{p_i p_j} \in V \}$. Intuitively, $V \blacklozenge W$ consists of edges in W that can be ruled out from $\text{RNG}(S)$ using the elimination rule implied by Lemma 2.1. We will apply this operation to sets W (second operand) which are orbits, i.e., they consist of edges of the same length. With this restriction in mind, we will show later how this operation may be efficiently implemented. Pseudo-code for the procedure Fine_Elimination(p_i) is presented below.

```

procedure Fine_Elimination( $p_i$ );
  begin
    Elim( $p_i$ ) := Coarse_Elim( $p_i$ );
    adj( $p_i$ ) :=  $\{ \overline{p_i p_j} : p_j \in S \}$ ;
    for  $l := 2$  to number_of_orbits( $p_i$ ) do
      begin

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 $V := \{ \overline{p_i p_j} \in \text{adj}(p_i) : |\overline{p_i p_j}| < \text{length of central}(p_i, l) \};$ 
Fine_Elim  $:= V \blacklozenge \text{orbit}(p_i, l);$ 
 $\text{orbit}(p_i, l) := \text{orbit}(p_i, l) - \text{Fine\_Elim};$ 
 $\text{Elim}(p_i) := \text{Elim}(p_i) \cup \text{Fine\_Elim}$ 
end
end procedure;

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Observe that this procedure applies the elimination rule of Lemma 2.1 to every edge in $\text{orbit}(p_i, l)$. According to the definition of the operation \blacklozenge only edges that are not in $\text{RNG}(S)$ are removed from $\text{orbit}(p_i, l)$. Therefore, for all $p_i \in S$, $\text{Elim}(p_i) \cap \text{RNG}(S) = \emptyset$. In addition, since each edge in $\text{orbit}(p_i, l)$ is tested against all shortest edges, if $\overline{p_i p_j} \notin \text{RNG}(S)$, then $\overline{p_i p_j} \in \text{Elim}(p_i)$.

Now, both of the above procedures can be combined into a final algorithm.

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Main_Algorithm
begin
  for all  $p_i \in S$  do
    begin
      Coarse_Elimination( $p_i$ );
      {removes some edges and partitions the others into orbits}
      Fine_Elimination( $p_i$ );
      {removes from orbits all edges not in  $\text{RNG}(S)$ }
    end;
     $G := \bigcup_{i=1}^N \bigcup_{l=1}^{\text{number\_of\_orbits}(p_i)} \text{orbit}(p_i, l)$ 
  end;

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Lemma 3.4. $G = \text{RNG}(S)$.

Proof. After the coarse and fine eliminations, for each $p_i \in S$,

$$\bigcup_{l=1}^{\text{number_of_orbits}(p_i)} \text{orbit}(p_i, l)$$

consists of edges in $\text{adj}(p_i) - \text{Elim}(p_i)$. In addition $\text{Elim}(p_i)$ contains all edges $\overline{p_i p_j}$ not belonging to $\text{RNG}(S)$. \square

The above lemma implies that the Main_Algorithm is correct, i.e., it constructs $\text{RNG}(S)$. The algorithm is organized in such a fashion that each iteration identifies (using $O(N)$ extra memory) the edges in $\text{RNG}(S)$ adjacent to p_i . Therefore the whole process needs no more than $O(N + \mu_3(S))$ memory. A subquadratic upper bound for $\mu_3(S)$ will be given in the next section; $\mu_3(S)$ is the size of $\text{RNG}(S)$.

The time complexity of the algorithm depends on the implementation of the \blacklozenge operation. Note that this operation is performed on sets of edges V and W where

the second operand contains edges of the same length d . Associate with each edge $\overline{p_i p_j} \in V$ a circular cone with its vertex at p_i , symmetric with respect to $\overline{p_i p_j}$, and with the vertex angle equal to $\arccos(|\overline{p_i p_j}|/(2 \times d))$. By Lemma 2.1, $\overline{p_i p_j}$ eliminates all edges in W that are in the interior of this cone. Consider a sphere with the origin at p_i and the radius d . Clearly, $\overline{p_i p_j}$ eliminates these edges in W which have their endpoints on the spherical cup formed by the intersection of the sphere with the cone associated with $\overline{p_i p_j}$. This observation reduces the problem at hand to a point location problem in the union of the spherical cups corresponding to cones associated with the vectors in V . A suitable stereographic projection transforms this problem to \mathbb{R}^2 where the problem becomes a point location problem for $O(N)$ points in the union of $O(N)$ circles, i.e., given N circles in the plane, determine if a given query per point p is contained in their union. This type of problem has been studied by several authors. In particular, Imai, Iri and Murota [8] have shown that utilizing the Voronoi diagrams in the Laguerre geometry the problem can be solved in $O(\log N)$ time per query point using $O(N \log N)$ preprocessing time. Aurenhammer [1] has utilized power diagrams to the same effect. A solution to the union of circles problem has also been given by Edelsbrunner and Seidel [6] who employed properties of arrangements of hyperplanes. See also Kedem et al. [12]. From these results we can conclude that the \diamond operation can be implemented as an algorithm that uses $O(N \log N)$ steps. This leads to the conclusion given in Theorem 1.1.

4. Upper bound on the size of RNG in \mathbb{R}^3

Let $\mu_d(S)$ denote the number of edges in $\text{RNG}(S)$ where S is a finite point set in \mathbb{R}^d . Define $\mu_d(N) = \max\{\mu_d(S) : |S| = N\}$.

From the well-known inclusion $\text{MST}(S) \subseteq \text{RNG}(S)$ where $\text{MST}(S)$ is a minimal spanning tree on S it follows that $\mu_d(N) \geq N - 1$ (see [18]). In the two-dimensional case $\mu_2(N) \leq 3N - 6$ since $\text{RNG}(S)$ is a subgraph of the Delaunay triangulation of S [18]. Previously no nontrivial upper bound on $\mu_d(N)$ has been known for $d \geq 3$. We will show that $\mu_3(N) < cN^{3/2+\varepsilon}$ where c is a constant and $\varepsilon > 0$. This fact will follow from recent results regarding an upper bound $F_3(N)$ on the number of times a particular distance, say the unit distance, occurs among N points in 3-dimensional Euclidean space. The problem of unit distance graphs was raised by Erdős and there is a large number of papers discussing the size of such graphs. In particular, recent results by Clarkson, Edelsbrunner, Guibas, Sharir and Welzl [5] show that $F_3(N) \leq cN^{3/2+\varepsilon}$, where $\varepsilon > 0$ is an arbitrary small real number. In fact they give this bound in a stronger form of $O(N^{3/2}(\lambda_6(N)/N)^{1/4})$, where λ_6 is related to the complexity of Davenport-Schinzel sequences. For more discussion on $F_3(N)$ and $F_d(N)$ in Euclidean spaces see also Chung [4].

Theorem 4.1. $\mu_3(N) \leq cN^{3/2+\varepsilon}$.

Proof. Consider a set $S = \{p_1, \dots, p_N\}$ of points in \mathbb{R}^3 . Let E_1, \dots, E_k be a partition of the set E of the edges in $\text{RNG}(S)$ such that for all $v, w \in E_i$, $|v| = |w|$. Let $S_i \subseteq S$ be a set of endpoints of the edges in E_i . The number of different lengths of the edges in $\text{RNG}(S)$ adjacent to p_i is bounded by a constant c_3 ; see the proof of Lemma 3.1. Therefore each point in S can appear in at most c_3 sets S_i and $|S_1| + \dots + |S_k| \leq c_3 N$. Since each of E_i consists of edges of the same length, by virtue of the results on $F_3(N)$, we have $|E_i| \leq c |S_i|^{3/2+\varepsilon}$. By induction, using the inequality $a^\alpha + x^\alpha \leq (a+x)^\alpha$ for $a, x > 0$, $\alpha \geq 1$ we have

$$\begin{aligned} |E_1 \cup \dots \cup E_k| &\leq c |S_1|^{3/2+\varepsilon} + \dots + c |S_k|^{3/2+\varepsilon} \leq c(|S_1| + \dots + |S_k|)^{3/2+\varepsilon} \\ &= \text{const} \times N^{3/2+\varepsilon}. \quad \square \end{aligned}$$

Note that the bound of Theorem 4.1 applies also to the supergraph of RNG constructed by the coarse elimination; see Section 3.1.

We believe that the size of the relative neighborhood graphs in \mathbb{R}^3 is smaller than the bound given in Theorem 4.1. On the other hand, a standard example shows that $\mu_4(N) = c \times N^2$. To this end, take $\frac{1}{2}N$ points of the form $(a, b, 0, 0)$ where $a^2 + b^2 = \frac{1}{2}$ and the other $\frac{1}{2}N$ points to be of the form $(0, 0, c, d)$ where $c^2 + d^2 = \frac{1}{2}$. The relative neighborhood graph of this set of points has $c \times N^2$ edges.

5. Concluding remarks

A deterministic algorithm constructing relative neighborhood graphs in \mathbb{R}^3 has been given. This algorithm is optimal with respect to the space complexity and its time complexity is $O(N^2 \log N)$. The algorithm uses two elimination phases which are based on a property of relative neighborhood graphs which relates lengths of edges with angles between them. This property links the problem to finding whether a given point is contained in the union of circles.

Also an interesting combinatorial problem related to the upper bound for the number $\mu_3(N)$ of edges in $\text{RNG}(S)$ has been discussed. Specifically, we have proven that $\mu_3(N) = O(N^{3/2+\varepsilon})$ and this bound has been derived from the combinatorial complexity of unit distance graphs. It would be interesting to find better bounds.

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