The critical radius in sampling-based motion planning

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Abstract
We develop a new analysis of sampling-based motion planning in Euclidean space with uniform random sampling, which significantly improves upon the celebrated result of Karaman and Frazzoli and subsequent work. In particular, we prove the existence of a critical connection radius proportional to \(\Theta(n^{-1/d})\) for \(n\) samples and \(d\) dimensions: below this value the planner is guaranteed to fail (similarly shown by Karaman and Frazzoli). More importantly, for larger radius values the planner is asymptotically (near-)optimal. Furthermore, our analysis yields an explicit lower bound of \(1 - \Theta(n^{-1})\) on the probability of success. A practical implication of our work is that asymptotic (near-)optimality is achieved when each sample is connected to only \(\Theta(1)\) neighbors. This is in stark contrast to previous work that requires \(\Theta(\log n)\) connections, which are induced by a radius of order \(\left(\frac{\log n}{n}\right)^{1/d}\). Our analysis applies to the probabilistic roadmap method (PRM), as well as a variety of “PRM-based” planners, including RRG, FMT*, and BTT. Continuum percolation plays an important role in our proofs. Lastly, we develop similar theory for all the aforementioned planners when constructed with deterministic samples, which are then sparsified in a randomized fashion. We believe that this new model, and its analysis, is interesting in its own right.

Keywords
Motion planning, sampling-based algorithms, probabilistic roadmaps, random geometric graphs, continuum percolation, probabilistic completeness, asymptotic optimality

1. Introduction
Motion planning is a fundamental problem in robotics concerned with allowing autonomous robots to efficiently navigate in environments cluttered with obstacles. Although motion planning has originated as a strictly theoretical problem in computer science (Halperin et al., 2016b), nowadays it is applied in various fields. Notably, motion planning arises in coordination of multiple autonomous vehicles (Frazzoli and Pavone, 2013), steering surgical needles (Baykal and Alterovitz, 2017), and planning trajectories of spacecrafts in orbit (Starek et al., 2016), to name just a few examples.

Motion planning is notoriously challenging from a computational perspective due to the continuous and high-dimensional search space it induces, which accounts for the structure of the robot, the physical constraints that it needs to satisfy, and the environment in which it operates. Low-dimensional instances of the problem can typically be solved by efficient and complete algorithms (Halperin et al., 2016a), and the same applies to several high-dimensional instances (see, e.g., Adler et al., 2015; Solovey et al., 2015). However, in general, motion planning is computationally intractable (see, e.g., Canny, 1988; Reif, 1979; Solovey and Halperin, 2016a).

Nowadays the majority of practical approaches for motion planning capture the connectivity of the free space by sampling (typically in a randomized fashion) configurations and connecting nearby samples, to form a graph data structure (see the detailed discussion in Section 2). Although such sampling-based planners are inherently incomplete, i.e., cannot detect situations in which a solution (collision-free path) does not exist, most have the desired property of being able to find a solution eventually, if one exists. That is, a planner is probabilistically complete (PC) if the probability of finding a solution tends to 1 as the number of samples \(n\) tends to infinity. Moreover, some

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recent sampling-based techniques are also guaranteed to return high-quality solutions that tend to the optimum as $n$ diverges: a property called asymptotic optimality (AO). Quality can be measured in terms of energy, length of the plan, clearance from obstacles, etc.

An important attribute of sampling-based planners, which dictates both the running time and the quality of the returned solution, is the number of neighbors considered for connection for each added sample. In many techniques this number is directly affected by a connection radius $r_n$: decreasing $r_n$ reduces the number of neighbors. This, in turn, reduces the running time of the planner for a given number of samples $n$, but may also reduce the quality of the solution or its availability altogether. Thus, it is desirable to come up with a radius $r_n$ that is small, but not to the extent that the planner loses its favorable properties of PC and AO.

### 1.1. Contribution

We develop a new analysis of the probabilistic roadmap method (PRM) (Kavraki et al., 1996) for uniform random sampling in Euclidean space, which relies on a novel connection between sampling-based planners and continuum percolation (see, e.g., Franceschetti and Meester, 2008). Our analysis is tight and proves the existence of a critical connection radius $r_n^{\gamma} = \gamma' n^{-1/d}$, where $\gamma' > 0$ is a constant (in particular, $0.4 \leq \gamma' \leq 0.6$ for all $d \geq 2$), and $d \geq 2$ is the dimension: if $r_n < r_n^{\gamma}$, then PRM is guaranteed to fail. Above the threshold, i.e., when $r_n > r_n^{\gamma}$, PRM is AO for the bottleneck cost, and asymptotically near optimal (AnO) with respect to the path-length cost. Furthermore, our analysis yields concrete bounds on the probability of success, which is lower-bounded by $1 - O(n^{-1})$. Note that this bound is comparable with that obtained in Starek et al. (2015) (see Section 2) although we show this for a much smaller radius.

Our analysis is not restricted to PRM and applies to a variety of planners that maintain PRM-like roadmaps, explicitly or implicitly. For instance, when $r_n$ is above the threshold, FMT$^*$ (Janson et al., 2015) is AnO with respect to the path-length cost, while BTT (Solovey and Halperin, 2017) is AO with respect to the bottleneck cost. RRG (Karaman and Frazzoli, 2011) behaves similarly for the two cost functions. Our results are also applicable to multi-robot motion planners such as the recently introduced dRRT$^*$ (Shome et al., 2019), and M$^*$ (Wagner and Choset, 2015) when applied to a continuous domain. See Figure 1 for additional PRM-based planners to which our analysis is applicable, and which are mentioned further on.

A practical implication of our work is that AO (or AnO), under the regime of uniform random sampling, can be achieved even when every sample is connected to $\Theta(1)$ neighbors. This is in stark contrast with previous work (e.g., Janson et al., 2015; Karaman and Frazzoli, 2011; Solovey and Halperin, 2016b; Kiril, 2018), which provided a rough estimate of this number, that is proportional to $O(\log n)$. Interestingly, our $\Theta(1)$ bound for uniform samples is comparable to the best known bound of for deterministic samples (Janson et al., 2017).

Lastly, we also study the asymptotic properties of the aforementioned planners when constructed using a deterministic point process, rather than uniform random sampling which we have mentioned so far. We introduce an analysis that shows that the graphs constructed using such deterministic samples can be sparsified in a randomized fashion while maintaining AO or AnO of the planners. We term this regime as semi-deterministic sampling. We believe that this new model, and its analysis, is interesting in its own right.

This paper is a revised extended version of a conference paper (Solovey and Kleinbort, 2018), that was presented at Robotics: Science and Systems, 2018. The present version includes proofs that were omitted in the conference version, additional experimental results, and a table with values of $\gamma'$. It also presents an extension of the analysis to FMT$^*$, BTT, RRG, and discusses the implications of our results to the multi-robot planners dRRT$^*$ and M$^*$. Another important new addition is the semi-deterministic sampling setting, which was described in the previous paragraph.

### 1.2. Organization

In Section 2 we discuss related work. Then we proceed to basic definitions and problem statement in Section 3. In that section, we also include a precise description of the robotic system our theory is developed for. Our central contribution (Theorem 1), which states the existence of a critical radius $r_n^\gamma$ with respect to PRM, is presented in Section 4. This section also contains an outline of the proof. In Section 5 we lay the foundations of the proof and discuss...
important aspects of continuum percolation, that would later be employed in the main proof, which appears in Section 6. In Section 7 we extend the analysis to \textit{FMT}\textsuperscript{*}, \textit{BTT}, \textit{RGG}, and discuss the implications of our results to the multi-robot planners \textit{dPRM}\textsuperscript{*} and \textit{M}\textsuperscript{*}. In Section 8 we present experimental work, which validates our theory. We conclude this paper with directions for future work (Section 9).

In the appendix we include proofs that were omitted from the main document and a table with values of \textit{\gamma}. We also present in the appendix a new analysis for the aforementioned planners when constructed with deterministic samples, which are then sparsified in a randomized fashion. We believe that this new model, and its analysis, is interesting in its own right.

2. Related work

This section is devoted to a literature review of sampling-based planners with emphasis on their theoretical analysis. Sampling-based techniques were first described in the mid 1990s and several prominent examples of that time, that are still used extensively today, include \textit{PRM} (Kavraki et al., 1996), \textit{RRT} (LaValle and Kuffner Jr., 2001), and \textit{EST} (Hsu et al., 1999). Those planners are also known to be PC (see Hsu et al., 1999; Kavraki et al., 1998; Michal et al., 2019; Ladd and Kavraki, 2004).

More recent work has been concerned with the quality of the returned solution. Nechushtan et al. (2010) were one of the first to address this matter in a mathematically rigorous manner. This work proved that \textit{RRT} can produce arbitrarily-bad (long) paths with non-negligible probability.

The influential work of Karaman and Frazzoli (2011) laid the theoretical foundations for analyzing quality in sampling-based planning. The authors introduced a set of techniques to prove AO. Using their framework, they showed that the following algorithms are AO: \textit{PRM}\textsuperscript{*}, which is a special case of \textit{PRM} (throughout this work we will refer to the more general algorithm \textit{PRM} rather than \textit{PRM}\textsuperscript{*}) with a specific value of the connection radius \(r_n\); an AO variant of \textit{RRT} (LaValle and Kuffner Jr., 2001) termed \textit{RRT}\textsuperscript{*}; \textit{RGG}, which can be viewed as a combination between \textit{RRT} and \textit{PRM}. The analysis in Karaman and Frazzoli (2011) establishes that \(r_n = \Theta\left((\log n/n)^{1/d}\right)\) guarantees AO, where the configuration space of the robot is assumed to be \([0, 1]^d\). This indicates that the expected number of neighbors used per vertex should be \(O(\log n)\). The authors also proved that for sufficiently small radii of order \(O(n^{-1/d})\) the planner is guaranteed to fail (asymptotically) in finding any solution.

Following the breakthrough of Karaman and Frazzoli (2011), other AO planners have emerged (see, e.g., Alterovitz et al., 2011; Arslan and Tsiotras, 2013; Gammell et al., 2015). Janson et al. (2015) introduced \textit{FMT}\textsuperscript{*}, which is a single-query planner that traverses an implicitly represented \textit{PRM} graph, and is comparable in performance to \textit{RRT}\textsuperscript{*}. The authors refined the proof technique of Karaman and Frazzoli (2011), which allowed them to slightly reduce the connection radius \(r_n\) necessary to \textit{FMT}\textsuperscript{*} and \textit{PRM} to achieve AO. We do mention that here again \(r_n = \Theta\left((\log n/n)^{1/d}\right)\). \textit{BFMT}\textsuperscript{*}, which is a bidirectional version of \textit{FMT}\textsuperscript{*}, was introduced by Starek et al. (2015). In this paper the authors also proved that the success rate of \textit{PRM}, \textit{FMT}\textsuperscript{*}, \textit{BFMT}\textsuperscript{*} can be lower bounded by \(1 - \Theta(n^{-\eta/d}\log^{-1/d} n)\), where \(\eta > 0\) is a tuning parameter. In this context, we also mention the work of Dobson et al. (2015), which bounds the success rate with an expression that depends on the amount of deviation from the optimum.

A recent work by Kiril (2018) developed a different method for analyzing sampling-based planners. It exploits a connection with \textit{random geometric graphs} (RGGs), which have been extensively studied (see, e.g., Penrose, 2003). Their work shows that one can slightly reduce the \textit{PRM} and \textit{FMT}\textsuperscript{*} radius obtained in Janson et al. (2015). Furthermore, the connection with RGGs yields additional analyses of different extensions of \textit{PRM}, which have not been analyzed before in a mathematically rigorous setting.

A number of methods have been developed to reduce the running time or space requirements of existing planners by relaxing AO constraints to AnO. For instance, \textit{LBT-RRT} (Salzman and Halperin, 2016) interpolates between the quick \textit{RRT} and the AO yet slower \textit{RGG}, according to a pre-defined parameter. \textit{MPLB} (Salzman and Halperin, 2015) can be viewed as a relaxation of \textit{FMT}\textsuperscript{*}. \textit{SPARS2} (Dobson and Bekris, 2014) and \textit{RSEC} (Salzman et al., 2014) perform a sparsification of \textit{PRM} in an online or offline fashion, respectively, to reduce the space footprint of the produced roadmap.

2.1. Extensions

The aforementioned papers deal mainly with the cost function of \textit{path length}. Two recent works (Solovey and Halperin, 2016b, 2017) considered the \textit{bottleneck-pathfinding} problem in a sampling-based setting and introduced the \textit{BTT} algorithm, which traverses an implicitly represented \textit{PRM} graph. The bottleneck-cost function, which arises for instance in high-clearance multi-robot motion (Solovey and Halperin, 2017) and Fréchet matching between curves (Holladay et al., 2017), is defined as follows: Every robot configuration \(x\) is paired with a value \(M(x)\), and the cost of a path is the maximum value of \(M\) along any configuration on the path. It was shown (Solovey and Halperin, 2016b, 2017) that \textit{BTT} is AO, with respect to bottleneck cost, for the reduced connection radius that was obtained by Kiril (2018).

The results reported until this point have dealt exclusively with holonomic robotic systems. Two recent papers by Schmerling et al. (2015a,b) developed the theoretical foundations of \textit{PRM} and \textit{FMT}\textsuperscript{*} -flavored planners when applied to robots having differential constraints. Li et al.
(2016) developed an AO algorithm that does not require a steering function, as PRM for instance does. Interestingly, the authors of the last paper also obtain bounds on the rate of convergence of their algorithm.

3. Preliminaries

We provide several basic definitions that will be used throughout the paper. Given two points $x, y \in \mathbb{R}^d$, denote by $\|x - y\|$ the standard Euclidean distance. Denote by $B_r(x)$ the $d$-dimensional ball of radius $r > 0$ centered at $x \in \mathbb{R}^d$ and $B_r(\Gamma) = \bigcup_{x \in \Gamma} B_r(x)$ for any $\Gamma \subseteq \mathbb{R}^d$. Similarly, given a curve $\pi : [0, 1] \to \mathbb{R}^d$ define $B_r(\pi) = \bigcup_{t \in [0, 1]} B_r(\pi(t))$. Given a subset $D \subseteq \mathbb{R}^d$ we denote by $|D|$ its Lebesgue measure. All logarithms are at base $e$.

Throughout the paper we will use the standard notation for asymptotic bounds: let $f = f(n), g = g(n)$ be two functions. The notation $f(n) = O(g(n))$ indicates that there exist constants $n_0 > 0$, $c > 0$, such that for all $n > n_0$, it holds that $f(n) \leq c \cdot g(n)$. Similarly, $f(n) = \Omega(g(n))$ if there exist constants $n_0' > 0$, $c' > 0$, such that for all $n < n_0'$, it holds that $f(n) \geq c' \cdot g(n)$. In addition, $f(n) = O(g(n))$ if and only if $f(n) = O(g(n))$ and $f(n) = \Omega(g(n))$. The notation $f(n) = \omega(g(n))$ indicates that $\lim_{n \to \infty} f(n)/g(n) \rightarrow \infty$, and $f(n) = \omega(g(n))$ indicates that $\lim_{n \to \infty} f(n)/g(n) \to 0$.

3.1. Motion planning

Denote by $C$ the configuration space of the robot, and by $\mathcal{F} \subseteq C$ the free space, i.e., the set of all collision-free configurations. Though our proofs may be extended to more complex robotic systems (see the discussion in Section 9), in this work we investigate the geometric (holonomic) setting of the problem in which no constraints are imposed on the motion of the robot. In addition, we assume that $C$ is some subset of the Euclidean space. In particular, $C = [0, 1]^d \subseteq \mathbb{R}^d$ for some fixed $d \geq 2$. We also assume that for any two configurations $x, x' \in C$ the robot is capable of following precisely the straight-line path from $x$ to $x'$.

Given start and target configurations $s, t \in \mathcal{F}$, the problem consists of finding a continuous path (curve) $\pi : [0, 1] \to \mathcal{F}$ such that $\pi(0) = s$, $\pi(1) = t$. That is, the robot starts its motion along $\pi$ on $s$, and ends in $t$, while remaining collision free. An instance of the problem is defined for a given $(\mathcal{F}, s, t)$, where $s, t \in \mathcal{F}$. We consider in this work the standard setting in which the robot’s environment, or workspace, is known a priori (although $\mathcal{F}$ is not), in contrast to the much more complex setting in which the robot needs to sense it environment to gather information about the locations of obstacles (see, e.g., Janson et al., 2018).

3.2. Cost function

It is usually desirable to obtain paths that minimize a given criterion. In this paper we consider the following two cost functions.

Definition 1. Given a path $\pi$, its length is

$$c_1(\pi) = \sup_{n \in \mathbb{N}, 0 = t_0 \leq t_1 \cdots \leq t_n = 1} \sum_{i=2}^{n} \| \pi(t_i) - \pi(t_{i-1}) \|$$

Definition 2. Given a path $\pi$, and a cost map $\mathcal{M} : C \to \mathbb{R}$, its bottleneck cost is

$$c_b(\pi, \mathcal{M}) = \max_{\tau \in [0, 1]} \mathcal{M}(\pi(\tau))$$

We proceed to describe the notion of robustness, which is essential when discussing properties of sampling-based planners. Given a subset $\Gamma \subseteq C$ and two configurations $x, y \in \Gamma$, denote by $\Pi_{x,y}^\Gamma$ the set of all continuous paths, whose image is in $\Gamma$, that start in $x$ and end in $y$, i.e., if $\pi \in \Pi_{x,y}^\Gamma$, then $\pi : [0, 1] \to \Gamma$ and $\pi(0) = x, \pi(1) = y$.

Definition 3. Let $(\mathcal{F}, s, t)$ be a motion-planning problem. A path $\pi \in \Pi_{s,t}^\mathcal{F}$ is robust if there exists $\delta > 0$ such that $B_\delta(\pi) \subseteq \mathcal{F}$. We also say that $(\mathcal{F}, s, t)$ is robustly feasible if there exists such a robust path.

Definition 4. The robust optimum with respect to $c_\ell$ is defined as

$$c_\ell^* = \inf \left\{ c_\ell(\pi) \middle| \pi \in \Pi_{s,t}^\mathcal{F}, \pi \text{ is robust} \right\}$$

The corresponding definition for the bottleneck cost is slightly more involved.

Definition 5. Let $\mathcal{M}$ be a cost map. A path $\pi \in \Pi_{s,t}^\mathcal{F}$ is $\mathcal{M}$-robust if it is robust and for every $\varepsilon > 0$ there exists $\delta > 0$ such that for every $x \in B_\delta(\pi)$, $\mathcal{M}(x) \leq (1 + \varepsilon) c_\ell(\pi, \mathcal{M})$. We also say that $\mathcal{M}$ is well behaved if there exists at least one $\mathcal{M}$-robust path.

Definition 6. The robust optimum with respect to $c_b$ is defined as

$$c_b^* = \inf \left\{ c_b(\pi, \mathcal{M}) \middle| \pi \in \Pi_{s,t}^\mathcal{F}, \pi \text{ is } \mathcal{M} - \text{robust} \right\}$$

3.3. Poisson point processes

We draw our main analysis techniques from the literature of continuum percolation, where point samples are generated with the following distribution. Thus we will use this point distribution in PRM, which would be formally defined in the following section.

Definition 7 (Franceschetti and Meester, 2008). A random set of points $\mathcal{X} \subseteq \mathbb{R}^d$ is a Poisson point process (PPP) of density $\lambda > 0$ if it satisfies the following conditions.

1. For mutually disjoint domains $D_1, \ldots, D_k \subseteq \mathbb{R}^d$, the random variables $|D_1 \cap \mathcal{X}|, \ldots, |D_k \cap \mathcal{X}|$ are mutually independent.
2. For any bounded domain $D \subseteq \mathbb{R}^d$ we have that for every $k \geq 0$,
\[ \Pr[|\mathcal{X} \cap D| = k] = e^{-\lambda |D|}(\lambda |D|)^k / k! \]

Informally, property (1) states that the number of points from \( \mathcal{X} \) in any two disjoint sets \( D_i, D_j \subset \mathbb{R}^d \) is independent. Another useful property that follows from (2) is that the expected number of points in a certain region is known precisely and corresponds to the volume of this region. That is, for any \( D \subset \mathbb{R}^d \) it holds that \( E(|\mathcal{X} \cap D|) = \lambda |D| \).

The following provides a simple recipe for generating PPP.

**Claim 1** (Franceschetti and Meester, 2008). Let \( N \) be a Poisson random variable with mean \( \lambda \). For a given \( z \in \mathbb{Z}^d \) draw a sample \( N_z \in N \) and define \( \mathcal{X}_z = \{ X_1, \ldots, X_{N_z} \} \) to be \( N_z \) points chosen independently and uniformly at random from \( z + [0, 1]^d \). Then \( \mathcal{X} = \bigcup_{z \in \mathbb{Z}^d} \mathcal{X}_z \) is a PPP of density \( \lambda \).

It will be convenient to think about PRM as a subset of the following RGG. We describe various properties of this graph in later sections.

**Definition 8** (Penrose, 2003). Let \( \mathcal{X} \subset \mathbb{R}^d \) be a PPP. Given \( r > 0 \), the RGG \( \mathcal{G}(\mathcal{X}; r) \) is an undirected graph with the vertex set \( \mathcal{X} \). Given two vertices \( x, y \in \mathcal{X} \), \( (x, y) \in \mathcal{G}(\mathcal{X}; r) \) if \( \| x - y \| \leq r \).

### 4. Analysis of PRM

In this section, we provide a mathematical description of PRM, which essentially maintains an underlying RGG with PPP samples. We proceed to describe our main contribution (Theorem 1), which is concerned with the conditions for which PRM converges to the (robust) optimum. We then provide an outline of the proof, in preparation for the following sections.

Recall that the configuration space of the robot is represented by \( \mathcal{C} = [0, 1]^d \), and the free space is denoted by \( \mathcal{F} \subset \mathcal{C} \). The motion-planning problem \((\mathcal{F}, s, t)\) will remain fixed throughout this section.

Recall that PRM accepts as parameters the number of samples \( n \in \mathbb{N} \) and a connection radius \( r_n \). Denote by \( \mathcal{X}_n \) a PPP with mean density \( n \). In relation to the definitions of the previous section, the graph data structure obtained by the preprocessing stage of PRM can be viewed as an RGG. For instance, when \( \mathcal{F} = \mathcal{C} \), the graph obtained by PRM is precisely \( \mathcal{G}(\mathcal{X}_n \cap [0, 1]^d; r_n) \). In the more general case, when \( \mathcal{F} \subset \mathcal{C} \), PRM produces the graph \( \mathcal{G}(\mathcal{X}_n \cap \mathcal{F}; r_n) \). As \( \mathcal{F} \) can be non-convex, we emphasize that the latter notation describes the maximal subgraph of \( \mathcal{G}(\mathcal{X}_n; r_n) \) such that vertices and edges are contained in \( \mathcal{F} \).

In the query stage, recall that PRM accepts two configurations \( s, t \in \mathcal{F} \), which are then connected to the preprocessed graph. Here we slightly diverge from the standard definition of PRM in the literature. In particular, instead of using the same radius \( r_n \) when connecting \( s, t \), we use the (possibly larger) radius \( r_n' \). The graph obtained after query is formally defined as follows.

**Definition 9.** The PRM graph \( \mathcal{P}_n \) is the union between \( \mathcal{G}(\mathcal{X}_n \cap \mathcal{F}; r_n) \) and the supplementary edges

\[ \bigcup_{y \in [s, t]} \{ (x, y) | x \in \mathcal{X}_n \cap B_{r_n}(y) \land xy \subset \mathcal{F} \} \]

**Remark.** We emphasize that the larger radius \( r_n' \) is only used when \( s, t \) are connected to the preprocessed RGG.

We reach our main contribution.

**Theorem 1.** Suppose that \((\mathcal{F}, s, t)\) is robustly feasible. Then there exists a critical radius \( r_n = \gamma' n^{-1/d} \), where \( \gamma' \) is a constant (see Appendix C), such that the following holds.

1. If \( r_n < r_n' \) and \( r_n' = \infty \), then PRM fails to find a solution a.a.s.
2. Suppose that \( r_n > r_n' \). There exists \( \beta_0 > 0 \) such that for \( y_n' = \beta_0 n \), where \( \beta > \beta_0 \) and any \( \epsilon > 0 \) the following holds with probability \( 1 - O(n^{-1}) \):
   1. \( \mathcal{P}_n \) contains a path \( \pi_n \in \Pi_{s, t}^F \) with \( c_s(\pi_n) \leq (1 + \epsilon)c_s^* \), where \( c_\pi \) is independent of \( n \);
   2. \( \mathcal{M} \) is well behaved, then \( \mathcal{P}_n \) contains a path \( \pi_n' \in \Pi_{s, t}^F \) with \( c_s(\pi_n', \mathcal{M}) \leq (1 + \epsilon)c_s^* \).

The theorem proves the existence of a critical connection radius \( r_n = \gamma' n^{-1/d} \), where \( \gamma' > 0 \) is a constant: if \( r_n < r_n' \), then PRM is guaranteed to fail. Above the threshold, i.e., when \( r_n > r_n' \), PRM is AO for the bottleneck cost, and ANO with respect to the path-length cost.

**Remark.** Theorem 1 also holds, with a slight modification, when the PPP is replaced with the standard binomial point process (BPP), in which the number of sampled points is fixed a priori. The only difference is that the probability of success should be reduced from \( 1 - O(n^{-1}) \) to \( 1 - O(n^{-1/2}) \). This follows from Lemma 1 of Friedrich et al. (2013), which states that if a property holds for PPP, then it also holds for BPP, albeit with slightly smaller probability.

### 4.1. Outline of proof

For the remainder of this section, we briefly describe our technique for proving this theorem, in preparation for the full proof, which is given in Section 6. The critical radius \( r_n' \) defined above, coincides with the percolation threshold, which determines the emergence of a connected component of \( \mathcal{G} \) that is of infinite size. In particular, if \( r_n < r_n' \), then \( \mathcal{G}(\mathcal{X}_n; r_n) \) breaks into tiny connected components of size \( O(\log n) \) each. Thus, unless \( s, t \) are infinitesimally close, no connected component can have both \( s \) and \( t \) simultaneously.

More interestingly, the radius of \( r_n' \) leads to the emergence of a unique infinite component of \( \mathcal{G}(\mathcal{X}_n; r_n) \). That is, in such a case one of the components of \( \mathcal{G}(\mathcal{X}_n; r_n) \) must contain an infinite number of vertices (Section 5.1). Denote this component by \( C_\infty \).
In contrast to \( G(X_n^r; r_n) \), which is defined for the unbounded space \( \mathbb{R}^d \), our motion-planning problem is bounded to \( C = [0, 1]^d \). Thus, the next step is to investigate the properties of \( G(X_n^r; r_n) \) when restricted to \( [0, 1]^d \) (Section 5.2). Denote by \( C_n \) the largest connected component of \( C_n \cap [0, 1]^d \). This structure plays a key role in our proof (see Section 6): with high probability (to be defined), there exist vertices of \( C_n \) that are sufficiently close to \( s \) and \( t \), respectively, so that a connection between the two vertices can be made through \( C_n \).

Of course, this overlooks the fact that some portions of \( C_n \) lie in forbidden regions of \( C \). Thus, we also have to take the structure of \( F \) into consideration. To do so, we rely on Penrose and Pisztora (1996) to prove that any small subset of \( [0, 1]^d \) must contain at least one point of \( C_n \) (see Lemmas 2 and 3). This allows us to trace the robust optimum (and collision-free) paths with points from \( C_n \).

The final ingredient, which allows to bound the path length along \( G(X_n^r; r_n) \cap [0, 1]^d \), is Theorem 4. It states that the distance over this graph is proportional to the Euclidean distance between the end points. This also ensures that the trace points from \( C_n \) can be connected with collision-free paths over the graph.

To conclude, in Section 5 we provide background on continuum percolation in unbounded and bounded domains, and prove two key lemmas (Lemmas 2 and 3). In Section 6, we return to the setting of motion planning and utilize the aforementioned results in the proof of Theorem 1.

5. Elements of continuum percolation

In this section, we describe some of the properties of the unbounded graph \( G(X_n^r; r_n) \) that will be employed in our analysis in the following section.

5.1. The basics

A fundamental question is when \( G \) contains an infinite connected component around the origin.

**Definition 10.** The percolation probability \( \theta(n, r) \) is the probability that the origin \( o \in \mathbb{R}^d \) is contained in a connected component of \( G(X_n^r; r_n) \) of an infinite number of vertices. That is, if \( C_o \) denotes the set of vertices connected to \( o \) in the graph, then \( \theta(n, r) = \Pr(\lvert C_o \rvert = \infty) \).

We say that a graph *percolates* if and only if \( \theta(n, r) > 0 \). Note that the selection of the origin is arbitrary, and the following result can be obtained for any \( x \in \mathbb{R}^d \) alternative to \( o \).

**Theorem 2** (Grimmett, 1999, Theorem 12.35). There exists a critical radius \( r_n^* = \gamma^* n^{-1/d} \), where \( \gamma^* \) is a constant, such that \( \theta(n, r_n) = 0 \) when \( r_n < r_n^* \), and \( \theta(n, r_n) > 0 \) when \( r_n > r_n^* \).

The following lemma states that the infinite connected component exists with probability strictly 0 or 1.

**Lemma 1.** Let \( \psi(n, r) \) be the probability that \( G(X_n^r; r) \) contains an infinite connected component, i.e., without conditioning on any specific additional vertex. Then \( \psi(n, r) = 0 \) when \( \theta(n, r) = 0 \) and \( \psi(n, r) = 1 \) when \( \theta(n, r) > 0 \).

**Proof.** Suppose that \( \theta(n, r) = 0 \), and for any \( x \in \mathbb{R}^d \), denote by \( \theta_x(n, r) \) the percolation probability of \( G(X_n^r \cup \{x\}; r) \), and note that \( \theta_x(n, r) = 0 \). Define \( Z_r = \{ r \cdot z \in \mathbb{Z}^d \} \), and observe that for any \( y \in \mathbb{R}^d \) there exists \( z \in Z_r \) such that \( y \in B_r(z) \). By definition, \( G(X_n^r \cup \{x\}; r) \) percolates, i.e., the infinite component touches \( B_r(x) \), if and only if there exists \( y \in C_n \) such that \( \| x - y \| \leq r \). By definition of \( Z_r \), if the latter event occurs then there exists \( z \in Z_r \) such that \( \| z - y \| \leq r \). Thus, by applying the union bound, and noting that a sum of countable number of zeros is still zero, we establish that

\[
\psi(n, r) = \Pr[\exists x \in \mathbb{R}^d, G(X_n^r \cup \{x\}; r) \text{ percolates}] = \Pr[\exists z \in Z_r, G(X_n^r \cup \{z\}; r) \text{ percolates}] \\
\leq \sum_{z \in Z_r} \theta_z(n, r) = 0
\]

For the other direction we employ Kolmogorov’s zero-one law (see, e.g., Bollobás and Riordan, 2006: Theorem 1, p.36). Informally, it states that an event, e.g., existence of an infinite connected component in \( G \), that occurs independently of any finite subset of independent random variables, e.g., points from \( X_n \), has a ability of either 0 or 1. Thus, as \( \psi(n, r) \equiv \theta(n, r) \) it immediately follows that \( \psi(n, r) = 1 \) when \( \theta(n, r) > 0 \).

The following theorem establishes that the infinite connected component is unique.

**Theorem 3** (Meester and Roy, 1994, Theorem 2.3). With probability 1, \( G(X_n^r; r) \) contains at most one infinite connected component.

5.2. Bounded domains

We study different properties of \( G(X_n^r; r) \) when it is restricted to the domain \( [0, 1]^d \). In case that \( \theta(n, r) > 0 \), we use \( C_\infty \) to refer to the infinite connected component of the unbounded graph \( G(X_n^r; r) \). Note that \( C_\infty \) exists (Lemma 1) and is unique (Theorem 3) with probability 1.

Denote by \( C_n \) the largest connected component of \( C_n \cap [0, 1]^d \). By definition, \( C_n \) is also a subgraph of \( G(X_n^r \cap [0, 1]^d; r_n) \). The following lemma shows that with high probability all the points from \( C_n \), that are sufficiently close to the center of \( [0, 1]^d \), are members of \( C_n \).

**Lemma 2.** Let \( r_n > r_n^* \). Define

\[
H_n = [0, 1]^d \setminus B_{1/\log n}(\mathbb{R}^d \setminus [0, 1]^d)
\]

Denote by \( \mathcal{E}_n \) the event that \( C_\infty \cap H_n \subset C_n \). Then there exist \( n_0 \in \mathbb{N} \) and \( \alpha > 0 \) such that for any \( n > n_0 \) it holds that...
having at least one point from $C_n$ in a (small) subregion of $H_n$. Note that the value $\beta$ corresponds to the value used in Theorem 1.

**Lemma 3.** Let $r_n > r'_n$. Define $H'_n \subset H_n$ to be a hypercube of side length

$$H'_n = \Theta \log^{-1/d} n$$

Denote by $\mathcal{E}_n^2$ the event that $H'_n \cap C_n \not= \emptyset$. Then there exists $n_0 \in \mathbb{N}$ and $\beta_0 > 0$ such that for any $n > n_0$, $\beta > \beta_0$ it holds that $\Pr[\mathcal{E}_n^2 | \mathcal{E}_n^2] \gg 1 - n^{-1}$.

**Proof.** Define $G_n = G(X_n, r_n) \cap H'_n \cap E[|X_n \cap H'_n|]$ and observe that $n' = n \cdot |H'_n| = \beta d \log^{\beta(d-1)} n$.

We treat $G_n$ as a subset of $\mathbb{R}^d$ in order to apply a rescaling argument. Observe that (scalar) multiplication of every point of $H'_n$ with $1/H'_n$ yields a translation of $[0, 1]^d$. We will use the superscript $1/H'_n$ to describe this rescaling to a given object. For instance, applying the same transformation on $G_n$ yields the graph $G_n^{1/H'_n}$, which has the same topology as $G_n$. Denote by $x_i \in H'_n$ the (lexicographically) smallest point of $H'_n$, i.e., $x_i = (1/\log n, \ldots, 1/\log n)$. Note that

$$G_n^{1/H'_n} - x_i^{1/H'_n} = G(X_n \cap H'_n \cap [0, 1]^d, r_n/H'_n)$$

where the minus sign in the left-hand side represents a translation by a vector. This implies that $G_n$, which is defined over $H'_n$ behaves as $G(X_n \cap H'_n; r_n)$. This allows to leverage Theorem 1 from Penrose and Pisztora (1996), which bounds the number of vertices from the unbounded component. In particular, there exists $\beta_0 > 0$ such that

$$\Pr[C_n \cap H'_n \cap H'' \not= \emptyset] \gg 1 - \exp\left(-\beta_0^{(d-1)} n^{d/d}\right)$$

$$= 1 - \exp\left(-\beta_0^{(d-1)} B^{d-1} \log n\right)$$

$$\gg 1 - \exp(-\log n) = 1 - n^{-1}$$

While this is an overkill for our purpose, it does the job in proving that $\Pr[C_n \cap H'_n \not= \emptyset] \gg 1 - n^{-1}$. As we assume that $\mathcal{E}_n^2$ holds (Lemma 2), it follows that $C_n \cap H'_n \not= \emptyset$ holds with probability at least $1 - n^{-1}$. □

The following statement allows to bound the graph distance between two connected vertices. We endow every edge of the graph with a length attribute that represents the Euclidean distance between the edges’ endpoints. For every two vertices $x, x'$ of $G$, dist($G, x, x'$) denotes the length of the shortest (weighted) path on $G$ between the two vertices.

**Theorem 4** (Friedrich et al., 2013, Theorem 3). Let $r_n > r_n^*$. There exists a constant $\xi \gg 1$, independent of $n$, such that $\Pr[\mathcal{E}_n^2] = 1 - O(n^{-1})$, where the event $\mathcal{E}_n^2$ is defined as follows: for any two vertices $x, x'$ in the same connected component of $G(X_n \cap [0, 1]^d; r_n)$, with $\|x - x'\| = o(r_n)$, it holds that dist($G_n, x, x'$) $\leq \xi \|x - x'\|$.
6. Proof of Theorem 1

Proofs for all the three settings of the main theorem are given individually in the following subsections.

6.1. Case (i)

Here we provide proof for the first (and easy) part of the theorem, which states that PRM fails when \( r_n < \gamma^{*} n^{-1/d} \), even with \( r''_n = \infty \). We mention that our proof is a simpler and shorter version of a similar proof that was given in Karaman and Frazzoli (2011) for a slightly different setting.

We show in the following that in the subcritical regime, i.e., \( r_n < r'_n \), the graph breaks into many small connected components. In particular, the probability of having the largest connected component of size \( m \) decays exponentially in \( m \). Denote by \( L(G) \) the size of the largest connected component in \( G \).

**Proposition 1.** Let \( r_n < r'_n \). Then \( L(G_n) = O(\log n) \text{ a.a.s.} \)

**Proof.** First, we start by stating that there exist a constant \( \zeta > 0 \) and an integer \( m_0 \) such that for all \( m \geq m_0 \) it holds that \( \Pr[L(G_n) \geq m] \leq ne^{-\xi m} \). This is a simplified version of Proposition 11.2 of Penrose (2003), which is given for \( n \) uniformly sampled points, to the case of PPP, which we have here. In particular, the original proof relies on a relation between these two distributions, which transforms uniform sampling into a PPP and induces an additional factor that is not necessary in our setting. In particular, the factor \( \exp(-\mu n) \) that appears in Equation 11.1 of Penrose (2003) should be eliminated.

To conclude the proof, we set \( m = \alpha \cdot \log n \), where \( \alpha > 1/\zeta \), similarly to the proof of Theorem 1.1, Equation 11.4 of Penrose (2003). Observe that \( ne^{-\xi m} \to 0 \) as \( n \) tends to \( \infty \).

Now, consider the configuration space \( C = [0,1]^d \) depicted in Figure 3 (similar examples can be devised for any \( d \geq 2 \)). The white and blue regions represent the free space. Observe that any path connecting \( s \) to \( t \) must go through the blue region, whose width is greater than 1/2. Also note that any point contained in the blue region cannot be connected by an edge from \( s \) or \( t \), which deems the large value of \( r''_n \) as irrelevant.

Suppose that \( G_n = G(X_n, r_n) \) contains a path \( \pi_n \) (black dashed curve) connecting \( s \) and \( t \), and denote by \( \pi''_n \subset \pi_n \) the subpath that is contained in the blue region. In addition, denote by \( s' \) and \( t' \) the first and last points along \( \pi''_n \), respectively. Obviously \( 1/2 \leq \left\| s' - t' \right\| \leq c(\pi_n) \). Then, the number of edges of \( G_n \), which induce \( \pi''_n \), is at least \( \left\| s' - t' \right\| / r_n = \Omega(n^{1/d}) \). However, this is in contradiction with the fact that every connected component of \( G_n \) is of logarithmic size in \( n \) (Proposition 1).

6.2. Case (ii)/(I)

We provide a full proof for the positive setting with length cost. Namely, we show that if \( r_n > r'_n \), then there exists \( \beta_i > 0 \) such that for \( r''_n = \frac{\beta \log^{1/4}(n)}{n} \), where \( \beta > \beta_0 \), and any \( \varepsilon > 0 \) the following holds with probability \( 1 - O(n^{-1}) \): \( P_n \) contains a path \( \pi_n \in \Pi_{r''} \) with \( c(\pi_n) \leq (1+\varepsilon)\xi c' \), where \( \xi \) is independent of \( n \). The proof for the bottleneck case, which appears later on, is very similar to the length case.

Suppose that \( r_n > r'_n \) and \( r''_n = \frac{\beta \log^{1/4}(n)}{n} \), \( \beta > \beta_0 \). For simplicity, we set \( r_n = \gamma n^{-1/d} \), where \( \gamma > \gamma^* \). By Lemma 1 and Theorem 3, \( G(X_n; r_n) \) contains a unique infinite connected component \( C_x \). Recall that \( C_n \) denotes the largest connected component of \( C_x \cap [0,1]^d \). Also note that \( r''_n = h'_n \), where \( h'_n \) is defined in Lemma 3.

Recall that \( c'_x \) denotes the robust optimum, with respect to path length (Definition 4). Fix \( \varepsilon > 0 \). By definition, there exists a robust path \( \pi_i \in \Pi_{r''} \) and \( \delta > 0 \) such that \( c(\pi_i) \leq (1+\varepsilon)c'_x \) and \( B_2(\pi_i) \subset F \). See the illustration in Figure 4.

We now define a sequence of \( k \) points \( p_1, \ldots, p_k \) along \( \pi_i \) that are separated by exactly \( \delta / 2\xi \) units, where \( \xi \) is as defined in Theorem 4. In particular, define \( k = \lceil c(\pi_i) \cdot 2\delta / \delta \rceil \), set \( p_1 = s, p_k = t \), and assign \( p_i \) along \( \pi_i \), such that \( c(\pi_i) \cdot (i-1) \delta / \delta \) \( \pi_i \) represents the subpath of \( \pi_i \) starting at \( p_{i-1} \) and ending at \( p_i \). Note that \( k \) is finite.

**Claim 2.** Denote by \( E^k_n \) the event that for all \( i \in [k] \) there exists \( q_i \in C_n \) such that \( q_i \in B_{r''}(p_i) \) and \( q_i \in C_n \). Then \( \Pr[E^k_n] \geq 1 - k \Pr(E^2_n) \text{ (see definition of } E^1_n, E^2_n \text{ in Lemmas 2 and 3, respectively).} \)

**Proof.** Define \( H'_n(x) \subset \mathbb{R}^d \) to represent a \( d \)-dimensional (axis-aligned) hypercube of side length \( h'_n \) that is centered in \( x \in \mathbb{R}^d \). Formally, \( H'_n(x) = x + h'_n \cdot [-1/2,1/2]^d \). Observe that \( H'_n(p_i) \subset H_n \) for \( n \) large enough. Also note that \( H'_n(p_i) \subset C_n \). Thus, the result follows from the union bound.

Suppose that \( E^1_n, E^2_n \) are satisfied. Let \( q_1, \ldots, q_k \in C_n \) be the points obtained from Claim 2. These points reside in a single connected component of \( G_n \). Define the path...
By Theorem 4 and the triangle inequality, it follows that
\[
    c_i(\pi_n) = \| s - q_1 \| + \text{dist}(G_n, q_1, q_k) + \| t - q_k \|
\]
\[
\leq 2r_n^d + \sum_{i=2}^{k} \text{dist}(G_n, q_{i-1}, q_i)
\]
\[
\leq o(1) + \sum_{i=2}^{k} \xi \| q_{i-1} - q_i \|
\]
\[
\leq o(1) + \xi \sum_{i=2}^{k} c_i(\pi_i^{d-1,i})
\]
\[
= o(1) + \xi c_i(\pi_e) \leq o(1) + (1 + \varepsilon)\xi c_i^e.
\]

We note that in order to eliminate the \( o(1) \) summand one can fix \( \varepsilon \) but work with \( \pi_e \) instead of \( \pi_n \), where \( \delta' < \varepsilon \). For simplicity, we chose to leave the proof as is.

To conclude, Claims 3 and 4 show the existence of a (collision-free) path \( \pi_n \) in \( \mathcal{P}_n \), whose length is at most \((1 + \varepsilon)\xi c_i^e\). It remains to bound the probability that \( \mathbf{e}_1^n, \mathbf{e}_3^n, \mathbf{e}_4^n \) hold simultaneously:

\[
    \Pr[\mathbf{e}_1^1 \land \mathbf{e}_3^1 \land \mathbf{e}_4^1] = \Pr[\mathbf{e}_3^1 \land \mathbf{e}_4^1 \land \mathbf{e}_1^1] \cdot \Pr[\mathbf{e}_1^1]
\]
\[
= (1 - \Pr[\mathbf{e}_3^1 \land \mathbf{e}_4^1 | \mathbf{e}_1^1]) \cdot \Pr[\mathbf{e}_1^1]
\]
\[
= (1 - \Pr[\mathbf{e}_3^1 \land \mathbf{e}_4^1 | \mathbf{e}_1^1]) \cdot \Pr[\mathbf{e}_1^1]
\]
\[
\geq (1 - (1 - \Pr[\mathbf{e}_3^1 | \mathbf{e}_1^1] - \Pr[\mathbf{e}_4^1 | \mathbf{e}_1^1]) \cdot \Pr[\mathbf{e}_1^1]
\]
\[
\geq (1 - O(n^{-1}) - k n^{-1}) \left(1 - \exp\left(-\alpha n^{-1/d} \log^{-1} n\right)\right)
\]
\[
= 1 - O(n^{-1}).
\]

6.3. Case (ii)(2)

We prove the AO of \( \text{PRM} \) for bottleneck cost, with appropriate values of \( r_n, r_n^d \). Namely, if \( r_n > r_n^d \) and \( \mathcal{M} \) is well behaved, then there exists \( \beta_0 > 0 \) such that for \( r_n^d = \theta \log^{(l-1)/d} n \), where \( \beta \geq \beta_0 \), and any \( \varepsilon > 0 \) the following holds with probability \( 1 - O(n^{-1}) \): \( \mathcal{P}_n \) contains a path \( \pi'_n \in \Pi_{x,t}^{\mathcal{F}} \) with \( c_{\phi}(\pi'_n, \mathcal{M}) \leq (1 + \varepsilon) c_\psi \). The proof follows very similar lines to that of the length cost. Fix \( \varepsilon > 0 \). For simplicity we assume that \( \varepsilon < 1 \) (the proof can be adapted to larger values of \( \varepsilon \)). By definition, there exists an \( \mathcal{M} \)-robust path \( \pi' \in \Pi_{x,t}^{\mathcal{F}} \) and \( \delta, \delta' > 0 \) such that:

(a) \( B_\psi(\pi') \subseteq \mathcal{F} \);
(b) \( c_\phi(\pi', \mathcal{M}) \leq (1 + \varepsilon/3) c_\psi \); and
(c) for all \( x \in B_\psi(\pi') : \mathcal{M}(x) \leq (1 + \varepsilon/3) c_\phi(\pi', \mathcal{M}) \).
Now, define $\delta^* = \min\{\delta, \delta'\}$. By substituting $\delta$ with $\delta^*$ and $\pi_n$ with $\pi'$ in the proof of Theorem 1(ii), it follows that if $E_n^1, E_n^2, E_n^3$ are satisfied, $P_n$ contains a path $\pi'_n \in \Pi^F_{\pi,t}$ such that $\text{Im}(\pi'_n) \subset B_0(\pi')$. This implies that

$$c_b(\pi'_n, M) \leq (1 + \varepsilon/3)c_b(\pi', M) \leq (1 + \varepsilon/3)\varepsilon c^*_b \leq (1 + \varepsilon)c^*_b$$

which concludes this proof.

7. Analysis of other planners

In this section, we describe the implications of Theorem 1 to additional planners, which are closely related to PRM. In particular, those planners maintain an implicit representation of a PRM graph, only a small portion of which is explicitly constructed. This allows to extend known properties concerning PRM to those planners. We mention that such a relation has been observed in previous work (Gammell et al., 2015; Solovey and Halperin, 2016b; Solovey et al., 2016).

In all the results in the following, the values $r_n$, $\gamma, \beta, \xi$ are identical to those in Theorem 1. Furthermore, for simplicity of presentation, we assume that $(F, s, t)$ is robustly feasible, and $M$ is well behaved, where relevant.

We first consider FMT* and BTT, and then proceed to RRG, whose analysis is more involved. In all the following algorithms we assume that the sampling method is a PPP.

7.1. The FMT* and the BTT planners

The following two corollaries are direct consequences of Theorem 1 as the two algorithms FMT* and BTT traverse an implicitly represented PRM graph while minimizing the cost functions $c_l$ and $c_b$, respectively.

**Corollary 1.** If $r_n < r_n^* \text{ and } r_n^* \to \infty$, then FMT* fails a.a.s. If $r_n > r_n^* \text{ and } r_n^* = \frac{\delta \log(1/4) - \delta}{\frac{\delta}{\tilde{n}}}$, then for any $\varepsilon > 0$ FMT* returns a path $\pi_n \in \Pi^F_{\pi,t}$ with $c_l(\pi_n, M) \leq (1 + \varepsilon)\xi c^*_b$ with probability $1 - O(n^{-1})$.

We remark that FMT* can ignore certain edges of the underlying PRM graph, which may result in a path of lower quality than that obtained by PRM. However, this situation only occurs for vertices which lie within distance of $r_n$ from $C \setminus F$ (see Janson et al., 2015: Remark 3.3). As our analysis of FMT in the previous section utilizes only vertices and edges that are far from obstacles our proofs extend to FMT* straightforwardly.

**Corollary 2.** If $r_n < r_n^* \text{ and } r_n^* \to \infty$, then BTT fails a.a.s. If $r_n > r_n^* \text{ and } r_n^* = \frac{\delta \log(1/4) - \delta}{\frac{\delta}{\tilde{n}}}$, then for any $\varepsilon > 0$ BTT returns a path $\pi_n \in \Pi^F_{\pi,t}$ with $c_b(\pi_n, M) \leq (1 + \varepsilon)c^*_b$ with probability $1 - O(n^{-1})$.

7.2. The RRG planner

We consider the incremental planner RRG, which can be viewed as a cross between RRT and PRM. Due to this relation we can extend the analysis of PRM to RRG. (The term "incremental" refers to planners that generate samples one after the other, and connect the current samples to previous ones.)

We introduce an incremental version of a PPP to extend our theory to RRG.

**Claim 5.** Let $N$ be a Poisson random variable with mean 1. At each iteration $i$ I draw a sample $N_i \in N$ and define $X_i = \{X_1, \ldots, X_{N_i}\}$ to be $N_i$ points chosen independently and uniformly at random from $[0, 1]^d$. The set $\mathcal{X} = \bigcup_{i=1}^{n} X_i$ obtained after $n$ iterations, is a PPP of density $\lambda = n$.

Note that the sum of $n$ independent and identically distributed (i.i.d.) Poisson random variables with means $\lambda_1, \lambda_2, \ldots, \lambda_n$ is a Poisson random variable with mean $\lambda = \sum_{i=1}^{n} \lambda_i$. Claim 5 follows directly from this property.

We now describe an adaptation of RRG for an incremental PPP. Given a start configuration $s \in F$ and a goal region $\mathcal{X}_{\text{goal}} \subseteq F$, RRG initializes a roadmap with a single node $s$. Let $\eta$ denote the constant used by RRG for local steering (see Karaman and Frazzoli, 2011). Similarly to PRM, RRG employs the connection radii $r_n^* \text{ and } r_n^*$. Let $N$ be a Poisson random variable with mean 1 and let $n_{i-1}$ denote the number of nodes of the constructed roadmap after $i-1$ iterations. At the $i$th iteration we draw a sample $N_i \in N$ and define $X_i = \{X_1, \ldots, X_{N_i}\}$ to be $N_i$ points chosen independently and uniformly at random from $[0, 1]^d$. RRG will iteratively process the samples $X_1, \ldots, X_{N_i}$, as follows: for a sample $X_k \in X_i$ RRG will first locate the nearest node $x_{n_{i-1}}$ to $X_k$. Let $x_{n_{i-1}}$ be a node at distance at most $\eta$ from $x_{n_{i-1}}$ at the direction of $X_k$. Here RRG will attempt to connect $x_{n_{i-1}}$ to $x_{n_{i-1}}$. If the connection attempt is successful, the new node $x_{n_{i-1}}$ will be added to the roadmap. Then, RRG will attempt to connect $x_{n_{i-1}}$ to all the existing nodes within a ball of radius $\min\{r_n^{\text{RRT}}, \eta\}$. It will additionally attempt to connect $x_{n_{i-1}}$ to $s$ if $\|x_{n_{i-1}} - s\| \leq \min\{r_n^{\text{RRT}}, \eta\}$. Note that for the $i$th iteration of RRG we set $n = n_{i-1}$.

The connection radii $r_n^{\text{RRT}}, r_n^{\text{RRG}}$ that will be used during the $i$th iteration are fixed, as both are functions of the number of nodes in the roadmap at the beginning of the iteration. Both radii will be set to decrease with the number of roadmap nodes, as in PRM. However, since we fix the radius at each iteration of RRG we use a slightly larger radius than the one obtained had we considered the current roadmap size. This will clearly keep all relevant connections and perhaps even add more.

**Theorem 5.** Suppose that $(F, s, \mathcal{X}_{\text{goal}})$ is robustly feasible (we extend Definition 3 to describe a robustly feasible path $(F, s, \mathcal{X}_{\text{goal}})$ for RRG). Let $R_n$ denote the roadmap constructed by RRG after $n$ iterations with
where \( r_n, r'_{n} \) are as in Theorem 1(ii), and \( \mu \) is any positive constant. Then for any \( \varepsilon > 0 \) the following holds with probability \( 1 - O(n^{-1}) \):

1. \( R_n \) contains a path \( \pi_n \in \Pi_{r_n}^F, \) with \( c_i(\pi_n) \leq (1 + \varepsilon)c_i^* \); and

2. if \( M \) is well behaved, then \( R_n \) contains a path \( \pi'_n \in \Pi_{r'}^F, \) with \( c_b(\pi'_n, M) \leq (1 + \varepsilon)c_b^* \).

### 7.2.1. Proof of Theorem 5

We provide proof only for \( c_i \), as the case for \( c_b \) is almost identical. In what follows, we consider a robust path \( \pi_e \in \Pi_{r_n}^F, \) for which \( c_i(\pi_e) \leq (1 + \varepsilon)c_i^* \). Define \( \delta > 0 \) denote the clearance of \( \pi_e \), and \( L \) denote its length. Finally, set \( \kappa = \min(\delta, \eta)/5 \).

Our proof uses Theorem 1 as a central ingredient, due to the following observation, which is formalized below: for \( n \) large enough, with probability \( 1 - O(e^{-n}) \), \( R_n \) contains a PRM roadmap \( \mathcal{P}_{n'} \) in the vicinity of \( \pi_e \), where \( n' \) is slightly smaller than \( n \). This in turn follows from the unique structure of \( RRG \), which can be viewed as a combination of \( PRM \) and \( RRT \), as \( RRG \) supplements the \( RRT \) tree with additional edges.

In particular, the following corollary can be deduced from the probabilistic completeness of \( RRT \) with samples from a PPP (see Appendix B), which does not rely on the parameters \( r_{n}^{RRG}, r'_{n}^{RRG} \).

**Claim 6.** It is possible to tile \( \pi_e \) with \( m = L/\kappa \) balls of radius \( \kappa \), such that after \( n \) iterations of \( RRG \), with probability at least \( 1 - a \cdot e^{-bn} \) for some constants \( a, b \in \mathbb{R}_{>0} \), every ball will contain at least one \( RRG \) vertex.

The following lemma formalizes the connection between \( RRG \) and \( PRM \): 

**Lemma 4.** Fix \( n \) and define \( n' = (1 + \mu)^d n \). Let \( R_n \) be an \( RRG \) graph constructed after \( n' \) iterations with connection radii \( r_n, r_n' \). With probability at least \( 1 - a \cdot e^{-bn} \), for some constants \( a, b' \in \mathbb{R}_{>0} \), the graph \( R_n \cap B_{e/2}(\pi) \) contains a \( PRM \) graph \( \mathcal{P}_{n'} \cap B_{e/2}(\pi) \), constructed with the radius \( r_{n'} \).

**Proof.** Consider the \( n - n' = n(1 - (1 + \mu)^{-d}) \) first iterations of \( RRG \). From Claim 6 we obtain that given a robust path \( \pi_n \), it can be tiled using a constant number of balls of radius \( \kappa \) such that with probability of at least \( 1 - ae^{-b(n-n')} = 1 - ae^{-bn} \) every ball will contain an \( RRG \) vertex after \( n \) iterations of \( RRG \), for some constants \( a, b \in \mathbb{R}_{>0} \), and \( b' = b(1 - (1 + \mu)^{-d}) \). Now, let us consider a new sample \( \pi' \in B_{e/2}(\pi) \) that is added in iteration \( n'' = n - n' \). By Claim 6, \( \mathcal{P}_{n'} \) includes (with certain probability) a node \( v \) that can be connected to \( \pi' \); that is, both \( \parallel x - v \parallel \leq \eta \) and the straight-line path from \( x \) to \( v \) is collision free. Thus, \( v \) is added as a node to \( RRG \) and then connected in a \( PRM \)-fashion to all its neighbors within a radius of \( r_{n''} \).

Note that the smallest value of the latter radius is when \( n'' = n \). Observe that

\[
r_{n''} = (1 + \mu)r_n = \gamma((1 + \mu)^{-d} n)^{-1/d} = \gamma(n')^{-1/d} = r_{n'}
\]

In addition,

\[
r_{n''} = (1 + \mu)r_n = \frac{\beta \log^{1/(d-1)} n}{n^{1/d}}
\]

This implies that all samples added after iteration \( n > n' \) are connected in a PRM fashion with a radius \( r_{n'} \). Thus, \( R_n \cap B_{e/2}(\pi) \) contains a \( PRM \) graph \( \mathcal{P}_{n'} \cap B_{e/2}(\pi) \), constructed with the radius \( r_{n'} \) with probability \( 1 - a \cdot e^{-bn} \), as required. \( \square \)

We are now ready to finish the proof of Theorem 5. Let us denote by \( B \) the event that the graph \( R_n \) contains a \( PRM \) graph \( \mathcal{P}_{n'} \cap B_{e/2}(\pi) \). Given that \( A \) occurs, we denote by \( B \) the event that the constructed roadmap \( R_n \cap B_{e/2}(\pi) \) maintains the good properties of \( PRM \), as defined in Theorem 1(ii). That is, \( \mathcal{P}_{n'} \cap B_{e/2}(\pi) \) contains a path \( \pi' \in \Pi_{r'}^F, \) with \( c_b(\pi') \leq (1 + \varepsilon)c_b^* \).

Given that \( A \) occurs, from Theorem 1 we obtain that \( B \) occurs on the \( PRM \)-like graph inside \( B_{e/2}(\pi) \) with probability of at least

\[
1 - O((n')^{-1}) = 1 - O((1 + \mu)^{-d} n^{-1}) = 1 - O(n^{-1})
\]

That is, \( \Pr[\mathcal{P}_{n'} \cap B_{e/2}(\pi)] \geq 1 - O(n^{-1}) \). In addition, Lemma 4 states that \( \Pr[\mathcal{P}_{n'}] \geq 1 - ae^{-bn} \) for some \( b' \). Note that we would like to bound the probability that both \( A \) and \( B \) occur simultaneously. From the definition of conditional probability we have that \( \Pr[A \cap B] = \Pr[\mathcal{P}_{n'} \cap B_{e/2}(\pi)] \cdot \Pr[A] \). Therefore,

\[
\Pr[A \cap B] \geq \left(1 - O(n^{-1})\right) \cdot \left(1 - ae^{-bn}\right)
\]

\[
\geq 1 - O(e^{-bn}) - O(n^{-1}) = 1 - O(n^{-1})
\]

### 7.3. Multi-robot planners

In this section, we briefly state the implications of our analysis of \( PRM \) for sampling-based multi-robot motion planning. In particular, we consider the planners \( M^* \) (Wagner and Choset, 2015), \( dRRT \) (Solovey et al., 2016), and \( dRRT^* \) (Shome et al., 2019).
The multi-robot setting involves \( m \geq 2 \) robots operating in a shared workspace. In this more challenging setting collisions between different robots must be avoided, in addition to standard robot–obstacle collisions. Single-robot sampling-based planners can be applied directly to the problem by considering the robot fleet as one highly complex robot. However, recently introduced planners that are tailored for the multi-robot case have proved to be much more effective than the aforementioned naive approach.

Such recent approaches include \( M^* \), \( \text{dRRT} \), and \( \text{dRRT}^* \). A common ground to these techniques is an implicit construction of a composite roadmap \( \mathcal{G} \) (Solovey et al., 2016), which results from a tensor product between \( m \) single-robot roadmaps \( G_1, \ldots, G_m \): for every \( 1 \leq i \leq m \) the graph \( G_i \) is a \( \text{PRM} \) embedded in the configuration space of robot \( i \). It was recently proved that if \( G_1, \ldots, G_m \) are AO, with respect to the individual robots for which they are defined, then \( \mathcal{G} \) is AO with respect to the multi-robot problem (Shome et al., 2019).

Thus, by Theorem 1 it follows that the underlying single-robot \( \text{PRM} \) graphs \( G_1, \ldots, G_m \) can be constructed using a smaller connection radius, while still guaranteeing the AO (or AnO) of \( \mathcal{G} \). This, in turn, guarantees AO (or AnO) of \( M^* \), \( \text{dRRT}^* \), and PC of \( \text{dRRT} \).

8. Experimental results

We present experiments demonstrating the effect of using different values of the connection radius \( r_n \) on the running time, cost of solution for the length cost \( c_l \), and success rate, when running the algorithms \( \text{PRM} \) and \( \text{FMT}^* \) on problems of dimensions up to 12. In particular, \( r_n \) ranges between the critical radius (Theorem 1) and previously obtained upper bounds from Janson et al. (2015) and Karaman and Frazzoli (2011).

We validate our theory for \( \text{PRM} \) (Theorem 1) and \( \text{FMT}^* \) (Corollary 1). We observe that smaller connection radii, than previously obtained bounds, still allow the planners to converge to high-quality, near-optimal paths. Furthermore, we identify situations in which using a radius that is close to \( r^*_r \) allows to obtain a high-quality solution more quickly. Moreover, although the resulting cost for the smaller radii can be slightly worse, we observe that postprocessing the paths using standard simplification methods yields solutions that are only marginally inferior to the best (postprocessed) solution. Specifically, in more difficult scenarios the advantage in using a smaller connection radius is more prominent; in some cases we obtain a reduction of 50% in running time, with an improved cost, and similar success rates when compared with the results obtained using the original \( \text{FMT}^* \) connection radius.

8.1. Implementation details

In our experiments, we used the Open Motion Planning Library (OMPL 1.3) (Sucan et al., 2012) on a 2.6 GHz x 2 Intel Core i5 processor with 16 GB of memory. Results were averaged over 50 runs and computed for dimensions up to 12.

The planners that we used are \( \text{PRM} \) and the batch variant of \( \text{FMT}^* \), which were adapted for samples from a PPP, where \( n \) is the expected number of samples. Specifically, given the expected number of samples \( n \), these variants generate a set of samples according to the recipe in Claim 1. The two planners use the connection radii \( r_{n,1}, r_{n,2}^* \). Note that \( r_{n,1}^* \) should be at least \( \frac{n}{\beta \log^{1/\alpha} n} \), but the exact value of \( \beta \) is unknown. For simplicity, we set \( r_{n,1}^* \) to be identical to \( r_{\text{post}}^* \), defined in Karaman and Frazzoli (2011). We emphasize that although we use an asymptotically smaller value, it still yields (empirically) convergence in cost. This suggests that the bound on \( r_{n,1}^* \) can be further reduced.

Given a scenario and a value \( n \), we define a set of \( k+1 \) increasing connection radii, \( \{r_0, \ldots, r_k\} \), as follows. We set the minimal connection radius to be \( r_0 = \gamma n^{-1/d} \), where \( \gamma = 1 \). Note that \( \gamma \) is larger than \( \gamma^* \) by a factor of roughly 2. The maximal connection radius, denoted by \( r_k = r_{\text{post}}^* \), is as defined in Janson et al. (2015). For each \( 1 \leq i \leq k \), we define \( r_i = r_0 + i \Delta \), where \( \Delta = (r_k - r_0)/k \). Now, for every scenario and number of samples \( n \) we run our planning algorithm with \( r_{n,i}^* \), and \( r_i \in \{r_0, \ldots, r_k\} \). Note that all our plots are for \( k = 10 \), and that in some experiments an additional radius \( r_{k+1} = r_{\text{post}}^*>r_{\text{post}}^* \) appears as well (see Karaman and Frazzoli, 2011). Figure 5 shows the colors and labeling that will be used throughout this section.

8.2. Results

8.2.1. Euclidean space. The scenario we consider (see Figure 6a) consists of a point robot moving in the obstacle-free unit \( d \)-dimensional hypercube. Therefore,
$F = C = [0, 1]^d$. We set the start and target positions of the robot to be $s = (0.1, \ldots, 0.1)$ and $t = (0.9, \ldots, 0.9)$, respectively.

We run PRM and plot (i) the overall running time, (ii) the normalized cost ($c_i$) of the obtained solution, where a value of 1 represents the best possible cost, and (iii) the portion of successful runs, all as a function of the expected number of samples $n$. Results are depicted in Figure 7.

The plots demonstrate the following trend: for each radius $r$ the cost obtained by PRM converges to some constant times the optimal cost, which is marked with the dashed red curve in the “Cost vs. $n$” plot. (We note that it is possible that for larger values of $n$ the cost values for different radii will eventually converge to the same value.) Clearly, $r_{FDB*}$ yields the best cost but at the price of increased running time. Here $r_{10} = r_{FDB*}$ obtains the next best cost, with improved running time, and so on. Note that for $d = 4$, already for $n = 1K$ a solution is found for all radii except $r_0$, whereas for $d = 8$ and $n = 5K$ a solution is found for all radii above $r_4$. It is important to note that there is a clear speedup in the running times of PRM when using $r_i$ for $i \leq 9$, over $r_{FMT}^*$, $r_{PRM*}$, with a slight penalty (a factor of roughly 2) in the resulting costs.

For the first set of experiments of PRM in an obstacle-free $d$-dimensional hypercube, we also study how the value of $r_i$ affects the size of the connected components. Note that Theorem 1 of Penrose and Pisztora (1996) states that $|C_n| = \Theta(n)$. In particular, we measure the size of the two...
largest connected components, denoted by $C_n$ and $C'_n$, as a function of both the radius $r_i$ and the expected number of samples $n$. The results for $d = 2$, $d = 12$ are summarized in Table 1. Already for $r_2$, $C_n$ is significantly larger than $C'_n$. However, for $r_0$ there is no clear difference. That is, we do not see in practice the expected emergence of the “huge” component. As $n$ increases, the proportion of $C_n$ increases as well, whereas that of $C'_n$ shrinks. Also note that for specific $r_i$, $n$ the maximal component is smaller for $d = 12$ than for $d = 2$.

8.2.2. General Euclidean space. We consider the following $d$-dimensional scenario (based on a scenario from Kiril
(2018)), for $d \in \{4, 8\}$, depicted in Figure 6b in three dimensions, and a point robot: $C = [0,1]^d$ is subdivided into $2^d$ sub-cubes by halving it along each axis. Each sub-cube contains a centered $d$-dimensional axis-aligned hypercubical obstacle that covers 25% of the sub-cube. The start position $s$ of the robot is placed on the diagonal between $(0, \ldots, 0)$ and $(1, \ldots, 1)$, such that it is equidistant from the origin and from the closest hypercubical obstacle. Here $t$ is selected similarly, with respect to $(1, \ldots, 1)$.

Here we use $\text{FMT}^\star$ with radii ranging from $r_0$ to $r_{10}$. We also ran $\text{PRM}$ which exhibited a similar behavior. Figure 8a presents the results for $d = 4$. We plot the average cost after simplifying the resulting paths in addition to the average original cost. We mention that there is a difference in cost between the various radii and that costs obtained using larger radii are often better. However, applying OMPL’s default path-simplification procedure yields paths with negligible differences in cost. We do mention, though, that the success rates of all radii deteriorate, when compared with $d = 4$.

8.2.3. General non-Euclidean space. We use the cubicles scenario (see Figure 6c) provided with the OMPL distribution (Sucan et al., 2012). Here the goal is to find a collision-free path for one or two L-shaped robots that need to exchange their positions. In the single-robot case, the robot needs to move between the two configurations depicted in red and green. In the two-robot case, these two configurations represent the start positions of the robots, where the goal is for each robot to finish at the start of the other robot. Since the robots are allowed to translate and rotate, then $d = 6$ or $d = 12$, respectively, and $C$ is non-Euclidean. Although our theory does not directly apply here, we chose to test it empirically for such a setting.
We ran our experiments with FMT\(^*\) using a connection radii ranging from \(r_0\) to \(r_{10} = r_{FMT}\). Initially, no solution was found in all runs. Indeed, as was mentioned in Kleinbort et al. (2016), this is not surprising, because the theory from which the radius values are derived assumes that the configuration space is Euclidean. In the same paper the authors proposed a heuristic for effectively using radius-based connections in motion planning. In our experiments (Figure 9b) we increased all radii by a multiplicative factor of 3 and 10, respectively, in order to increase the success rates. We mention that this yielded similar behavior to that when using the connection scheme suggested by Kleinbort et al. (2016).

Results for the single-robot case are depicted in 9a. In this case, even with 25,000 samples already with \(r_3\) (green) we were able to obtain high success rates. The running time was slightly better than that of \(r_{10}\), whereas the cost was higher than that obtained using \(r_{10}\). However, after applying the smoothing procedure the difference in cost decreased significantly.

Let us proceed to the two-robot setting. Observe in Figure 9b, that the variance in cost after path simplification is again significantly smaller than the variance in the original cost. Clearly, smaller radii exhibit shorter running times, but with smaller success rates. However, because the success rate improves as the number of samples \(n\) increases, one could use the smaller radii with larger \(n\) and still obtain a solution with comparable cost and success rate in shorter running time. Indeed, using \(r_3\) (green) with 75,000 samples we obtain a solution whose cost (after simplification) is slightly better than the cost obtained using \(r_{10}\) with 42,000 samples (after simplification). Moreover, the running time of the former is roughly 50% of the time taken for the latter, and both obtain similar success rates. This indicates that in such settings, one could benefit from using smaller radii in terms of favorable running times and obtain a comparable or even better cost with similar success rates.

9. Future work

In this work, we leveraged techniques from percolation theory to develop stronger analysis of \(PRM\)-based sampling-based motion planners. In the hope of providing mathematically rigorous presentation, while still being accessible, we chose to focus the discussion on simplified, possibly unrealistic, robotic systems. In particular, we assumed a holonomic system having a Euclidean configuration space.

Our immediate future goal is to extend our theory to non-Euclidean spaces, such as those arising from rigid bodies translating and rotating in space. The next challenge will be to extend the model to robots with differential constraints. While the latter task seems daunting, we should keep in mind that the state space of such systems can be modeled as a differential manifold. This may allow to locally apply our Euclidean-space techniques to analyze manifold spaces. Indeed, a similar approach has already been considered in previous work (Schmerling et al., 2015a,b).

The next algorithmic challenge is to consider robotic systems which for precise steering, i.e., solving the two-point boundary value problem, cannot be performed, at least not efficiently (see the discussion in Li et al., 2016: Section 1.3). In such cases, \(PRM\)-based planners (as we considered here), or \(RRT\)^* -based techniques, cannot be applied. We pose the following question: Is it possible to extend existing planners to work in the absence of a steering function, while maintaining their theoretical properties? We plan to tackle this question in the near future.

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Notes

1 AnO means that the cost of the solution tends to at most a constant factor times the optimum, compared with AO in which this constant is equal to one.
2 Let \(A_1, A_2, \ldots\) be random variables in some probability space and let \(B\) be an event depending on \(A_n\). We say that \(B\) occurs asymptotically almost surely (a.a.s.) if \(\lim_{n \to \infty} \Pr(B|A_n) = 1\)

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A.1. The basics

We start with some basic definitions. For a given $n$ define $\mathcal{L}_n = \mathbb{Z}_d$ and $\mathcal{L}_n = \mathbb{Z}_d$. That is, $\mathcal{L}_n$ contains all the points of a rectangular lattice, with edge length $\mathcal{L}_n$. Observe that $\mathcal{G}(\mathcal{L}_n; \mathcal{L}_n)$ is a deterministic graph which connects by an edge every two points $x, x' \in \mathcal{L}_n$ that differ in exactly one coordinate and the distance between them is exactly $\mathcal{L}_n$. We proceed to define a randomized subset of $\mathcal{G}(\mathcal{L}_n; \mathcal{L}_n)$:

**Definition 11.** Given $0 \leq p \leq 1$, the graph $\mathcal{G}_p = \mathcal{G}(\mathcal{L}_n; \mathcal{L}_n; p)$ is defined in the following manner. Every point of $\mathcal{L}_n$ is added independently with probability $p$ to $\mathcal{G}_p$ as a vertex. Two vertices $x, x' \in \mathcal{G}_p$ are connected by an edge if $|x - x'| = 2\mathcal{L}_n$.

This corresponds to a well-studied model of site percolation. Similarly to the continuous model studied in the previous section, we say that $\mathcal{G}_p$ percolates if it contains an infinite connected component $C^p_0$ that includes the origin $o$. Also denote by $C^p_0$ the largest connected component of $\mathbb{Z}_d$.

The following theorem corresponds to a combination of Theorem 2, Lemma 1, and Theorem 3 for the continuous case.

**Theorem 6** (Franceschetti and Meester, 2008). There exists a critical value $0 < p^* < 1$ such that for $p < p^*$ the graph $\mathcal{G}_p$ percolates with probability zero. If $p > p^*$ then $\mathcal{G}_p$ percolates with non-zero probability. Furthermore, if $p > p^*$ then $\mathcal{G}_p$ is unique with probability 1. See Appendix C for values of $p^*$.

The next two lemmas serve as discrete versions of Lemmas 2 and 3, respectively. Their proofs are very similar to the original proofs and are therefore omitted. We do mention that they rely on Theorems 5 and 4 of Penrose and Pisztora (1996), respectively.

**Lemma 5.** Suppose that $p > p^*$. Denote by $\mathcal{E}'_n 1$ the event that $C^p_0 \cap \mathcal{H}_n \subset C^p_0$, where $\mathcal{H}_n$ is as defined in Lemma 2. Then there exists $n_0 \in \mathbb{N}$ and $\alpha > 0$ such that for any $n > n_0$ it holds that

$$\Pr[\mathcal{E}'_n 1] \geq 1 - \exp\left(-\alpha n^{1/d} \log^{-1} n\right)$$

**Lemma 6.** Suppose that $p > p^*$. Define $H'_n \subset \mathcal{H}_n$ to be a hypercube of side length $h'_n = \beta' \log^{1/(d-1)} n / n^{1/d}$. Denote by $\mathcal{E}'_n 2$ the event that $H'_n \cap C^p_0 \neq \emptyset$. Then there exist $n_0 \in \mathbb{N}$ and $\beta'_0 > 0$ such that for any $n > n_0, \beta' > \beta'_0$ it holds that $\Pr[\mathcal{E}'_n 2 | \mathcal{E}'_n 1] \geq 1 - n^{-1}$.

Appendix A. Theory for (semi-)deterministic sampling

So far we have studied the theoretical properties of various sampling-based planners constructed with the PPP sample set $X_n$. In this appendix, we study similar aspects but for a different sampling regime. In particular, we replace $X_n$ with $\mathcal{L}_n$ which represents a rectilinear lattice, also known as a Sukharev sequence. A recent paper by Janson et al. (2017) studied the theoretical properties of PRM and FMT when constructed with $\mathcal{L}_n$ (and other deterministic low-dispersion sequences). Here we introduce a new analysis which shows that PRM constructed with $\mathcal{L}_n$ can be further sparsified while maintaining AO. This result can also be extended to FMT*, BT, RRG, and other planners, as was done in Section 7 for the setting of PPP.
Now we proceed to the discrete counterpart of Theorem 4. For any \(v \in \mathbb{R}^d\) define \(x(v) = \arg\min_{x \in L_v} \|x - v\|\). The following is a simplified version of Antal and Pisztora (1996: Theorem 1.1).

**Theorem 7.** Suppose that \(p > p^*\). There exists a positive constant \(\xi' \geq 1\) such that for any two fixed points \(v, v' \in \mathbb{R}^d\) it holds that

\[
\Pr\{x(v), x(v') \in C_v, \text{dist}(G^*_n, x(v), x(v')) \\
\geq \xi' \|x(v) - x(v')\|\} = o(1)
\]

For the subcritical regime \((p < p^*)\), we have the following theorem. Define

\[
S_k = \left\{ x \in L_v | \|x\|_{\tilde{r}^*_n} \leq k \right\}
\]

That is, \(S_k\) represents all the points in \(L_v\) that can be reached from \(o\) by at most \(k\) edges of length \(\tilde{r}^*_n\) over the corresponding rectilinear lattice. Denote by \(\partial S_k\) the boundary of \(S_k\), i.e., all the points that are exactly \(k\) edges from \(o\). Let \(A_k\) be the event that \(G^*_n\) contains a path from \(o\) to some vertex in \(\partial S_k\).

**Theorem 8** (Grimmert, 1999, Theorem 5.4). If \(p < p^*\), there exists a constant \(c>0\) such that \(\Pr[A_k] < e^{-ck}\) for all \(k\).

This theorem implies that in the subcritical regime it is improbable that \(G^*_n\) has a path connecting two lattice points \(x(v), x(v')\) as defined above, as such a path must have \(o(1)\) edges.

**A.2. Back to motion planning**

To conclude this section, we define a sparsified version of PRM defined on the deterministic samples \(L_v\), and state its theoretical properties. The proof is omitted as it is almost identical to that obtained in Section 4, with the difference that it now uses theorems and lemmas presented earlier this section, rather than those from Section 5.

Recall that PRM uses the two connection radii \(r_n, r'_n\). For simplicity, we fix \(r_n = \tilde{r}^*_n\).

**Definition 12.** The sparsified PRM graph \(\mathcal{P}^*_n\) is the union between \(G^*_n \cap \mathcal{F}\) and the supplementary edges

\[
\bigcup_{(x, y) \in \mathcal{F}} \{(x, y) | x \in G^*_n \cap B_{\hat{d}}(y) \text{ and } xy \subset \mathcal{F}\}
\]

**Theorem 9.** Suppose that \((\mathcal{F}, s, t)\) is robustly feasible. Then there exists \(p^* > 0\) such that the following holds.

(i) If \(p < p^*\) and \(r'_n = \infty\), then PRM fails (to find a solution) a.a.s.

(ii) Suppose that \(p > p^*\). There exists \(\beta'_0 > 0\) such that for \(r'_n = \beta' \log_{\frac{1}{p}} p\), where \(\beta' > \beta'_0\) and any \(\epsilon > 0\) the following holds a.a.s.: 

1. \(\mathcal{P}^*_n\) contains a path \(\pi_n \in \Pi_{r'_n}^F\) with \(c_0(\pi_n) \leq (1 + \epsilon)\xi' \cdot \ell_{\pi_n}\).

2. If \(\mathcal{M}\) is well behaved, then \(\mathcal{P}^*_n\) contains a path \(\pi'_n \in \Pi_{r'_n}^F\) with \(c_b(\pi'_n, M) \leq (1 + \epsilon)c'_b\).

**Appendix B. Probabilistic completeness of RRT with samples from a PPP**

This section is devoted to the proof of Claim 6. A probabilistic completeness proof of RRT, for the setting of uniform random sampling, is given in Michal et al. (2019: Section 3). We describe here how to modify this proof for samples from a PPP, which are of interest in the present paper.

Note that RRT is an incremental planning algorithm, which generates samples one after the other, and connects the current sample to previous ones. Therefore, it should employ the incremental PPP sampling defined in Claim 5.

We wish to apply the line of arguments as for the uniform-sampling case described in Michal et al. (2019: Theorem 1). This theorem assumes that a path \(\pi\) from start to target exists, whose clearance is \(\delta > 0\) and length is \(L\). The path \(\pi\) is covered with a set of \(m\) balls of radius \(r\), where \(m = L/r\) and \(r\) is a constant depending on both the steering parameter \(\eta\) of RRT and the clearance \(\delta\). The proof describes the process as a Markov chain with transition probability \(p\), where \(p\) is the probability to produce an RRT vertex in the \((i + 1)\)th ball \(B_{\delta}(x_{i+1})\) given that there exists an RRT vertex in the \(i\)th ball \(B_{\delta}(x_i)\). It then shows that the probability that RRT will not produce samples within all \(m\) balls decays to zero exponentially as the number of samples tends to infinity. To adapt the proof for samples from a PPP, it suffices to outline how to compute the transition probability \(p\).

Recall that at each iteration \(i\) we draw a sample \(X_i \in \mathcal{N} \in N\), where \(N\) is a Poisson random variable with mean 1, and define \(X_i = \{X_i, X_{i+1}, \ldots, X_{i+N}\}\). \(N_i\) denotes the number of points chosen independently and uniformly at random from \([0, 1]^d\). Let \(\mathcal{A}\) denote the event that \(N_i \geq 1\), and let \(\mathcal{B}\) denote the event that at least one of \(X_i, X_{i+1}, \ldots, X_{i+N}\) falls inside \(B_{\delta}(x_{i+1})\). Since \(p = \Pr[\mathcal{A} \cap \mathcal{B}] = \Pr[\mathcal{A}] \cdot \Pr[\mathcal{B} | \mathcal{A}]\), we will bound \(\Pr[\mathcal{A}], \Pr[\mathcal{B} | \mathcal{A}]\). Observe that

\[
\Pr[\mathcal{A}] = \Pr[N_i \geq 1] = 1 - \Pr[N_i = 0] = 1 - e^{-1}
\]

and note that \(\Pr[\mathcal{B} | \mathcal{A}] \geq |B_{\delta}||\mathcal{F}|\). Therefore,

\[
p = \Pr[\mathcal{A} \cap \mathcal{B}] \geq (1 - e^{-1}) \cdot |B_{\delta}|/|\mathcal{F}|
\]

The rest of the proof of Claim 6 follows the same lines as Michal et al. (2019: Theorem 1).

**Appendix C. Critical values**

In this section, we provide for reference a list of (estimates) of the values \(\gamma^*\) and \(p^*\).
C.1. Continuum percolation

Table 2 presents an estimate of $v^*$ in continuum percolation. It is directly derived from (Torquato and Jiao, 2012: Table I), where an estimate of the critical node degree $\Delta(n, r_n)$ is derived. We mention that for $d$ large enough $v^* \sim \frac{1}{2b_d}$ (Torquato and Jiao, 2012), where $b_d$ is volume of the Lebesgue measure of the unit ball in $\mathbb{R}^d$.

C.2. Lattice percolation

Table 3 presents the value of $p^*$ for $d = 2$, the best known estimates for $4 \leq d \leq 13$. Numbers in round brackets are single standard deviations.